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The AFLOW Library of Crystallographic Prototypes: Part 3

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Abstract

The *AFLOW Library of Crystallographic Prototypes* has been extended to include a total of 1,100 common crystal structural prototypes (510 new ones with Part 3), comprising all of the inorganic crystal structures defined in the seven-volume *Strukturbericht* series published in Germany from 1937 through 1943. We cover a history of the *Strukturbericht* designation system, the evolution of the system over time, and the first comprehensive index of inorganic *Strukturbericht* designations ever published.

Keywords: Crystal Structure, Space Groups, Wyckoff Positions, Lattice Vectors, Basis Vectors, Database

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Prototypes

***P*1 (1)** .....
1. $NaC_5H_{11}O_8S$: A5B11CD8E_aP26_1_5a_11a_a_8a_a 35

***P*$\bar{1}$ (2)** .....
1. $Ni(NO_3)_2(H_2O)_6$: A12B2CD12_aP54_2_12i_2i_i_12i 38
2. $Co_2B_2O_5$: A2B2C5_aP18_2_2i_2i_5i 42
3. Kyanite (Al_2SiO_5, $S0_1$): A2B5C_aP32_2_4i_10i_2i 44
4. α-$Ho_2Si_2O_7$: A2B7C2_aP44_2_4i_14i_4i 47
5. $Co_3(SeO_3)_3{\cdot}H_2O$: A3B2C10D3_aP36_2_ah2i_2i_10i_3i 51
6. δ-WO_3: A3B_aP32_2_12i_4i 54
7. Chalcanthite ($CuSO_4{\cdot}5H_2O$, $H4_{10}$): AB10C9D_aP42_2_ae_10i_9i_i 58
8. Boric Acid (H_3BO_3, $G5_1$): AB3C3_aP28_2_2i_6i_6i 62
9. Wollastonite ($CaSiO_3$): AB3C_aP30_2_3i_9i_3i 65
10. Albite ($NaAlSi_3O_8$, $S6_8$): ABC8D3_aP26_2_i_i_8i_3i 68
11. TaTi (BCC SQS-16): AB_aP16_2_4i_4i 71

***P*2$_1$ (4)** .....
1. $W_2O_3(PO_4)_2$: A11B2C2_mP60_4_22a_4a_4a 73
2. $Li_2SO_4{\cdot}H_2O$ ($H4_8$): A2B2C5D_mP20_4_2a_2a_5a_a 77
3. Ca_3UO_6: A3B6C_mP20_4_3a_6a_a 79

***C*2 (5)** .....
1. Bassanite [$CaSO_4(H_2O)_{0.5}$, $H4_7$]: A2B2C9D2_mC90_5_ab2c_3c_b13c_3c 81
2. $NbAs_2$: A2B_mC12_5_2c_c 85
3. $D0_{15}$ ($AlCl_3$) (*obsolete*): AB3_mC16_5_c_3c 87
4. $Rb_2CaCu_6(PO_4)_4O_2$: AB6C18D4E2_mC62_5_a_2b2c_9c_2c_c 89
5. *C*2 $(Ba,Ca)CO_3$: ABC3_mC10_5_b_a_ac 92

***Pm* (6)** .....
1. Ta_5Ti_{11} (BCC SQS-16): A5B11_mP16_6_2abc_2a3b3c 94

***Pc* (7)** .....
1. $Na_2Ca_6Si_4O_{15}$: A6B2C15D4_mP54_7_6a_2a_15a_4a 96
2. Low-Temperature Mo_8O_{23}: A8B23_mP124_7_16a_46a 100
3. Calaverite ($AuTe_2$): AB2_mP12_7_2a_4a 107

***Cm* (8)** .....
1. Monoclinic Co_4Al_{13}: A13B4_mC102_8_17a11b_8a2b 109
2. $TaTi_3$ (BCC SQS-16): AB3_mC32_8_4a_12a 113
3. $TaTi_3$ (BCC SQS-16): AB3_mC32_8_4a_4a4b 115
4. $F5_{11}$ (KNO_2) (*obsolete*): ABC2_mC8_8_a_a_b 117

***Cc* (9)** .....
1. Nacrite [$Al_2Si_2O_5(OH)_4$, $S5_4$]: A2B4C9D2_mC68_9_2a_4a_9a_2a 119
2. Chrysotile ($Mg_3Si_2O_5(OH)_4$): A3B5C4D2_mC56_9_3a_5a_4a_2a 122
3. $Cs_6W_{11}O_{36}$: A6B36C11_mC212_9_6a_36a_11a ... 125

***P*2$_1$/*m* (11)** .....
1. $K_2S_2O_5$ ($K0_1$): A2B5C2_mP18_11_2e_e2f_2e 131
2. $ZrSe_3$: A3B_mP8_11_3e_e 133
3. *y*-$Y_2Si_2O_7$: A7B2C2_mP22_11_3e2f_2e_ab 135
4. Barytocalcite ($BaCa(CO_3)_2$): AB2CD6_mP20_11_e_2e_e_2e2f 137
5. O(OH)Y: ABC_mP6_11_e_e_e 139

***C*2/*m* (12)** .....
1. $Al_{13}Fe_4$: A13B4_mC102_12_dg8i5j_4ij 141
2. Os_4Al_{13}: A13B4_mC34_12_b6i_2i 144
3. Bischofite ($MgCl_2{\cdot}6H_2O$, $J1_7$): A2B12CD6_mC42_12_i_2i2j_a_ij 146
4. Tremolite ($Ca_2Mg_5Si_8O_{22}(OH)_2$, $S4_2$): A2B2C5D24E8_mC82_12_h_i_agh_2i5j_2j 148
5. β-Ga_2O_3: A2B3_mC20_12_2i_3i 151
6. $K_2Ti_2O_5$: A2B5C2_mC18_12_i_a2i_i 153
7. CaC_2-III: A2B_mC12_12_2i_i 155
8. Tolbachite ($CuCl_2$): A2B_mC6_12_i_a 157
9. δ-Ni_3Sn_4 ($D7_a$): A3B4_mC14_12_ai_2i 159
10. $LiOH{\cdot}H_2O$ ($B36$): A3BC2_mC24_12_ij_h_gi 161
11. Staurolite ($Al_5Fe_2O_{10}(OH)_2Si_2$): A5B2C10D2E2_mC84_12_acghj_bdi_5j_2i_j 163
12. Manganese-leonite [$K_2Mn(SO_4)_2{\cdot}4H_2O$, $H4_{23}$]: A8B2CD15E2_mC112_12_2i3j_j_ad_g4i5j_2i 167
13. Monoclinic $FeTlSe_2$: AB2C_mC16_12_g_2i_i 171

14. $NbTe_2$: AB2_mC18_12_ai_3i 173
15. $D2_2$ ($MgZn_5$?) (*problematic*): AB5_mC48_12_2i_ac5i2j 175
16. Chrysotile ($H_4Mg_3Si_2O_9$, $S4_5$): AB6C11D6E4_mC112_12_e_gi2j_i5j_2i2j_2j 178
17. Sr_2NiTeO_6: AB6C2D_mC40_12_ad_gh4i_j_bc 182
18. Ta_2PdSe_6: AB6C2_mC18_12_a_3i_i 184
19. Sanidine ($KAlSi_3O_8$, $S6_7$): AB8C4_mC52_12_i_gi3j_2j 186
20. $MnPS_3$: ABC3_mC20_12_g_i_ij 189
21. $AlNbO_4$: ABC4_mC24_12_i_i_4i 191
22. SiAs: AB_mC24_12_3i_3i 193

***P2/c* (13)** .....

1. High-Temperature Mo_8O_{23}: A8B23_mP62_13_4g_c11g 195
2. Huanzalaite ($MgWO_4$, $H0_6$): AB4C_mP12_13_f_2g_e 199

***P2$_1$/c* (14)** .....

1. $Cs_{11}O_3$: A11B3_mP56_14_11e_3e 201
2. Azurite [$Cu_3(CO_3)_2(OH)_2$, $G7_4$]: A2B3C2D8_mP30_14_e_ce_e_4e 205
3. $HgCl_2{\cdot}2HgO$: A2B3C2_mP14_14_e_ae_e 208
4. Orpiment (As_2S_3, $D5_f$): A2B3_mP20_14_2e_3e .. 210
5. Monoclinic Cu_2OSeO_3: A2B4C_mP28_14_abe_4e_e 213
6. $Sb_4O_5Cl_2$: A2B5C4_mP22_14_e_c2e_2e 216
7. Ca_2UO_5: A2B5C_mP32_14_2e_5e_ab 218
8. Gd_2SiO_5 (RE_2SiO_5 X1): A2B5C_mP32_14_2e_5e_e 221
9. Sanguite ($KCuCl_3$): A3BC_mP20_14_3e_e_e 224
10. γ-WO_3: A3B_mP32_14_6e_2e 226
11. $K_2Ni(CN)_4$: A4B2C4D_mP22_14_2e_e_2e_a 230
12. $KICl_4{\cdot}H_2O$ ($H0_{10}$): A4BCD_mP28_14_4e_e_e_e 232
13. γ-$Y_2Si_2O_7$[‖]: A4BC_mP24_14_4e_e_e 235
14. $K_2Pt(SCN)_6{\cdot}2H_2O$: A6B4C2D6E2FG6_mP54_14_3e_2e_e_3e_e_a_3e .. 238
15. K_2NbF_7 ($K6_2$): A7B2C_mP40_14_7e_2e_e 242
16. Manganese-leonite 110 K [$K_2Mn(SO_4)_2{\cdot}4H_2O$]: A8B2CD12E2_mP100_14_8e_2e_ad_12e_2e 246
17. Co_2Al_9 ($D8_d$): A9B2_mP22_14_a4e_e 252
18. Tutton salt [$Cu(NH_4)_2(SO_4)_2{\cdot}6H_2O$, $H4_4$]: AB20C2D14E2_mP78_14_a_10e_e_7e_e 255
19. Parawollastonite ($CaSiO_3$, $S3_3$ (II)): AB3C_mP60_14_3e_9e_3e 260
20. β-B_2H_6[*]: AB3_mP16_14_e_3e 264
21. B_2H_6 ($P2_1/c$)[*]: AB3_mP16_14_e_3e 266
22. $KAuBr_4{\cdot}2H_2O$ ($H4_{19}$): AB4C2D_mP32_14_e_4e_2e_e 268
23. Anhydrous $KAuBr_4$: AB4C_mP24_14_ab_4e_e ... 271
24. Ammonium Persulfate [$(NH_4)_2S_2O_8$, $K4_1$][†‖]: AB4C_mP24_14_e_4e_e 274
25. Monazite ($LaPO_4$)[†‖]: AB4C_mP24_14_e_4e_e 277
26. Sr_2MnTeO_6: AB6C2D_mP20_14_a_3e_e_d 280
27. Cryolite (Na_3AlF_6, $J2_6$): AB6C3_mP20_14_a_3e_de 282
28. KNO_2 III[‡]: ABC2_mP16_14_e_e_2e 285
29. Manganite (γ-MnO(OH), $E0_6$)[‡]: ABC2_mP16_14_e_e_2e 287
30. $AgMnO_4$ ($H0_9$)[‖]: ABC4_mP24_14_e_e_4e 289
31. Nahcolite ($NaHCO_3$, $G0_{12}$): ABCD3_mP24_14_e_e_e_3e 292
32. Cu(OH)Cl: ABCD_mP16_14_e_e_e_e 295
33. Arsenopyrite (FeAsS, $E0_7$): ABC_mP12_14_e_e_e 297
34. α-ICl[§]: AB_mP16_14_2e_2e 299
35. LiAs[§]: AB_mP16_14_2e_2e 301
36. ϵ-1,2,3,4,5,6-Hexachlorocyclohexane (C_6Cl_6): AB_mP24_14_3e_3e 303
37. Pararealgar (AsS)[¶]: AB_mP32_14_4e_4e 306
38. Realgar (AsS, B_l)[¶]: AB_mP32_14_4e_4e 309

***C2/c* (15)** .....

1. Ag_2PbO_2: A2B2C_mC20_15_ad_f_e 312
2. Catapleiite ($Na_2ZrSi_3O_9{\cdot}2H_2O$): A2B3C9D3E_mC144_15_2f_bcdef_9f_3f_ae 314
3. Na_2PrO_3: A2B3C_mC48_15_aef_3f_2e 319
4. Eudidymite ($BeHNaO_8Si_3$): A2B4C2D17E6_mC124_15_f_2f_f_e8f_3f 322
5. ζ-Nb_2O_5 (B-Nb_2O_5): A2B5_mC28_15_f_e2f 327
6. Muscovite ($KH_2Al_3Si_3O_{12}$, $S5_1$): A2BC10D2E4_mC76_15_f_e_5f_f_2f 329
7. $Rb_2C_2O_4{\cdot}H_2O$: A2BC4D2_mC36_15_f_e_2f_f 332

‖ γ-$Y_2Si_2O_7$, Ammonium Persulfate [$(NH_4)_2S_2O_8$, $K4_1$], Monazite ($LaPO_4$), and $AgMnO_4$ ($H0_9$) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

* β-B_2H_6 and B_2H_6 ($P2_1/c$) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

† Ammonium Persulfate [$(NH_4)_2S_2O_8$, $K4_1$] and Monazite ($LaPO_4$) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

‡ KNO_2 III and Manganite (γ-MnO(OH), $E0_6$) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

§ α-ICl and LiAs have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

¶ Pararealgar (AsS) and Realgar (AsS, B_l) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

8. Alluaudite [$NaMnFe_2(PO_4)_3$]: A2BCD12E3_mC76_15_f_e_b_6f_ef 334
9. ThC_2 (C_g): A2B_mC12_15_f_e 338
10. Clinocervantite (β-Sb_2O_4): A2B_mC24_15_2f_ce 340
11. $(CdSO_4)_3{\cdot}8H_2O$ ($H4_{20}$): A3B16C20D3_mC168_15_ef_8f_10f_ef 342
12. Chlorite [$Mg_3(Mg_2Al)(Si_3Al)O_{10}(OH)_8$, $S5_5$]: A3B5C4D2_mC112_15_a3ef_5f_4f_2f 347
13. Y_2SiO_5 (RE_2SiO_5 X2): A5BC2_mC64_15_5f_f_2f 351
14. α-$Zn_2V_2O_7$: A7B2C2_mC44_15_e3f_f_f 354
15. Manganese-leonite 185 K [$K_2Mn(SO_4)_2{\cdot}4H_2O$]: A8B2CD12E2_mC200_15_8f_2f_ce_2e11f_2f 357
16. $BaNi(CN)_4{\cdot}4H_2O$ ($H4_{22}$): AB4C4D4E_mC56_15_e_2f_2f_2f_a 363
17. Gypsum ($CaSO_4{\cdot}2H_2O$, $H4_6$): AB4C6D_mC48_15_e_2f_3f_e 366
18. Ta_2NiSe_5: AB5C2_mC32_15_e_e2f_f 369
19. Pyrophyllite [$AlSi_2O_5(OH)$, $S5_6$]: AB5CD2_mC72_15_f_5f_f_2f 371
20. Titanite ($CaTiSiO_5$, $S0_6$): AB5CD_mC32_15_e_e2f_e_b 374
21. $KFeS_2$ ($F5_a$): ABC2_mC16_15_e_e_f 376
22. Diopside [$CaMg(SiO_3)_2$, $S4_1$]: ABC6D2_mC40_15_e_e_3f_f 378
23. β-Ga (*obsolete*): A_mC4_15_e 381

$P222_1$ (17)

1. $NaNbO_3$: ABC3_oP40_17_abcd_2e_abcd4e 383

$P2_12_12$ (18)

1. γ-TeO_2: A2B_oP12_18_2c_c 386
2. Diamminetriamidodizinc Chloride ($[Zn_2(NH_3)_2(NH_2)_3]Cl$): AB12C5D2_oP40_18_a_6c_b2c_c 388

$P2_12_12_1$ (19)

1. Morenosite ($NiSO_4{\cdot}7H_2O$, $H4_{12}$): A14BC11D_oP108_19_14a_a_11a_a 391
2. Wülfingite (ϵ-$Zn(OH)_2$, $C31$): A2B2C_oP20_19_2a_2a_a 396
3. Ferroelectric $NH_4H_2PO_4$: A6BC4D_oP48_19_6a_a_4a_a 398
4. β-Arabinose [$(CH_2O)_{20}$]: AB2C_oP80_19_5a_10a_5a 401
5. $NaAlCl_4$: AB4C_oP24_19_a_4a_a 405
6. NaP: AB_oP16_19_2a_2a 407

$C222_1$ (20)

1. $D0_7$ (CrO_3) (*obsolete*): AB3_oC16_20_a_bc 409
2. $AlPO_4$ "low cristobalite type": AB4C_oC24_20_b_2c_a 411
3. Tl_2AlF_5 ($K3_3$): AB5C2_oC32_20_b_a2bc_c 413

C222 (21)

1. $HoSb_2$: AB2_oC6_21_a_k 415

F222 (22)

1. Predicted Phase IV $Cd_2Re_2O_7$: A2B7C2_oF88_22_k_bdefghij_k 417

$Pca2_1$ (29)

1. Mercury (II) Azide [$Hg(N_3)_2$]: AB6_oP28_29_a_6a 420
2. Low-Temperature $(NH_3CH_3)Al(SO_4)_2{\cdot}12H_2O$: ABC30DE20F2_oP220_29_a_a_30a_a_20a_2a 422

$Pmn2_1$ (31)

1. Orthorhombic Co_4Al_{13}: A13B4_oP102_31_17a11b_8a2b 430
2. $Mg(ClO_4)_2{\cdot}6H_2O$ ($H4_{11}$): A2B6CD8_oP34_31_2a_2a2b_a_4a2b 435
3. B_4SrO_7: A4B7C_oP24_31_2b_a3b_a 438
4. $D8_7$ (Shcherbinaite, V_2O_5) (*obsolete*): A5B2_oP14_31_a2b_b 440

Pba2 (32)

1. $Mo_{17}O_{47}$: A17B47_oP128_32_a8c_a23c 442

$Pna2_1$ (33)

1. Possible δ-$Gd_2Si_2O_7$: A2B7C2_oP44_33_2a_7a_2a 447
2. CaB_2O_4 (III): A2BC4_oP84_33_6a_3a_12a 450
3. Cervantite (α-Sb_2O_4): A2B_oP24_33_4a_2a 455
4. CsB_4O_6F: A4BCD6_oP48_33_4a_a_a_6a 457
5. $LiGaO_2$: ABC2_oP16_33_a_a_2a 460
6. γ-$LiIO_3$: ABC3_oP20_33_a_a_3a 462

Pnn2 (34)

1. $MnF_{2-x}(OH)_x$: A2BC2_oP10_34_c_a_c 464

$Cmc2_1$ (36)

1. Si_2N_2O: A2BC2_oC20_36_b_a_b 466
2. Bi_2GeO_5: A2BC5_oC32_36_b_a_a2b 468
3. Ni_3Si_2: A3B2_oC80_36_4a4b_2a3b 470
4. Bertrandite ($Be_4Si_2O_7(OH)_2$, $S4_6$): A4B7C2D2_oC60_36_2b_a3b_2a_b 473
5. MoP_2: AB2_oC12_36_a_2a 476
6. α-Potassium Nitrate (KNO_3) II: ABC3_oC80_36_2ab_2ab_2a5b 478

Amm2 (38)

1. Ta_3Ti_{13} (BCC SQS-16): A3B13_oC32_38_ac_a2bcdef 481
2. Ta_3Ti_5 (BCC SQS-16): A3B5_oC32_38_abce_abcdf 483
3. $NaNb_6O_{15}F$: ABC6D15_oC46_38_b_b_2a2d_2ab4d2e 485

Ama2 (40)

1. $Rb_2Mo_2O_7$: A2B7C2_oC88_40_abc_2b6c_a3b 488
2. Orthorhombic CrO_3: AB3_oC16_40_b_a2b 491

Aba2 (41)

1. Santite ($KB_5O_8{\cdot}4H_2O$, $K3_5$): A5B8CD12_oC104_41_a2b_4b_a_6b 493

Fdd2 (43)

1. Ag_2O_3[**]: A2B3_oF40_43_b_ab 497
2. Natrolite ($Na_2Al_2Si_3O_{10}{\cdot}2H_2O$, $S6_{10}$): A2B4C2D12E3_oF184_43_b_2b_b_6b_ab 499
3. Blossite (α-$Cu_2V_2O_7$): A2B7C2_oF88_43_b_a3b_b 504
4. Archerite (KH_2PO_4): A2BC4D_oF64_43_b_a_2b_a 507
5. Cs_2Se: A2B_oF24_43_b_a 510
6. Zr_2Al_3[**]: A3B2_oF40_43_ab_b 512

***Imm2* (44)** .....
1. Hemimorphite ($Zn_4Si_2O_7(OH)_2{\cdot}H_2O$, $S2_2$): A2B5CD2_oI40_44_2c_abcde_d_e 514
2. Ferroelectric $NaNO_2$ ($F5_5$): ABC2_oI8_44_a_a_c 516
3. $AgNO_2$ ($F5_{12}$): ABC2_oI8_44_a_a_d 518
4. $B30$ (MgZn?): AB_oI48_44_6d_ab2cde 520

***Ima2* (46)** .....
1. $Nb_2Zr_6O_{17}$: A2B17C6_oI100_46_ab_b8c_3c 523

***Pmma* (51)** .....
1. Parkerite ($Ni_3Bi_2S_2$): AB2C_oP8_51_e_be_f 527
2. $LiNb_6O_{15}F$: ABC6D15_oP46_51_f_d_2e2i_aef4i2j 529

***Pnna* (52)** .....
1. Carnallite [$Mg(H_2O)_6KCl_3$]: A3B12CDE6_oP276_52_d4e_18e_ce_de_2d8e 532

***Pmna* (53)** .....
1. Eriochalcite ($CuCl_2{\cdot}2H_2O$, $C45$): A2BC4D2_oP18_53_h_a_i_e 544
2. NH_4HF_2 ($F5_8$): A2BC_oP16_53_eh_ab_g 546

***Pbam* (55)** .....
1. Orthorhombic $Sr_4Ru_3O_{10}$: A10B3C4_oP68_55_2e2fgh2i_adef_2e2f 548
2. $Nb_2Pd_3Se_8$: A2B3C8_oP26_55_h_ag_2g2h 551
3. $K_2HgCl_4{\cdot}H_2O$ ($E3_4$): A4BCD2_oP32_55_ghi_f_e_gh 553
4. $Ru_{11}B_8$: A8B11_oP38_55_g3h_a3g2h 556
5. $HoMn_2O_5$: AB2C5_oP32_55_g_fh_eghi 559

***Pccn* (56)** .....
1. Calciborite (CaB_2O_4 II): A2BC4_oP56_56_2e_e_4e 562

***Pbcm* (57)** .....
1. $D0_{10}$ (WO_3) (*obsolete*): A3B_oP16_57_a2d_d 566
2. $SrUO_4$: A4BC_oP24_57_cde_d_a 569
3. Lueshite ($NaNbO_3$): ABC3_oP40_57_cd_e_cd2e 571

***Pnnm* (58)** .....
1. Kotoite ($Mg_3(BO_3)_2$): A2B3C6_oP22_58_g_af_gh 574
2. Andalusite (Al_2SiO_5, $S0_2$): A2B5C_oP32_58_eg_3gh_g 576
3. Protoanthophyllite ($H_2Mg_7Si_8O_{24}$): A2B7C24D8_oP82_58_g_ae2f_2g5h_2h 579
4. In_4Se_3: A4B3_oP28_58_4g_3g 584
5. Adamite [$Zn_2(AsO_4)(OH)$, $H2_7$]: ABC5D2_oP36_58_g_g_3gh_eg 586
6. InS: AB_oP8_58_g_g 589

***Pmmn* (59)** .....
1. RuB_2: A2B_oP6_59_f_a 591
2. NH_4NO_3 IV ($G0_{11}$): A4B2C3_oP18_59_ef_ab_af 593
3. Shcherbinaite (V_2O_5) (*revised*): A5B2_oP14_59_a2f_f 595

***Pbcn* (60)** .....
1. CaB_2O_4 I ($E3_2$): A2BC4_oP28_60_d_c_2d 597
2. ζ-Fe_2N[×]: A2B_oP12_60_d_c 600
3. α-PbO_2[×]: A2B_oP12_60_d_c 602
4. Cr_5O_{12}: A5B12_oP68_60_c2d_6d 604
5. Columbite ($FeNb_2O_6$, $E5_1$): AB2C6_oP36_60_c_d_3d 608

***Pbca* (61)** .....
1. Ca_2RuO_4: A2B4C_oP28_61_c_2c_a 611
2. Tellurite (β-TeO_2, $C52$): A2B_oP24_61_2c_c 614
3. $(TiCl_4{\cdot}POCl_3)_2$: A7BCD_oP80_61_7c_c_c_c 616
4. Hambergite [$Be_2BO_3(OH)$, $G7_2$]: AB2CD4_oP64_61_c_2c_c_4c 620
5. Enstatite ($MgSiO_3$, $S4_3$): AB3C_oP80_61_2c_6c_2c 623
6. COCl: ABC_oP24_61_c_c_c 627

***Pnma* (62)** .....
1. Topaz ($Al_2SiO_4F_2$, $S0_5$): A2B2C4D_oP36_62_d_d_2cd_c 629
2. Norbergite [$Mg(F,OH)_2 \cdot Mg_2SiO_4$, $S0_7$]: A2B3C4D_oP40_62_d_cd_2cd_c 632
3. Arcanite (K_2SO_4, $H1_6$): A2B4C_oP28_62_2c_2cd_c 635
4. Anthophyllite ($Mg_5Fe_2Si_8O_{22}(OH)_2$, $S4_4$): A2B5C22D2E8_oP156_62_d_c2d_2c10d_2c_4d 638
5. Sillimanite (Al_2SiO_5, $S0_3$): A2B5C_oP32_62_bc_3cd_c 645
6. $K_2S_3O_6$ ($K5_1$): A2B6C3_oP44_62_2c_2c2d_3c 648
7. Danburite ($CaB_2Si_2O_8$, $S6_3$): A2BC8D2_oP52_62_d_c_2c3d_d 651
8. $C53$ ($SrBr_2$) (*obsolete*): A2B_oP12_62_2c_c 654

[**] Ag_2O_3 and Zr_2Al_3 have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

[×] ζ-Fe_2N and α-PbO_2 have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

9. Cs_2Sb: A2B_oP24_62_4c_2c 656
10. $RhCl_2(NH_3)_5Cl$ ($J1_8$): A3B15C5D_oP96_62_cd_3c6d_3cd_c 658
11. NH_4I_3 ($D0_{16}$): A3B_oP16_62_3c_c 663
12. Original β-WO_3 (*obsolete*): A3B_oP32_62_ab4c_2c 665
13. P_4Se_3: A4B3_oP112_62_8c4d_4c4d 668
14. Mo_4P_3: A4B3_oP56_62_8c_6c 673
15. $K_2SnCl_4{\cdot}H_2O$ ($E3_5$): A4BC2D_oP32_62_2cd_b_2c_a 676
16. $K_2SnCl_4{\cdot}H_2O$: A4BC2D_oP32_62_2cd_c_d_c 679
17. $VOSO_4$: A5BC_oP28_62_3cd_c_c 682
18. Possible δ-$Y_2Si_2O_7$: A7B2C2_oP44_62_3c2d_2c_d 684
19. $K_4[Mo(CN)_8]{\cdot}2H_2O$ ($F2_1$): A8B4C4DE8F2_oP108_62_4c2d_2d_2cd_c_4c2d_d 687
20. $SbCl_5{\cdot}POCl_3$: A8BCD_oP44_62_4c2d_c_c_c 692
21. Autunite $\{Ca[(UO_2)(PO_4)]_2(H_2O)_{11}\}$: AB22C23D2E2_oP200_62_c_11d_3c10d_d_d 695
22. Atacamite ($Cu_2(OH)_3Cl$): AB2C3D3_oP36_62_c_ac_cd_cd 704
23. NH_4CdCl_3 ($E2_4$): AB3C_oP20_62_c_3c_c 707
24. Berthierite ($FeSb_2S_4$, $E3_3$): AB4C2_oP28_62_c_4c_2c 709
25. Chalcocyanite ($CuSO_4$): AB4C_oP24_62_a_2cd_c 711
26. Rynersonite (Orthorhombic $CaTa_2O_6$): AB6C2_oP36_62_c_2c2d_d 713
27. Copper (II) Azide [$Cu(N_3)_2$]: AB6_oP28_62_c_6c 716
28. Diaspore (AlOOH, $E0_2$): ABC2_oP16_62_c_c_2c 718
29. α-Potassium Nitrate (KNO_3) I°: ABC3_oP20_62_c_c_cd 720
30. NH_4NO_3 III ($G0_{10}$)°: ABC3_oP20_62_c_c_cd 723
31. Aragonite ($CaCO_3$, $G0_2$)°: ABC3_oP20_62_c_c_cd 726
32. Epididymite ($BeHNaO_8Si_3$, $S4_7$): ABCD8E3_oP112_62_d_2c_d_4c6d_3d 728
33. NH_4ClBrI ($F5_{14}$): ABCD_oP16_62_c_c_c_c 734
34. MnCuP: ABC_oP12_62_c_c_c 736
35. η-NiSi (B_d): AB_oP8_62_c_c 738

***Cmcm* (63)**
1. $Cu_2Pb(SeO_3)_2Br_2$: A2B2C6DE2_oC52_63_g_e_fh_c_f 740
2. Pseudobrookite (Fe_2TiO_5, $E4_1$)††: A2B5C_oC32_63_f_c2f_c 742
3. $MgCuAl_2$ ($E1_a$): A2BC_oC16_63_f_c_c 744
4. $S0_4$ (Staurolite, $Fe(OH)_2Al_4Si_2O_{10}$) (*obsolete*): A4BC12D2_oC76_63_eg_c_f3gh_g 746
5. Pd_5Pu_3: A5B3_oC32_63_cfg_ce 749
6. $ZrTe_5$: A5B_oC24_63_c2f_c 751
7. Lepidocrocite (γ-FeO(OH), $E0_4$): AB2C2_oC20_63_c_f_2c 753
8. Na_2CrO_4 ($H1_8$): AB2C4_oC28_63_c_bc_fg 755
9. $ThFe_2SiC$: AB2CD_oC20_63_b_f_c_c 757
10. Mn_3As ($D0_d$): AB3_oC16_63_c_3c 759
11. Re_3B: AB3_oC16_63_c_cf 761
12. Ta_2NiS_5††: AB5C2_oC32_63_c_c2f_f 763
13. V_3AsC: ABC3_oC20_63_c_b_cf 765
14. Si_{24} Clathrate: A_oC24_63_3f 767

***Cmca* (64)**
1. Base-centered orthorhombic $Sr_4Ru_3O_{10}$: A10B3C4_oC68_64_2dfg_ad_2d 769
2. $Na_2Mo_2O_7$: A2B2C7_oC88_64_ef_df_3f2g 772

***Cmmm* (65)**
1. Li_2PrO_3: A2B3C_oC12_65_h_bh_a 775
2. $Mg(NH_3)_2Cl_2$ ($E1_3$): A2B8CD2_oC26_65_h_r_a_i 777
3. Nb_3O_7F: A3B8_oC22_65_ag_bd2gh 779

***Cmma* (67)**
1. $NH_4H_2PO_2$ ($F5_7$): A2BC2D_oC24_67_m_a_n_g ..781

***Fddd* (70)**
1. Thenardite [Na_2SO_4 (V), $H1_7$]: A2B4C_oF56_70_g_h_a 783
2. Mg_2Cu (C_b): AB2_oF48_70_g_fg 785

***Immm* (71)**
1. High-Temperature Cryolite (Na_3AlF_6): AB6C3_oI20_71_a_in_cj 787
2. $CsFeS_2$ (100 K): ABC2_oI16_71_g_i_eh 789
3. CsO: AB_oI8_71_g_i 791

***Ibam* (72)**
1. Ga_2Mg_5 ($D8_g$): A2B5_oI28_72_j_bfj 793

***Imma* (74)**
1. $Zn(NH_3)_2Cl_2$ ($E1_2$): A2B6C2D_oI44_74_h_ij_i_e .795
2. $CeCu_2$: AB2_oI12_74_e_h 797
3. $LiCuVO_4$: ABC4D_oI28_74_a_d_hi_e 799

***P*4$_2$ (77)**
1. Gwihabaite [NH_4NO_3 (V)]: A4B2C3_tP72_77_8d_ab2c2d_6d 801

***I*$\bar{4}$ (82)**
1. Kesterite [$Cu_2(Zn,Fe)SnS_4$]: A2BCD4_tI16_82_ac_b_d_g 805

***P*4/*n* (85)**
1. Bromocarnallite ($KMg(H_2O)_6(Cl,Br)_3$, $E2_6$): A3B6CD_tP44_85_bcg_3g_ac_e 807
2. $MoPO_5$: AB5C_tP14_85_c_cg_b 810

***P*4$_2$/*n* (86)**
1. $PNCl_2$ ($E1_4$): A2BC_tP32_86_2g_g_g 812
2. $Nd_4Re_2O_{11}$: A4B11C2_tP68_86_2g_ab5g_g 815
3. $NaSb(OH)_6$ ($J1_{11}$): AB6C_tP32_86_d_3g_c 819

°α-Potassium Nitrate (KNO_3) I, NH_4NO_3 III ($G0_{10}$), and Aragonite ($CaCO_3$, $G0_2$) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

††Pseudobrookite (Fe_2TiO_5, $E4_1$) and Ta_2NiS_5 have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

4. β-$LiIO_3$: ABC3_tP40_86_g_g_3g 822

***I*4/*m* (87)** .....

1. Marialite Scapolite [$Na_4Cl(AlSi_3)_3O_{24}$, $S6_4$]: AB4C24D12_tI82_87_a_h_2h2i_hi 825
2. Sr_2NiWO_6: AB6C2D_tI20_87_a_eh_d_b 828

***I*4$_1$/*a* (88)** .....

1. $Na_4Ge_9O_{20}$: A9B4C20_tI132_88_a2f_f_5f 830
2. Copper (I) Azide (CuN_3): AB3_tI32_88_d_cf 835
3. Scheelite ($CaWO_4$, $H0_4$): AB4C_tI24_88_b_f_a .. 837

***P*42$_1$2 (90)** .....

1. $G7_5$ ($PbCO_3 \cdot PbCl_2$, Phosgenite) (*obsolete*): AB2C3D2_tP16_90_c_f_ce_e 839

***P*4$_1$2$_1$2 (92)** .....

1. Retgersite (α-$NiSO_4{\cdot}6H_2O$, $H4_5$): A12BC10D_tP96_92_6b_a_5b_a 841
2. Paratellurite (α-TeO_2): A2B_tP12_92_b_a 846

***I*4$_1$22 (98)** .....

1. Phase III $Cd_2Re_2O_7$: A2B7C2_tI44_98_f_bcde_f 848

***P*4*bm* (100)** .....

1. $F5_4$ (NH_4ClO_2) (*obsolete*): ABC2_tP8_100_b_a_c 851
2. NH_4NO_3 II ($G0_9$): ABC3_tP10_100_b_a_bc 853

***P*4*cc* (103)** .....

1. VSe_2O_6: A6B2C_tP72_103_abc5d_2d_abc 855

***I*4*mm* (107)** .....

1. BaNiSn$_3$: ABC3_tI10_107_a_a_ab 859

***P*$\bar{4}$2*m* (111)** .....

1. $E3_1$ (β-Ag_2HgI_4) (*obsolete*): A2BC4_tP7_111_f_a_n 861

***P*$\bar{4}$2$_1$*m* (113)** .....

1. Ammonium Chlorite (NH_4ClO_2): AB4CD2_tP16_113_c_f_a_e 863

***P*$\bar{4}$2$_1$*c* (114)** .....

1. $C_{19}Sc_{15}$: A19B15_tP68_114_bc4e_ac3e 865
2. $Ag_2SO_4{\cdot}4NH_3$ ($H4_{17}$): A2B12C4D4E_tP46_114_d_3e_e_e_a 869

***P*$\bar{4}$*m*2 (115)** .....

1. $F6_1$ (Chalcopyrite, $CuFeS_2$) (*obsolete*): ABC2_tP4_115_a_c_g 872

***I*$\bar{4}$*m*2 (119)** .....

1. Phase II $Cd_2Re_2O_7$: A2B7C2_tI44_119_i_bdefgh_i 874
2. Tetragonal $TlFeS_2$: AB2C_tI8_119_c_e_a 877

***I*$\bar{4}$*c*2 (120)** .....

1. $BeSO_4{\cdot}4H_2O$ ($H4_3$): AB8C8D_tI72_120_c_2i_2i_b 879

***I*$\bar{4}$2*m* (121)** .....

1. $SrCu_2(BO_3)_2$: A2B2C6D_tI44_121_i_i_ij_c 882
2. $C17$ (Fe_2B) (*obsolete*): AB2_tI12_121_ab_i 884
3. K_3CrO_8: AB3C8_tI24_121_a_bd_2i 886
4. α-V_3S: AB3_tI32_121_g_f2i 888

***I*$\bar{4}$2*d* (122)** .....

1. Mercury Cyanide [$Hg(CN)_2$, $F1_1$]: A2BC2_tI40_122_e_d_e 890
2. KH_2PO_4 ($H2_2$): A4BC4D_tI40_122_e_b_e_a 893
3. $NH_4H_2PO_4$: A8BC4D_tI56_122_2e_b_e_a 896
4. NaS_2: AB2_tI48_122_cd_2e 899

***P*4/*mmm* (123)** .....

1. NH_4HgCl_3 ($E2_5$): A3BC_tP5_123_cg_a_d 902
2. K_2PtCl_4 ($H1_5$): A4B2C_tP7_123_j_e_a 904
3. $E6_1$ ($Sr(OH)_2(H_2O)_8$) (*obsolete*): A8B2C_tP11_123_r_f_a 906
4. $E6_2$ [$SrO_2(H_2O)_8$] (*possibly obsolete*): A8B2C_tP11_123_r_h_a 908
5. $TlAlF_4$ ($H0_8$): AB4C_tP6_123_d_eh_a 910
6. δ-CuTi ($L2_a$): AB_tP2_123_a_d 912

***P*4/*mcc* (124)** .....

1. $CaO_2(H_2O)_8$: AB8C2_tP22_124_a_n_h 914

***P*4/*nnc* (126)** .....

1. Vesuvianite ($Ca_{10}Al_4(Mg,Fe)_2Si_9O_{34}(OH)_4$, $S2_3$): A4B10C2D34E4F9_tP252_126_k_ce2k_f_h8k_k_d2k 916
2. $Ag[Co(NH_3)_2(NO_2)_4]$ ($J1_9$): ABC4D2E8_tP32_126_a_b_h_e_k 926

***P*4/*mbm* (127)** .....

1. $Pd(NH_3)_4Cl_2{\cdot}H_2O$ ($H4_9$): A2BC4D_tP16_127_h_d_i_a 929
2. Phosgenite [$Pb_2Cl_2(CO_3)$]: AB2C3D2_tP32_127_g_eh_gk_k 931

***P*4/*mnc* (128)** .....

1. Chiolite ($Na_5Al_3F_{14}$, $K7_5$): A3B14C5_tP44_128_ac_ehi_bg 934
2. Apophyllite ($KCa_4Si_8O_{20}F{\cdot}8H_2O$, $S5_2$): A4BC16DE28F8_tP116_128_h_a_2i_b_g3i_i 937

***P*4/*nmm* (129)** .....

1. $CaBe_2Ge_2$: A2BC2_tP10_129_ac_c_bc 943
2. Meta-autunite (I) [$Ca(UO_2)_2(PO_4)_2{\cdot}6H_2O$, $H5_{10}$]: AB4C6DE_tP26_129_c_j_2ci_a_c 945
3. NH_4Br ($B25$): AB4C_tP12_129_c_i_a 948
4. LaOAgS: ABCD_tP8_129_b_c_a_c 950

***P*4/*ncc* (130)** .....

1. $Sr(OH)_2(H_2O)_8$: A18B10C_tP116_130_2c4g_2c2g_a 952
2. α-WO_3: A3B_tP16_130_cf_c 957

***P*4$_2$/*mnm* (136)** .....

1. Zr_3Al_2: A2B3_tP20_136_j_dfg 960
2. $ZrFe_4Si_2$: A4B2C_tP14_136_i_g_b 962
3. $K_2CuCl_4{\cdot}2H_2O$ ($H4_1$): A4BC4D2E2_tP26_136_fg_a_j_d_e 964
4. $Nd_2Fe_{14}B$: AB14C2_tP68_136_f_ce2j2k_fg 966

***I*4/*mmm* (139)** .....

1. $Sr_4Ti_3O_{10}$: A10B4C3_tI34_139_c2eg_2e_ae 970

2. $TlCo_2S_2$[∂‡‡]: A2B2C_tI10_139_d_e_a 972
3. $ThCr_2Si_2$[∂‡‡]: A2B2C_tI10_139_d_e_a 974
4. Au_2Nb_3: A2B3_tI10_139_e_ae 976
5. CaC_2-I ($C11_a$): A2B_tI6_139_e_a 978
6. $K_2OsO_2Cl_4$ ($J1_5$): A4B2C2D_tI18_139_h_d_e_a .. 980
7. K_2NiF_4: A4B2C_tI14_139_ce_e_a 982
8. $K_3TlCl_6{\cdot}2H_2O$ ($J3_1$): A6B2C3D_tI168_139_egikl2m_ejn_bh2n_acf 984
9. $Sr_3Ti_2O_7$: A7B3C2_tI24_139_aeg_be_e 988
10. Fe_8N ($D2_g$): A8B_tI18_139_deh_a 990
11. Li_2CN_2[⊗‡‡]: AB2C2_tI10_139_a_d_e 992
12. $H5_9$ [Autunite, $Ca(UO_2)_2(PO_4)_2{\cdot}10\frac{1}{2}H_2O$] (*obsolete*)[⊗‡‡]: AB2C2_tI10_139_a_d_e 994
13. $AuCsCl_3$ ($K7_6$): AB3C_tI20_139_ab_eh_d 996
14. "Martensite Type" FeC_x ($x \leq 0.06$) ($L'2_0$): AB_tI4_139_b_a 998

***I4/mcm* (140)** .....
1. V_4SiSb_2: A2BC4_tI28_140_h_a_k 1000
2. KHF_2 ($F5_2$): A2BC_tI16_140_h_d_a 1002
3. $Pu_{31}Rh_{20}$: A31B20_tI204_140_b2gh3m_ac2fh3l .. 1004
4. $NH_4Pb_2Br_5$ ($K3_4$): A5BC2_tI32_140_bl_a_h 1010
5. Cs_3CoCl_5 ($K3_1$): A5BC3_tI36_140_cl_b_ah 1012
6. U_6Mn ($D2_c$): AB6_tI28_140_a_hk 1014

***I4₁/amd* (141)** .....
1. $BaCd_{11}$: AB11_tI48_141_a_bdi 1016

***I4₁/acd* (142)** .....
1. Analcime ($NaAlSi_2O_6{\cdot}H_2O$, $S6_1$): A2B2C3D12E4_tI184_142_f_f_be_3g_g 1019
2. Cd_3As_2: A2B3_tI160_142_deg_3g 1025

***P3* (143)** .....
1. La_3BWO_9 ($P3$): AB3C9D_hP28_143_2a_2d_6d_bc 1030

***P3₁* (144)** .....
1. $RbNO_3$ (IV): AB3C_hP45_144_3a_9a_3a 1033

***P$\bar{3}$* (147)** .....
1. Na_2SO_3 ($G3_2$): A2B3C_hP12_147_abd_g_d 1037

***R$\bar{3}$* (148)** .....
1. Dolomite [$MgCa(CO_3)_2$, $G1_1$]: A2BCD6_hR10_148_c_a_b_f 1039
2. $K_2Sn(OH)_6$ ($H6_2$): A6B2C6D_hR15_148_f_c_f_a 1041
3. $Ni(H_2O)_6SnCl_6$ ($I6_1$): A6B6CD_hR14_148_f_f_b_a 1044
4. Li_7TaO_6: A8B6C_hR15_148_cf_f_a 1047

***P321* (150)** .....
1. $SrCl_2{\cdot}(H_2O)_6$: A2B12C6D_hP21_150_d_2g_ef_a 1049
2. $Cs_3As_2Cl_9$ ($K7_3$): A2B9C3_hP14_150_d_eg_ad .. 1052
3. Paralstonite ($BaCa(CO_3)_2$): AB2CD6_hP30_150_e_c2d_f_3g 1054
4. KSO_3 ($K1_1$): AB3C_hP30_150_ef_3g_c2d 1057
5. Steklite [$KAl(SO_4)_2$, $H3_2$]: ABC8D2_hP12_150_b_a_dg_d 1060

***R32* (155)** .....
1. $KBe_2BO_3F_2$: AB2C2DE3_hR9_155_b_c_c_a_e .. 1063

***R3m* (160)** .....
1. SbI_3S_{24}: A3B24C_hR28_160_b_2b3c_a 1065
2. Fe_3PO_7: A3B7C_hR11_160_b_a2b_a 1068
3. Cronstedtite {$Fe(Fe,Si)[(OH)_2,O]O_3$, $S5_7$}: AB3C2D_hR7_160_a_b_2a_a 1070
4. Low-Temperature $GaMo_4S_8$: AB4C8_hR13_160_a_ab_2a2b 1072
5. $KBrO_3$ ($G0_7$)[□]: ABC3_hR5_160_a_a_b 1074
6. γ-Potassium Nitrate (KNO_3)[□]: ABC3_hR5_160_a_a_b 1076

***R3c* (161)** .....
1. α-BaB_2O_4 (Low-Temperature): A2BC4_hR42_161_2b_b_4b 1078

***P$\bar{3}$1m* (162)** .....
1. $I1_3$ ($SrCl_2{\cdot}(H_2O)_6$) (*obsolete*)[§§]: A2B6C_hP9_162_d_k_a 1083
2. Rosiaite ($PbSb_2O_6$)[§§]: A6BC2_hP9_162_k_a_d .. 1085

***P$\bar{3}$1c* (163)** .....
1. $NaSbF_4(OH)_2$ ($J1_{12}$): A6BC_hP16_163_i_b_c ... 1087
2. Colquiriite ($LiCaAlF_6$): ABC6D_hP18_163_d_b_i_c 1090

***P$\bar{3}$m1* (164)** .....
1. Predicted Li_2MgH_{16} 300 GPa: A16B2C_hP19_164_2d2i_d_b 1093
2. Ce_2O_2S[¶¶]: A2B2C_hP5_164_d_d_a 1095
3. Brucite [$Mg(OH)_2$][¶¶]: A2BC2_hP5_164_d_a_d .. 1097

[∂] $TlCo_2S_2$ and $ThCr_2Si_2$ have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[‡‡] $TlCo_2S_2$, $ThCr_2Si_2$, Li_2CN_2, and $H5_9$ [Autunite, $Ca(UO_2)_2(PO_4)_2{\cdot}10\frac{1}{2}H_2O$] (*obsolete*) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

[⊗] Li_2CN_2 and $H5_9$ [Autunite, $Ca(UO_2)_2(PO_4)_2{\cdot}10\frac{1}{2}H_2O$] (*obsolete*) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[□] $KBrO_3$ ($G0_7$) and γ-Potassium Nitrate (KNO_3) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[§§] $I1_3$ ($SrCl_2{\cdot}(H_2O)_6$) (*obsolete*) and Rosiaite ($PbSb_2O_6$) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

[¶¶] Ce_2O_2S and Brucite [$Mg(OH)_2$] have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

4. $K_2Pt(SCN)_6$ ($H6_3$)[××]: A2BC6_hP9_164_d_a_i .. 1099
5. Bararite (Trigonal $(NH_4)_2SiF_6$, $J1_6$)[××]: A6B2C_hP9_164_i_d_a 1101
6. K_2GeF_6 ($J1_{13}$)[××]: A6BC2_hP9_164_i_a_d 1103
7. Jacutingaite (Pt_2HgSe_3): AB2C3_hP12_164_d_ae_i 1105
8. $D0_{13}$ ($AlCl_3$) (*obsolete*): AB3_hP4_164_b_ad 1107
9. Nevskite (BiSe): AB_hP12_164_c2d_c2d 1109

***R*$\bar{3}$*m* (166)**
1. $B_{13}C_2$ "B_4C" ($D1_g$): A13B2_hR15_166_b2h_c .. 1111
2. $MnBi_2Te_4$[°°]: A2BC4_hR7_166_c_a_2c 1114
3. Shandite ($Ni_3Pb_2S_2$): A3B2C2_hR7_166_d_ab_c 1116
4. Chabazite ($Ca_{1.4}Sr_{0.3}Al_{3.8}Si_{8.3}O_{24}{\cdot}13H_2O$, $S3_4$ (I)): A5B21C24D12_hR62_166_a2c_ehi_fg2h_i 1118
5. $CaSi_2$ ($C12$): AB2_hR6_166_c_2c 1124
6. Rhombohedral $CuTi_2S_4$: AB4C2_hR28_166_2c_2c2h_abh 1126
7. $CaCu_4P_2$[°°]: AB4C2_hR7_166_a_2c_c 1129
8. $CaUO_4$: AB4C_hR6_166_b_2c_a 1131
9. $TaTi_7$ (BCC SQS-16): AB7_hR16_166_c_c2h ... 1133
10. Rhombohedral Delafossite ($CuFeO_2$): ABC2_hR4_166_a_b_c 1136
11. β-Potassium Nitrate (KNO_3): ABC6_hR8_166_a_b_h 1138
12. K(SH) ($B22$): AB_hR2_166_a_b 1140

***R*$\bar{3}$*c* (167)**
1. $Zr_{21}Re_{25}$: A25B21_hR92_167_b2e3f_e3f 1142
2. β-BaB_2O_4 (High-Temperature): A2BC4_hR42_167_f_ac_2f 1151
3. $CrCl_3(H_2O)_6$ ($J2_2$): A3BC6_hR20_167_e_b_f ... 1156
4. FeF_3 ($D0_{12}$): A3B_hR8_167_e_b 1159
5. Rinneite ($K_3NaFeCl_6$): A6BC3D_hR22_167_f_b_e_a 1161
6. $Cs_3Tl_2Cl_9$ ($K7_2$): A9B3C2_hR28_167_ef_e_c 1164

***P*6$_3$ (173)**
1. Crancrinite ($Na_6Ca_2Al_6Si_6O_{24}(CO_3)_2$, $S3_3$ (I)): A3BCD3E15F3_hP52_173_c_b_b_c_5c_c 1168
2. La_3CuSiS_7: AB3C7D_hP24_173_a_c_b2c_b 1172
3. La_3BWO_9 ($P6_3$): AB3C9D_hP28_173_a_c_3c_b . 1175
4. α-$LiIO_3$: ABC3_hP10_173_b_a_c 1178
5. $LiKSO_4$ ($H1_4$): ABC4D_hP14_173_a_b_bc_b 1180

***P*6$_3$/*m* (176)**
1. $Rh_{20}Si_{13}$: A10B7_hP34_176_c3h_b2h 1182
2. Th_7S_{12} ($D8_k$): A3B2_hP20_176_2h_ah 1185
3. β-Si_3N_4: A4B3_hP14_176_ch_h 1187
4. Fluorapatite [$Ca_5F(PO_4)_3$, $H5_7$]: A5BC12D3_hP42_176_fh_a_2hi_h 1189
5. $Fe_2(CO)_9$ ($F4_1$): A9B2C9_hP40_176_hi_f_hi 1192
6. $K_3W_2Cl_9$ ($K7_1$): A9B3C2_hP28_176_hi_af_f 1195

***P*6$_2$22 (180)**
1. Hg_2O_2NaI: A2BCD2_hP18_180_f_c_b_i 1198

***P*6$_3$22 (182)**
1. $BaAl_2O_4$ ($H2_8$): A2BC6_hP18_182_f_b_gh 1200
2. $E2_3$ ($LiIO_3$) (*obsolete*): ABC3_hP10_182_c_b_g 1202

***P*6$_3$*mc* (186)**
1. $Zn_2Mo_3O_8$: A3B8C2_hP26_186_c_ab2c_2b 1204
2. $Nd(BrO_3)_3{\cdot}9H_2O$ ($G2_2$): A3B9CD9_hP44_186_c_3c_b_cd 1207
3. Swedenborgite ($NaBe_4SbO_7$, $E9_2$): A4BC7D_hP26_186_ac_b_a2c_b 1210
4. $C27$ (CdI_2) (*questionable*): AB2_hP6_186_b_ab 1213
5. $LiClO_4{\cdot}3H_2O$ ($H4_{18}$): AB6CD7_hP30_186_b_d_a_b2c 1215
6. Cd(OH)Cl ($E0_3$): ABCD_hP8_186_b_b_a_a 1218

***P*$\bar{6}$*m*2 (187)**
1. Cr-233 Quasi-One-Dimensional Superconductor ($K_2Cr_3As_3$): A3B3C2_hP16_187_jk_jk_ck 1220
2. Cs_7O: A7B_hP24_187_ai2j2kn_j 1222

***P*$\bar{6}$2*m* (189)**
1. ZrNiAl: ABC_hP9_189_g_ad_f 1225

***P*$\bar{6}$2*c* (190)**
1. $CsSO_3$ ($K1_2$): AB3C_hP20_190_ac_i_f 1227
2. Bastnäsite [$CeF(CO_3)$]: ABCD3_hP36_190_h_g_af_hi 1230

***P*6/*mmm* (191)**
1. $TiBe_{12}$ (approximate, $D2_a$): A12B_hP13_191_cdei_a 1233
2. Hexagonal WO_3: A3B_hP12_191_gl_f 1235

***P*6$_3$/*mcm* (193)**
1. $D0_6$ (Tysonite, LaF_3) (*obsolete*): A3B_hP24_193_ack_g 1238
2. Ti_5Ga_4: A4B5_hP18_193_bg_dg 1241

***P*6$_3$/*mmc* (194)**
1. Proposed 300 GPa HfH_{10}: A10B_hP22_194_bhj_c 1243
2. Magnetoplumbite ($PbFe_{12}O_{19}$): A12B19C_hP64_194_ab2fk_efh2k_d 1246
3. Pt_2Sn_3 ($D5_b$): A2B3_hP10_194_f_bf 1249
4. β-Alumina (Al_2O_3, $D5_6$): A2B3_hP60_194_3fk_cdef2k 1251

[××] $K_2Pt(SCN)_6$ ($H6_3$), Bararite (Trigonal $(NH_4)_2SiF_6$, $J1_6$), and K_2GeF_6 ($J1_{13}$) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

[°°] $MnBi_2Te_4$ and $CaCu_4P_2$ have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

5. $S3_4$ (II) (Catapleiite, $Na_2Zr(SiO_3)_3 \cdot H_2O$) (*obsolete*): A3B2C9D3E_hP36_194_g_f_hk_h_a 1254
6. ReB_3: A3B_hP8_194_af_c 1257
7. $Cs_3Cr_2Cl_9$: A9B2C3_hP28_194_hk_f_bf 1259
8. $Na_{0.74}CoO_2$: AB2C2_hP10_194_a_bc_f 1261
9. $EuIn_2P_2$: AB2C2_hP10_194_a_f_f 1263
10. Lu_2CoGa_3: AB3C2_hP24_194_f_k_bh 1265
11. Hexagonal Delafossite ($CuAlO_2$): ABC2_hP8_194_a_c_f 1268
12. $LiZn_2$ (C_k)[∂∂]: AB_hP4_194_a_c 1270
13. Fe_2N (approximate, $L'3_0$)[∂∂]: AB_hP4_194_c_a ..1272

***P*2$_1$3 (198)** .....
1. Cubic Cu_2OSeO_3: A2B4C_cP56_198_ab_2a2b_2a 1274
2. Na_2CaSiO_4 ($S6_6$): AB2C4D_cP32_198_a_2a_ab_a 1277
3. α-Carnegieite ($NaAlSiO_4$, $S6_5$): ABC4D_cP28_198_a_a_ab_a 1280

***I*2$_1$3 (199)** .....
1. $C26_a$ (NO_2) (*obsolete*): AB2_cI36_199_b_c 1282

***Pn*$\bar{3}$ (201)** .....
1. $Bi_3Ru_3O_{11}$: A3B11C3_cP68_201_be_efh_g 1285

***Fm*$\bar{3}$ (202)** .....
1. $K_3Co(NO_2)_6$ ($J2_4$): AB3C6D12_cF88_202_a_bc_e_h 1289

***Im*$\bar{3}$ (204)** .....
1. $LaFe_4P_{12}$: A4BC12_cI34_204_c_a_g 1292
2. $NaMn_7O_{12}$: A7BC12_cI40_204_bc_a_g 1294
3. NO_2 (Modern, $C26$): AB2_cI36_204_d_g 1296

***Pa*$\bar{3}$ (205)** .....
1. $Zn(BrO_3)_2 \cdot 6H_2O$ ($J1_{10}$)[⊗⊗]: A2B6C6D_cP60_205_c_d_d_a 1299
2. $H6_4$ [$Ni(NO_3)_2(NH_3)_6$] (*obsolete*)[⊗⊗]: A2B6CD6_cP60_205_c_d_a_d 1302
3. $Pb(NO_3)_2$ ($G2_1$): A2B6C_cP36_205_c_d_a 1306
4. CaB_2O_4 (IV): A2BC4_cP84_205_d_ac_2d 1309
5. SnI_4 ($D1_1$): A4B_cP40_205_cd_c 1313
6. $NaSbF_6$: A6BC_cP32_205_d_b_a 1316
7. ZrP_2O_7 High-Temperature ($K6_1$): A7B2C_cP40_205_bd_c_a 1319
8. $NaCr(SO_4)_2 \cdot 12H_2O$ Alum: AB12CD8E2_cP96_205_a_2d_b_cd_c 1322
9. γ-Alum [$AlNa(SO_4)_2 \cdot 12H_2O$, $H4_{15}$]: AB24CD20E2_cP192_205_a_4d_b_c3d_c 1327
10. α-Alum [$KAl(SO_4)_2 \cdot 12H_2O$, $H4_{13}$]: AB24CD28E2_cP224_205_a_4d_b_2c4d_c 1335
11. β-Alum [$Al(NH_3CH_3)_2(SO_4)_2 \cdot 12H_2O$, $H4_{14}$]: AB2C36D2E20F2_cP252_205_a_c_6d_c_c3d_c ... 1344

***P*4$_3$32 (212)** .....
1. Maghemite (γ-Fe_2O_3, $D5_7$): A2B3_cP60_212_bcd_ace 1354

***P*4$_1$32 (213)** .....
1. Al_2Mo_3C: A2BC3_cP24_213_c_a_d 1358
2. Mg_3Ru_2: A3B2_cP20_213_d_c 1361

***F*$\bar{4}$3*m* (216)** .....
1. Zunyite [$Al_{13}(OH,F)_{18}Si_5O_{20}Cl$, $S0_8$]: A13BC18D20E5_cF228_216_dh_b_fh_2eh_ce 1363
2. Murataite [$(Y,Na)_6(Zn,Fe)_5Ti_{12}O_{29}(O,F)_{10}F_4$]: A16B40C12D6E5_cF316_216_eh_e2g2h_h_f_be .. 1367
3. $Sm_{11}Cd_{45}$: A45B11_cF448_216_bd4efg5h_ac2eh . 1371
4. Hg_2TiCu Inverse Heusler: AB2C_cF16_216_b_ad_c 1376
5. $GaMo_4S_8$: AB4C8_cF52_216_a_e_2e 1378
6. High-Temperature Cubic $KClO_4$ ($H0_5$): ABC4_cF24_216_b_a_e 1380
7. AlN (cF40): AB_cF40_216_de_ce 1382

***I*$\bar{4}$3*m* (217)** .....
1. Tennantite ($Cu_{12}As_4S_{13}$): A4B24C13_cI82_217_c_deg_ag 1384
2. AlN (cI16): AB_cI16_217_c_c 1387

***P*$\bar{4}$3*n* (218)** .....
1. Hauyne [$(Na_{0.5}Ca_{0.3}K_{0.2})_8(Al_6Si_6O_{24})(SO_4)_{1.5}$, $S6_9$]: A3B4C4D4E16F4G3_cP76_218_c_e_e_e_ei_e_d .. 1389
2. Sodalite [$Na_4(AlSiO_4)_3Cl$, $S6_2$]: A3BC4D12E3_cP46_218_d_a_e_i_c 1394

***I*$\bar{4}$3*d* (220)** .....
1. Eulytine ($Bi_4(SiO_4)_3$, $S1_5$): A4B12C3_cI76_220_c_e_a 1397
2. Mayenite ($12CaO \cdot 7Al_2O_3$, $K7_4$, C12A7): A7B12C19_cI152_220_bc_2d_ace 1400
3. $Al(PO_3)_3$ ($G5_2$): AB9C3_cI208_220_c_3e_e 1405
4. AlN (cI24): AB_cI24_220_a_b 1412

***Pm*$\bar{3}$*m* (221)** .....
1. γ-Fe_4N ($L'1_0$): A4B_cP5_221_bc_a 1414
2. Predicted High-Pressure $YCaH_{12}$: AB12C_cP14_221_a_h_b 1416
3. NH_4NO_3 I ($G0_8$): AB_cP2_221_a_b 1418

***Pn*$\bar{3}$*m* (224)** .....
1. Dodecatungstophosphoric Acid Hexahydrate [$H_3PW_{12}O_{40} \cdot 6H_2O$]: A27B52CD12_cP184_224_dl_eh3k_a_k 1420
2. Mg_3P_2 ($D5_5$): A3B2_cP10_224_d_b 1427
3. $H_3PW_{12}O_{40} \cdot 3H_2O$: A3B40CD12_cP112_224_d_e3k_a_k 1429
4. 12-phosphotungstic acid [$H_3PW_{12}O_{40} \cdot 5H_2O$ ($H4_{16}$)]: A5B40CD12_cP116_224_cd_e3k_a_k 1434

***Fm*$\bar{3}$*m* (225)** .....

[∂∂] $LiZn_2$ (C_k) and Fe_2N (approximate, $L'3_0$) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

[⊗⊗] $Zn(BrO_3)_2 \cdot 6H_2O$ ($J1_{10}$) and $H6_4$ [$Ni(NO_3)_2(NH_3)_6$] (*obsolete*) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

1. LaH_{10} High-T_c Superconductor: A10B_cF44_225_cf_b 1439
2. Double Perovskite (Ba_2MnWO_6): A2BC6D_cF40_225_c_a_e_b 1441
3. $Cu_3[Fe(CN)_6]_2{\cdot}xH_2O$ ($J2_5$, $x \approx 3$): A6B9CD2E6_cF96_225_e_bf_a_c_e 1443
4. Co_9S_8 ($D8_9$): A9B8_cF68_225_af_ce 1446
5. $(NH_4)_3AlF_6$ ($J2_1$): AB30C16D3_cF200_225_a_ej_2f_bc 1448
6. $L1_a$ (disputed $CuPt_3$): AB7_cF32_225_b_ad 1451
7. Sulphohalite [$Na_6ClF(SO_4)_2$, $H5_8$]: ABC6D8E2_cF72_225_b_a_e_f_c 1453

***Fd$\bar{3}$m* (227) .....**

1. γ-Ga_2O_3: A11B4_cF120_227_acdf_e 1455
2. Predicted Li_2MgH_{16} High-Temperature Superconductor (250 GPa): A16B2C_cF152_227_eg_d_a 1458
3. $Mg_3Cr_2Al_{18}$: A18B2C3_cF184_227_fg_d_ac 1461
4. $Zn_{22}Zr$: A22B_cF184_227_cdfg_a 1464
5. $H_3PW_{12}O_{40}{\cdot}29H_2O$ ($H4_{21}$): A29B40CD12_cF656_227_ae2fg_e3g_b_g 1467
6. $G7_3$ [Northupite, $Na_3MgCl(CO_3)_2$] (*obsolete*): A2BCD3E6_cF208_227_e_c_d_f_g 1475
7. $D6_2$ (Sb_2O_4) (*obsolete*): A2B_cF96_227_abf_cd 1478
8. Senarmontite (Sb_2O_3, $D6_1$): A3B2_cF80_227_f_e 1481
9. $H5_6$ [Tychite, $Na_6Mg_2SO_4(CO_3)_4$)] (*obsolete*): A4B2C6D16E_cF232_227_e_d_f_eg_a 1484
10. Cubic CuPt ($L1_3$ (I), $D4$): AB_cF32_227_c_d ... 1488

***Im$\bar{3}$m* (229) .....**

1. α-AgI ($B23$): A21B_cI44_229_bdh_a 1490

***Ia$\bar{3}$d* (230) .....**

1. $Ca_3Al_2(OH)_{12}$ ($J2_3$): A2B3C12D12_cI232_230_a_c_h_h 1492

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1. Introduction

Crystal structure classification began near the dawn of X-ray Crystallography with early works by Wyckoff [1], Ewald's 1927 *Handbuch der Physik* article [2], and the Landolt-Börnstein 1927 supplement [3]. In the next few decades, these catalogs proliferated, including the *Strukturbericht* (Structure Reports) series [4], Pearson's *Handbook of Lattice Spacings and Structures of Metals and Alloys* [5, 6], Smithells's *Metals Reference Book* [7, 8], and Wyckoff's updated *Structure of Crystals* [9]. In later years, *Pearson's Handbook* [10] and *Smithells Metals Reference Book* [11] acquired new editors and even today produce periodic editions. Subsequent works, such as the *Gmelin Handbook* [12], also appear regularly. As online access became the norm at the beginning of this century, electronic resources appeared, including the *American Mineralogist Crystal Structure Database* [13], which compiles the structures of minerals from a variety of sources, and the SpringerMaterials website [14], an extension of *Pearson's Handbook* that also incorporates the Landolt-Börnstein series.

Computational material-property databases have also become available: for example AFLOW (Automatic FLOW for Materials Discovery) [15–17], Novel Materials Discovery (NoMaD) [18], the Materials Project [19], and the Open Quantum Materials Database (OQMD) [20].

While these print and electronic databases contain a vast number of crystal structures, many are not optimized for use in materials computations, *e.g.*, by providing structural information in a format which can be read by one of the common density functional computer codes. In addition, while most of these resources make their data available in standard formats, such as space group/Wyckoff positions or a Crystallographic Information File (CIF) [21], these forms still pose a challenge for novice researchers. Therefore, it is desirable to have a database that addresses both of these concerns.

The AFLOW Library of Crystallographic Prototypes (hereafter AFLOW Prototype Encyclopedia or Prototype Encyclopedia for brevity) [22, 23] has been designed with these objectives in mind. Integrated with the AFLOW computational framework, it can provide structural input for a wide variety of materials computational programs, including VASP [24], Quantum ESPRESSO [25], FHI-AIMS [26], ABINIT [27], and ELK [28] as well as in CIF format. Each structure's web page provides all the necessary crystallographic information, including space group, lattice vectors, and Wyckoff positions. In addition, the formulas for the primitive vectors and basis vectors of the structure are shown. This is particularly useful for students and novice researchers, as it shows the relationship between the standard crystallographic coordinates needed for computational work and the Cartesian positions of the atom in the primitive and/or conventional unit cell. If the desired chemical com-

position for the structure is not the prototype, the chemistry and primitive cell dimensions can be changed by the user before generating the structural input. Finally, the structure may be viewed as a primitive cell, conventional cell, or supercell at a variety of angles and magnifications via JSmol [29].

This article is about the ongoing development of the AFLOW Prototype Encyclopedia. In Part 1 [22], we introduced the Prototype Encyclopedia, which began as an updated version of the now-defunct Naval Research Laboratory "*Structure of Crystals*" web site. This was a collection of crystal structures that we found useful as a starting point for computational studies of materials. One feature of the site was that it included a number of well-known *Strukturbericht* designations, from the German effort to classify common crystal structures [4]. This article also provided a brief introduction of the crystallography needed to understand crystal structures, as well as a description of the entries in the AFLOW Prototype Encyclopedia for each structure. In all, 288 structures were included in the initial Prototype Encyclopedia.

Part 2 [23] continued the discussion of crystallography by reviewing the enantiomorphic space groups, which are pairs of space groups that are mirror images of each other. As part of our teaching mission, we also described the Wigner-Seitz version of the primitive unit cell of a crystal, and discussed the 17 two-dimensional "plane groups," the analog of the 230 three-dimensional space groups. Another 302 structures were added to the database to bring the total to 590, with at least one structure shown in each of the 230 space groups.

In this article, Part 3, we add 510 new structures to the encyclopedia. These include all of the inorganic *Strukturbericht* designations which had not previously been cataloged by us. The *Strukturbericht* classification scheme was one of the earliest attempts to group elements[1] and compounds into common crystal structures. However, despite its historical importance there is no available history of its development, and until now there has been no comprehensive index of all the *Strukturbericht* designations that were defined in the series or added by others.

Recently, M. J. Mehl published a brief history of the development of the *Strukturbericht* designations [31]. In Part 3, we improve and extend that work, beginning with the introduction of these labels by Ewald [2] through the publication of the seven *Strukturbericht* volumes and the post-World War II extensions to the scheme made by Smithells [8], Pearson [5], and others. The *Strukturbericht* classifications changed over the run of the series, and several structures were published under multiple designations. Here, we track all of those changes.

The article is organized as follows: in Sections 2 and 3, we present common prototype classification schemes, followed by a short history of the development and presentation of *Strukturbericht* designations. As the *Strukturbericht* designations changed over time, Section 4 discusses the history of the major categories and sub-categories, as well as our editorial decisions as to which label should be used for a given prototype. Section 5 presents what we believe to be the first comprehensive list of *Strukturbericht* designations for elements and inorganic compounds, including all references to labels which we could find in the literature, with a link to the corresponding online pages in the AFLOW Prototype Encyclopedia. Section 6 describes AFLOW software modules that will be used to identify and classify new prototypes for future iterations of this work. Additionally, Section 7 introduces an online form for users to suggest new prototypes to enrich the Encyclopedia. Section 8 provides information about AFLOW Schools: hands-on workshops for students and researchers to learn about structural analysis methods — especially those discussed in Parts 1, 2, and 3 of this work — as well as other topics related to computational materials science. Section 9 summarizes our conclusions. This is followed by a list of all 510 new prototypes we have included in the Prototype Encyclopedia since the publication of Part 2.

2. Crystallographic Classification Schemes

Just before the First World War, Max von Laue suggested that diffraction of X-rays through crystals would produce a pattern which could be used to determine the crystal structure [32]. Shortly thereafter, this technique was also applied by Bragg and Bragg [33] and Hull [34] to analyze the structures of diamond and graphite, respectively. Over the next decade, the structures of hundreds of elements and compounds were determined [2, 3, 35, 36].

It was soon obvious that nature is in the habit of repeating itself. For example, the metallic elements mostly take close-packed face-centered-cubic (fcc) or hexagonal (hcp) structures, or the nearly close-packed body-centered cubic (bcc) structure. (The major exceptions being manganese, which takes on two unique atomic arrangements, and the actinides.) Of the remaining elements, the early group 14 all take on the diamond structure. We only see other patterns when we go to the far right-hand side of the periodic table.

Similar trends occur in compounds. Many AB type binaries have ground states in the sodium chloride, cesium chloride or zincblende structures, A_2B compounds can frequently have the positions found in fluorite (CaF_2), and metallic compounds of the form A_2BC often appear in the Heusler structure. This continues into more complicated chemistries. Data-mining sources such as AFLOW-ICSD

[1] Of interest is the work of Davey [30], who found that tungsten was too opaque for standard X-ray diffraction techniques, and so diluted his sample with ten parts wheat flour in order to get a diffraction pattern.

show that a large number of compounds exist in a relatively small number of crystal structures [37].

This duplication of structures leads to the assumption that most elements and compounds exist in a relatively small number of crystallographic forms, and that compounds having similar structures are somehow related to one another. If true, then it should be possible to classify crystal structures according to their periodicity and arrangement of atoms in the unit cell. We now show four common classification schemes.

The simplest crystal structure classification is by their space group and the occupied Wyckoff positions within that space group [38]. This is certainly efficient, and we have used a variant of it to index crystal structures [22, 23]. Unfortunately, this description is neither particularly informative nor unique. The ground states of gallium, iodine, and phosphorus are all in space group *Cmca* (#64), and all the atoms occupy a single ($8f$) Wyckoff position [39], but their structures are all different: the iodine atoms form dimers, while gallium and phosphorus have very different conventional unit cells.

A second way of defining classes of crystal structures is to group all similar elements or compounds into one class, and refer to it by one "prototype". Ionic compounds such as NaCl, MgO, BaS, and many others all take the same structure: fcc space group $Fm\bar{3}m$ (#225), with one atom on the ($4a$) Wyckoff site and the other on the ($4b$) site. This is universally known as the "sodium chloride" or "rock salt" structure [35]. This approach works well if there is an agreement on a single prototype for each structure, but can lead to confusion if different researchers use different prototypes: for example, ϵ-FeN_2 and β-VN_2 are both claimed as prototypes for the same structure [10, 40], while the mineral rosiaite ($PbSb_2O_6$), with similar lattice parameters and atomic coordinates [41], is considered the prototype of the ternary form.

A third method was introduced by Pearson [6]: specify the lattice type (cubic, hexagonal, tetragonal, etc.), followed by the number of atoms in the conventional unit cell. These *Pearson Symbols* are a compact version of the first method, and suffer from the same problems.

A fourth method is derived from the second: assign a label to each structure class and refer to it by that label. These came to be called *Strukturbericht* designations, for reasons discussed below. This scheme can prove unwieldy, as there are a potentially infinite number of possible structures, making it impossible for any labeling scheme to keep up, especially if there is no systematic method for choosing the labels. It does have the advantage that there is a unique label for every structure class, independent of chemical composition.

In practice, one identifies the structure using several of these methods. Of the four, the *Strukturbericht* scheme is the most obscure, and so we now turn our attention to it.

3. A Brief History of *Strukturbericht* Symbols [2]

This *Strukturbericht* scheme was introduced by Ewald in his review of X-ray crystallography for the 1927 edition of the *Handbuch der Physik* [2]. The 1927 supplement to the 1923 edition of Landolt-Börnstein [3] also contains a shortened version of Ewald's list, presumably taken from the *Handbuch*, but we have not been able to determine which one was actually published first.

The structural cataloging method devised by Ewald was rather simple:

- Purely elemental structures were given designations beginning with the letter A, followed by a number. Thus, the face-centered cubic lattice, prototype copper, was given the designation $A1$, body-centered cubic (tungsten) $A2$, hexagonal close-packed (magnesium) $A3$, diamond $A4$, and so on. Each new structure was labeled with the next available number, making the scheme rather arbitrary, as there may be no relation between structure An and $An + 1$.

- Binary compounds with composition AB were given designations starting with B: $B1$ (NaCl), $B2$ (CsCl), $B3$ (zincblende ZnS), $B4$ (wurtzite ZnS), *etc.* Thus, we can easily distinguish between the two forms of zinc sulfide.

- Binary compounds with composition AB_2 had designations starting with C, *e.g.*, fluorite (CaF_2) is $C1$.

- Binary compounds A_mB_n start with D. The original $D1$ structure was corundum (Al_2O_3).

- Ternary and higher compounds that do not contain radicals such as SiO_4 or OH have labels beginning with E. The only example was $E1$ ($KMgF_3$).

- The letter F was used for compounds with linear radicals, such as $F1$ [K(CN)].

- "Flat" (two dimensional) radicals start with G: $G1$, calcite ($CaCO_3$).

- "Complex" (three dimensional) radicals were assigned to H: $H1$, anhydrite ($CaSO_4$).

The *Strukturbericht* scheme was refined and extended by various literature reviews. The earliest crystallographic critical review that we have found was published by Wyckoff as part of his seminal 1924 work, *The Structure of Crystals* [1], followed by the 1931 second edition [35], the 1934 supplement edition [42], and culminating in the six-volume second edition of *The Structure of Crystals* from 1963 through 1971 [9].

[2] A somewhat different version of this section appeared in Ref. [31].

In 1931, Ewald and C. Hermann [36] published an exhaustive review of the structural literature from 1913-1928, "*Strukturbericht* (Structure Reports) 1913-1928" (we will refer to this as SB-I) as a supplement to the journal *Zeitschrift für Kristallographie*. In their introduction, they say that the purpose was (translated from German)

> ... to provide a compilation and critical review of *all* [our emphasis] previously published articles on X-ray structure determination. The structures are arranged according to their chemical formulas, at the same time a summary by "types" provides an overview over the known structures and their representatives. Unlike other published structural compositions, the present one stands out by discussing individual works, not just the results of the structures of the crystal types. It is expected that this will improve the accuracy of the structural determinations, the exact nature of the material used, the corrections to be made to older reports due to new experiences, and that other similar points can be assessed better.

The given dates show that this was intended to cover the field of structural studies from the beginning. This volume was primarily a critical review of the current X-ray crystallographic literature, but it also grouped crystal structures into types using the classification scheme defined in the *Handbuch*, leading to the structural prototypes becoming known as *Strukturbericht* designations.

The second (SB-II) and third (SB-III) volumes of *Strukturbericht* were published in 1937, encompassing the years 1928-1932 and 1933-1935, respectively. Each was designed to cover all published crystal structures, or at least all structures which the editors could obtain, for the indicated timeframe. Later volumes (SB-IV through SB-VII) span one year of research each, with the seventh and final volume, including 1939, published in 1943 during the Second World War. At the end of the war, the International Union of Crystallography (IUCr) continued this project, publishing it as *Structure Reports*, in English, through 1990. The first IUCr volume contained research published in 1945-1946 and was designated as volume 10 [43]. Volumes 8 and 9, covering 1940-1941 [44] and 1942-1944, respectively, were published later, providing continuity with the original *Strukturbericht*.

Volumes SB-I-VII modified and extended Ewald's classification scheme, but after the war, it was realized that the number of *Strukturbericht* symbols would keep growing uncontrollably and so the practice of giving new structures symbols was discontinued by the IUCr. A few new symbols were introduced by C. J. Smithells in his 1955 *Metals Reference Book* [8], and by W. B. Pearson in his 1958 *Handbook of Lattice Spacings and Structures of Metals and Alloys* [5]. Several other symbols have been added by other authors, noted in Section 5.

Even though the generation of new symbols ceased, the ones found in the *Strukturbericht* volumes, Smithells, and Pearson are still in use. For example, compounds with the $A15$ structure (Cr_3Si or β-W) became known as the first "high temperature" (20 K) superconductors [45]; and there are frequent references to "$B1$" (NaCl) or "$B2$" (CsCl) structures. Despite this, there is no complete index of all the *Strukturbericht* symbols: an index was published as part of the table of contents for SB-II, but this was not continued into the later volumes. In 1976, Trotter and Bree [46] published a complete index of all the compounds mentioned in all seven volumes, but it was by formula and by author, with no reference to the *Strukturbericht* designations. Smithells, Pearson, and other works such as Villars [10], and Parthé *et al.* [12] have published partial lists, but these are weighted toward the intermetallics and mostly ignore silicates (*Strukturbericht S*) and other non-metallic categories.

While we do not intend to write a complete history of *Strukturbericht* designations, it is worthwhile to give a brief description of the changes in the system that occurred throughout its run. The remainder of this section summarizes the changes to the system made in each volume. Following that, Section 4 delves more deeply into the changes made in each *Strukturbericht* category. Finally, we provide the first complete list of inorganic *Strukturbericht* designations in Section 5. As part of this effort we added all of the inorganic *Strukturbericht* designations as entries in the AFLOW Prototype Encyclopedia, leaving out the very few where the definition was so vague that we could not determine a structure.

3.1. *Strukturbericht* Band (Volume) I 1913-1928

SB-I [36] classified structures using a modified version of the scheme presented in Ewald's *Handbuch* article. This included the introduction of three new *Strukturbericht* categories:

- Type L, "alloys" (*Legierungen*): These are ordered structures based on the fcc and bcc lattices, such as Cu_3Au (*Strukturbericht* $L1_2$), which can be considered as an fcc lattice with an ordered substitution of gold atoms for copper, and the Heusler prototype $AlCu_2Mn$ ($L2_1$), a decoration of the bcc lattice.
- Type M, "solid solutions" (*Mischkristalle*). Presumably this was to be used for solid solutions, but no M-type structures were ever defined.
- Type O, "organic:" These are categories of organic crystals cataloged in a manner similar to the inorganic categories A-L. Type O structures were defined in every volume of *Strukturbericht*, but the labels never became common and we will not discuss them in this article.

In addition to the new categories, SB-I subdivided Ewald's original categories. For example, while type D still

represented binary compounds of the form A_mB_n, the type was divided into ranges, with $D1 - D10$ reserved for compounds AB_3, $D11 - D20$ for AB_n with $n > 3$, and finally, $D50 - D100$ for the more general compounds. Thus, the original $D1$, Al_2O_3, now became $D51$, while $D1$ was used for ammonia, NH_3. Similar divisions were made in the F, G, H and L categories.

The volume itself was divided by *Strukturbericht* category. Each section began with a description of the structures associated with each *Strukturbericht* designation and ended with a review of the literature for these and similar structures. These reviews would be present in all *Strukturbericht* volumes.

3.2. *Strukturbericht* Band II, 1928-1932

SB-II [47] was published in 1937 and covered the years 1928 through 1932. By then, Ewald, protesting the firing of all Jewish professors at the Technische Hochschule Stuttgart, resigned his post and emigrated to the United Kingdom [48], so the volume was edited by Carl Hermann, Otto Lohrmann, and Hans Philipp.

The new publication included a table of contents with the only index of *Strukturbericht* symbols ever published in the original volumes. Scheme changes developed in the *Handbuch* and SB-I were made. The most notable was in the format of the labels for classes D and above, due to the large number of new structures being found. This caused a problem in the D category, where designations $D1$ through $D10$ were reserved for structures with stoichiometry AB_3. By 1932, thirteen unique structures of this type were known, exceeding the category limit. To remedy this, the editors changed the SB-I $D1 - D100$ letter-number format to letter-number-subscript, where now the letter-number combination reveals the stoichiometry of the crystal and the subscript the specific structure. Thus, ammonia (NH_3) formerly $D1$, became $D0_1$; skutterudite, formerly $D2$, became $D0_2$. (Eventually the $D0$ range would extend to $D0_{24}$, Ni_3Ti.) This practice was applied to all of the categories beyond D.

SB-II also introduced the first fourteen entries in the rather general E category, including $E0_1$, PbFCl. Other E entries included chalcopyrite ($CuFeS_2$, $F6_1$), cubic perovskite ($CaTiO_3$, $G0_5$), and ilmenite ($FeTiO_3$, $G0_4$); were reclassified to $E1_1$, $E2_1$, and $E2_2$, respectively, presumably because they were insufficiently radical.

Other new categories were added:

- Type I, compounds with BC_6 radicals. These were mostly renamed from type H, where in SB-I $H50$ and above had been reserved for large radicals. Now, *e.g.*, K_2PtCl_6, originally $H6_1$, was reclassified as $I1_1$, though the SB-II table of contents listed it under both designations. Compounds with radicals of the form BC_m ($m < 6$), remained in the H category, although many SiO_5 silicates were co-listed in the new S category.
- Type K, "complex radicals," including $K0_1$ ($K_2S_2O_5$), $K1_1$ ($K_2S_2O_6$), and $K1_2$ ($Cs_2S_2O_6$). At least that is how they were listed in the text. In the SB-II table of contents, they were listed as $K1_1$, $K2_1$, and $K2_2$, respectively. (We use the names in the text rather than the names in the table of contents below.)
- Type S, the silicates, which were now moved to a new category. The silicates defined in SB-I had mostly been given H designations, but now were also listed under S, at least in the index of SB-II. Thus, kyanite (Al_2SiO_5, originally labeled $H5_1$ or $H51$, acquired the new label $S0_1$. In all, thirty-two compounds were identified, all but six of these new to SB-II.

3.3. *Strukturbericht* Band III, 1933-1935

SB-III was the last multi-year volume of the series, published in 1937 [49]. Edited by C. Gottfried and F. Schlossberger, it covered papers from 1933 through 1935.

This volume did not include any new *Strukturbericht* categories, but, perhaps to avoid confusion between the letters in all Latin alphabets, category I was renamed as category J. Since some of the previously defined I compounds were never mentioned in this or later volumes, we will attempt to resolve this confusion in Sections 4 and 5.

3.4. The War Years: *Strukturbericht* Band IV-VII

Each of the remaining volumes covered a single year and did not introduce any new categories. Remarkably, most were published while World War II was in progress, despite the fact that many German scientists, following Ewald, had fled the country, or like Hermann, had been imprisoned by the Nazi regime [50]. In spite of the War, these later volumes were distributed throughout the world. In 1943 in the United States, Edwards Brothers (Ann Arbor) published all seven volumes "in the Public Interest by Authority of the Alien Property Custodian," an office created by Presidential order to manage enemy alien property [51].

- SB-IV (covering 1936) [52], edited by Gottfried, was published in 1938.
- SB-V (1937) [53], again edited by Gottfried, was published in 1940.
- SB-VI (1938) [54], edited by K. Herrmann (not to be confused with C. Hermann), was published in 1941.
- SB-VII (1939) [55] — the last volume — was again edited by Herrmann and published in 1943, while the Eastern Front with the Soviet Union was collapsing, the Allies were invading Italy, and preparations for D-Day were underway. SB-VII begins by noting that scandium is face-centered cubic and so is an $A1$ structure, and ends with an entry for cronstedtite, $Fe^{II}_4Fe^{III}_2[(OH)_8|Fe^{III}_2Si_2O_{10}]$, assigning it the designation $S5_7$. This is the final inorganic *Strukturbericht* designation given in the original series.

3.5. Post-War: Continuation and Extension of *Strukturbericht*

The classification of compounds into *Strukturbericht* designations discontinued in the IUCr's post-war *Structure Reports*, but two authors created new designations as part of their own structural compilations. In 1949, Colin J. Smithells published his first *Metals Reference Book* [7], which included a partial index of *Strukturbericht* designations and the crystal structure associated with each. This list concentrated on metallic systems, and so left out some elements and compounds. This was not just a list of previously defined types, however. Smithells listed two new compounds: $ThSi_2$, and $TeCu_2$, provisionally giving them the *Strukturbericht* designation C_x. In the second edition of *Metals Reference* [8], Smithells extended this, adding many new *Strukturbericht* designations. While still following the letter-number-subscript scheme started in *Strukturbericht* II, the subscripts for the new entries were alphabetic rather than numeric. Thus, $ThSi_2$ became *Strukturbericht* type C_c and $TeCu_2$ became C_g. Reflecting the advances in physics in the war and post-war era, the list included the unique structures of the actinides: A_a (Pa), A_b (β-U), A_c (α-Np), A_d (β-Np), etc. Altogether Smithells added seventy-five new structures to the list.

In 1958, Pearson published the first edition of his influential *Handbook of Lattice Spacings and Structures of Metals and Alloys* [5]. It included most of Smithells' *Strukturbericht* index and added ten new structures, using the letter-number-alphabetic subscript format, with subscripts chosen to avoid overlap with Smithells. Pearson soon realized that further extensions would become unwieldy. In the next edition of his Handbook [6], he dropped the *Strukturbericht* index and listed structures by their Pearson symbols, which are not arbitrary and have the advantage of making the crystal type and number of atoms readily apparent, at the cost of having numerous structures of wildly different types using the same symbol. He still, however, gave the *Strukturbericht* designation for systems he was describing, a practice that continues in the modern editions of *Pearson's Handbook*.

3.6. Modern indices of compounds

Despite the demise of the *Strukturbericht* labeling system, it is still useful to have lists which classify crystal structures. Several authors and groups have done this over the years:

- Trotter and Bree [46] published an index to the original seven volumes of *Strukturbericht*, including both organic and inorganic structures. Regrettably, it did not include an index by *Strukturbericht* designation.
- Modern versions of *Pearson's Handbook* [10] and *Smithells Metals Reference Book* [11] include indices of metallic structures by *Strukturbericht* designation, Pearson symbol, and space group.
- The *Gmelin Handbook of Inorganic and Organometallic Chemistry* [12] also includes a list of *Strukturbericht* designations, as well as the Pearson symbol, space group, and occupied Wyckoff positions for each structure. Unlike the Pearson and Smithells Handbooks, this list includes many non-metallic systems. The list also notes when two *Strukturbericht* designations actually refer to the same structure. This happened several times in the original works, undoubtedly due to the frequent changes of editors and the lack of computers to keep accurate databases.
- The modern versions of Smithells, Pearson and Gmelin also include other structures not given a *Strukturbericht* designation.

Electronic databases are also available:

- *The American Mineralogist Crystal Structure Database* [13] gives structural information for a large number of minerals (including mineral names), but does not include the *Strukturbericht* designation. It does list the source publication, and when possible links to its PDF version. The site includes all structures published in *American Mineralogist*, *The Canadian Mineralogist*, *European Journal of Mineralogy*, and *Physics and Chemistry of Minerals*, as well as other sources.
- *Springer Materials* [14] maintains a constantly updated list of crystal structures. Edited by P. Villars, this is essentially an electronic form of *Pearson's Handbook*, but it includes non-metallic systems as well, unfortunately neglecting *Strukturbericht* designations.

None of these lists and databases give a comprehensive list of *Strukturbericht* symbols. In 1995, R. Benjamin Young, a summer intern at the U.S. Naval Research Laboratory in Washington D.C., and M. J. Mehl created an online database of *Strukturbericht* and related structures, in what was called the *Crystal Lattice Structures* web site, currently offline. In 2016, Mehl joined with the Curtarolo Research Group at Duke University to create the AFLOW Prototype Encyclopedia as an updated version of of the original website.

The original database contained 288 structures, including many, but not nearly all, of the inorganic *Strukturbericht* designations. It was later expanded [23] to 590 structures, and with this publication increases the count to well over 1,000, including all of the inorganic *Strukturbericht* designations from the original volumes, Smithells, Pearson, and other sources (see Section 5). We will continue to add new structures (Section 6), and invite readers to suggest prototypes (Section 7).

AFLOW Prototype Encyclopedia lists structures by *Strukturbericht* (if any), Pearson symbol, space group and prototype material. Each entry includes:

- The *Strukturbericht* designation of the structure, if one has been given,
- the Pearson symbol
- the space group,
- occupied Wyckoff positions,
- primitive vectors for the lattice,
- basis vectors for each atom in the primitive unit cell,
- a rotatable view of the primitive or conventional unit cell, which can be expanded to show multiple unit cells,
- a standardized Crystallographic Information File (CIF), and
- a structure file which can be used as input for a variety of electronic structure codes.

4. *Strukturbericht* categories and subcategories.

The original *Strukturbericht* scheme developed by Ewald was modified and extended over the years, and until now there has been no comprehensive list of *Strukturbericht* categories or designations. This section and the next correct this omission. Here we describe all the primary (A, B, C, ...) *Strukturbericht* categories, while the next section will list every *Strukturbericht* designation.

Primary categories from D through S are broken into smaller secondary categories, *e.g.*, $D0$, $E1$, $F2$, which we will describe as well. As the various primary and secondary categories changed over time, we offer a brief description of the changes in each category introduced over the course of the *Strukturbericht* publication run.

***A*:** Elements – this is essentially unchanged from Ewald's original *Handbuch* article.

***B*:** Binary compounds with stoichiometry AB – again unchanged from the original list.

***C*:** Binary compounds with stoichiometry AB_2 – also unchanged.

***D*:** Binary compounds with stoichiometry A_mB_n, beyond AB and AB_2.

In the *Handbuch*, this was just a list of compounds, in no particular order. SB-I changed this by introducing ranges corresponding to a given stoichiometry, so that compounds with the same m and n were in the same secondary category. Originally, these categories held at most ten compounds, *e.g.*, $D1 - D10$ for AB_3. In SB-II this was changed to the letter-numeral-subscript format described in the previous section:

***D0*:** (originally $D1 - D10$) Binary compounds with stoichiometry AB_3.

***D1*:** ($D11 - D20$) Binary compounds with stoichiometry AB_4. In SB-I, it was defined for all AB_n structure with $n > 3$, but all the examples given were AB_4.

***D2*:** This should have been the range $D21 - D30$, but this was not listed in SB-I. Starting with SB-II, it was used for binary compounds of stoichiometry AB_n, $n > 4$.

***D3*:** Binary compounds with stoichiometry $(AB)_2$.

***D4*:** Binary compounds with stoichiometry $(AB_n)_2$, $n > 1$.

Obviously, these last two categories can also be indexed in B, C or $D0 - D2$ by reducing the stoichiometries. The stoichiometry for $D3$ (A_2B_2) reduces to that of B (AB). Similarly, the stoichiometry for $D4$ can be reduced (A_2B_{2n} to AB_n), which is equivalent to C when $n = 2$, $D0$ when $n = 3$, $D1$ when $n = 4$, and $D2$ when $n > 4$. Therefore, only one compound was defined in $D3$ and $D4$.

The remaining secondary categories in D were listed together in SB-I as $D50 - D100$, and only implicitly divided into other categories in later volumes. This is our best interpretation of the method:

***D5*:** Binary compounds with stoichiometry A_2B_3.

***D6*:** Binary compounds with stoichiometry A_2B_4 – this is violated in SB-I by the first entry $D61$, but it seems to be latter editors' intent to define it in this way.

***D7*:** Binary compounds with stoichiometry A_3B_4.

***D8*:** Binary compounds with stoichiometry A_mB_n that do not fit into previous categories.

***D10*:** This was only defined post-war by Pearson, who used it for compounds with stoichiometry A_3B_7.

***E*:** Ternary and more complex compounds.

In the original *Handbuch* article, the E label was to be used for general multi-atom compounds that contained radicals. No E structures were introduced in SB-I, but subsequent volumes defined several secondary types, all involving ternary compounds:

***E0*:** Ternary compounds with stoichiometry ABC.

***E1*:** Ternary compounds with stoichiometry ABC_2.

***E2*:** Ternary compounds with stoichiometry ABC_n, $n > 2$.

***E3*:** Ternary compounds with stoichiometry AB_2C_n, $1 < n < 5$.

Pearson [5] gives $CdAl_2S_4$ the designation $E3$, without any subscript. It should be defined as

$E3_a$, but as he did not put any other structures into this subcategory, we will leave it as is.

E4: Ternary compounds with stoichiometry AB_2C_5.

E5: Ternary compounds with stoichiometry AB_2C_n, $5 < n < 8$.

E6: Ternary compounds with stoichiometry AB_2C_8.

E7: Never defined.

E8: Ternary compounds with stoichiometry $A_2B_2C_n$.

E9: Other ternary and quaternary compounds.

***F*:** Compounds with two- and three- atom radicals.

Originally, compounds of the form $A_m(BC)_n$, where "BC" form a radical, were to be given designations in the range $F1-F50$, with higher numbers for $A_m(BC_2)_n$ and $A_m(BCD)_n$ compounds. With the publication of SB-II, this became:

F0: Compounds with stoichiometry $A(BC)$.

F1: Compounds with stoichiometry $A(BC)_2$.

F2: Compounds with stoichiometry $A(BC)_m$, $m > 2$.

F3: Never defined.

F4: Compounds with stoichiometry $A_2(BC)_m$.

F5: Compounds with stoichiometry $A_m(BC_2)$ or $A_m(BCD)$.

F6: Additional compounds with stoichiometry $A_m(BC_2)$ or $A_m(BCD)$.

$F6$ was only used in SB-I for $F61$ (now $F6_1$), $CuFeS_2$ chalcopyrite.

Here, as well as in categories G-J, the letter "A" may represent more than one atom.

***G*:** Compounds with radicals of the form (BC_3):

G0: Compounds with stoichiometry $A(BC_3)$.

G1: Compounds with stoichiometry $AD(BC_3)$.

G2: Compounds with stoichiometry $A(BC_3)_2$.

G3: Compounds with higher order (BC_3) radicals.

G4: Never defined.

G5: Compounds with BO_3 and PO_3 radicals.

***H*:** Compounds with radicals of the form $(BC_n, n > 3)$.

This was originally a catch-all category, but in SB-II, compounds which would have been labeled as $H6$ and up were moved to a new class, I, and most silicates (SiO_n) were moved to S. This means that some structures have multiple *Strukturbericht* listings, so we will describe the complete H class here. The SB-II table of contents listed some compounds in both the H and I categories:

H0: Compounds with stoichiometry $A(BC_4)$.

H1: Compounds with stoichiometry $A_2(BC_4)$ or $AD(BC_4)$.

H2: More complex compounds with only one (BC_4) radical.

H3: Compounds with multiple (BC_4) radicals.

H4: Hydrated compounds with (BC_4) radicals.

H5: Compounds with (BC_5) radicals.

H6: Compounds with stoichiometry $A_2(BC_6)$.

This became $I1$ in SB-II and $J1$ in SB-III. $H6_2$ and $H6_4$ were only referred to as $I1_2$ and $I1_4$ in the SB-II table of contents, so we use the $H6$ designation in the list below. The examples given for $H6_3$ and $I1_3$ are rather different, so we use $H6_3$ for the original prototype $K_2Pt(SCN)_6$ and $I1_3$ for $SrCl_2{\cdot}(H_2O)_6$ as given in SB-II.

H7: Compounds with stoichiometry $A_3(BC_6)$.

This became $I2$ in SB-II and $J2$ in SB-III.

***I*:** Compounds with radicals of the form $A_m(BC_6)$.

This was introduced in SB-II, essentially transferring $H6$ to $I1$, $H7$ to $I2$, etc.:

I1: Compounds with stoichiometry $A_2(BC_6)$.

I2: Compounds with stoichiometry $A_3(BC_6)$.

I3: Never defined.

I4: Never defined.

I5: Never defined.

I6: Compounds with stoichiometry $A_6(BC_6)$.

The only structure listed here is $Ni(H_2O)_6SnCl_6$ ($I6_1$), which should have been subcategory $I5$ if the scheme laid out for $I1$ and $I2$ was used. It may have been the editors' intent to use this for hydrated systems. The $I6_1$ structure was never listed as $J6_1$, so we leave it in the I category in the AFLOW Prototype Encyclopedia.

***J*:** Compounds with radicals of the form $A_m(BC_6)$.

As noted above, this is the same category as I in SB-II, renamed in SB-III. Some structures in the I category were never renamed, and so we have left them with the I designation in Section 5:

J1: Compounds with stoichiometry $A_2(BC_6)$.

J2: Compounds with stoichiometry $A_3(BC_6)$.

J3: Compounds with stoichiometry $A_3(BC_6) \cdot mH_2O$.

The only structure with this designation is $K_3TlCl_6{\cdot}\ 2H_2O$, from which we infer the editors' intent.

***K*:** Compounds with more complicated radicals.

The intended range of these labels is open to question, as many of the entries have only one or two compounds:

- ***K*0:** Compounds with radicals of the form B_2C_n, $n < 6$.
- ***K*1:** Compounds with radicals of the form B_2C_n, $n = 6$.

  There is some confusion in this category. SB-II defines $K1_1$ $K_2S_2O_6$ and $K1_2$ $Ca_2S_2O_6$ in the text but uses $K2_1$ and $K2_2$ for the same structures in the table of contents. We use the labels in the actual text.
- ***K*2:** Never defined, except in the index of SB-II.
- ***K*3:** Compounds with radicals of the form B_mC_5.
- ***K*4:** Compounds with radicals of the form B_2C_n, $n > 5$.
- ***K*5:** Compounds with radicals of the form B_3C_n.
- ***K*6:** Compounds with radicals of the form B_mC_7.
- ***K*7:** Compounds with radicals of the form B_mC_n, $m < n$, $n > 7$.

***L*:** Alloys.

These are ordered intermetallic systems. To determine the subcategory, start with an elemental lattice, say fcc ($A1$), construct a supercell of that lattice, and populate that supercell with various atomic types, so that on average the structure looks like the original fcc lattice, even though it is an ordered crystal. This category could in principle be used with any of the lattices found in part A, *e.g.*, zincblende $B3$, a derivative of the diamond structure, and could have been listed as $L4_1$. In practice, though, only a few such structures were defined:

- ***L*1:** Supercells of the face-centered cubic (fcc) structure, $A1$.
- ***L*2:** Supercells of the body-centered cubic (bcc) structure, $A2$.
- ***L*3:** Supercells of the hexagonal-close-packed (hcp) structure, $A3$.
- ***L*6:** Supercells of the body-centered tetragonal (bct) structure, $A6$.

There are also labels such as $L'1_0$, Fe_4N. These indicate interstitial additions to the supercell. In this case, the atoms in Fe_3N are on fcc sites, and the extra iron atom is in the octahedral interstitial between the iron atoms.

***S*:** Silicates.

These compounds were broken off from the *H* category in SB-II, and some of the compounds have both *H* and *S* labels:

- ***S*0:** Silicates with isolated SiO_4 tetrahedra and other anions.
- ***S*1:** Silicates with isolated $(SiO_4)_n$ tetrahedra, $n > 1$.
- ***S*2:** Silicates with connected groups of SiO_4 tetrahedra.

  In these compounds, the oxygen ions are shared between two tetrahedra.
- ***S*3:** Silicates with one-dimensional Si-O chains.
- ***S*4:** Silicates with two-dimensional Si-O networks.
- ***S*5:** Silicates with three-dimensional Si-O networks.
- ***S*6:** Other silicates.

5. Index of *Strukturbericht* symbols in the Prototype Encyclopedia

Here we list all of the "inorganic" (not beginning with *O*) *Strukturbericht* designations we have found in the literature, with a link to the corresponding web page in the Prototype Encyclopedia. Our primary source is the seven volumes of *Strukturbericht*, but we include designations from other sources as well.

The list has five columns:

SB: This is the *Strukturbericht* designation appearing in the literature. When a structure has multiple designations, we choose one of these, usually the latest, to be the primary designation. If the current designation is the primary one, it will appear in bold and italics, otherwise it will be in italics (unbolded).

In several cases, the literature did not give sufficient information to allow us to reconstruct the structure. In those cases the designation is in italics and there is no primary designation nor any page in the Prototype Encyclopedia.

Prototype: This is the chemical formula of the prototype for this *Strukturbericht* designation. When available, the common name for this structure is also listed here. In most cases we use the prototype from the source in the next column. When we use another prototype, this column will have two entries: the first is the compound we use as the prototype, and the second is the compound defined as the prototype in the literature.

In some cases, later experiments have shown that a structure defined in the literature is incorrect. We will still have a Prototype Encyclopedia entry for these structures, but they will be marked "obsolete." Here

we indicate an obsolete structure by listing the prototype in italics. If data for the currently accepted structure is available, there will be a link to it from the obsolete structure's Prototype Encyclopedia page.

Source: This is where we found the definition of the current *Strukturbericht* designation. The entry, or entries, consist of a label defining the source followed by the page number, where available. We use the following abbreviations:

SB-I – SB-VII: The original *Strukturbericht* volumes [36, 47, 49, 52–55].

Smi: *Smithells Metals Reference,* 2^{nd} *Edition* (1955) [8]. This is the source of forty-five designations not found in the original *Strukturbericht* series.

Per: *Pearson's Handbook of Lattice Spacings and Structures of Metals and Alloys* (1958) [5] adds ten more designations, and repeats some of Smithells's list. Pearson's list is repeated in later editions of the Handbook, but often not in an easily searchable form.

Per91: *Pearson's Handbook* (1991 edition) [10].

Gmelin: *Gmelin Handbook* (1993) [12].

Hart: G. L. W. Hart and coworkers (2009, 2010, 2011, 2013) [56–59].

This reference here is the source which defines the *Strukturbericht designation*. The source we used to obtain the actual crystal structure is given in the prototype's Encyclopedia entry.

Other SB: We list other *Strukturbericht* designations for the current structure, as well as their sources. If the designation in the SB column is not the primary one, the primary will be highlighted in bold here. These duplications occurred because of shifting definition of the main *Strukturbericht* categories in the first three volumes, but there were also several editorial errors, including missing *Strukturbericht* designations ($C20$) and multiple definitions of the $S3_3$ and $S3_4$ structures.

AFLOW: The AFLOW label for this structure. These are only the primary designations. In the online version of this article, clicking on the structure label will open the corresponding web page in the AFLOW Prototype Encyclopedia.

SB	Prototype	Source	Other SB	AFLOW
***A*1**	*fcc* Cu	SB-I 13		A_cF4_225_a
***A*2**	*bcc* W	SB-I 15		A_cI2_229_a
***A*3**	*hcp* Mg	SB-I 16		A_hP2_194_c
***A*3′**	α-La	Per91		A_hP4_194_ac
***A*4**	diamond	SB-I 19		A_cF8_227_a
***A*5**	β-Sn	SB-I 21		A_tI4_141_a
***A*6**	In	SB-I 23		A_tI2_139_a.In
***A*7**	α-As	SB-I 25		A_hR2_166_c.alpha-As
***A*8**	γ-Se	SB-I 27		A_hP3_152_a
***A*9**	hexagonal graphite	SB-I 28		A_hP4_194_bc
***A*10**	α-Hg	SB-I 31		A_hR1_166_a.alpha-Hg
***A*11**	α-Ga	SB-I 738		A_oC8_64_f.alpha-Ga
***A*12**	α-Mn	SB-II 2		A_cI58_217_ac2g
***A*13**	β-Mn	SB-II 3		A_cP20_213_cd
***A*14**	I_2	SB-II 5		A_oC8_64_f.I
***A*15**	β-W / Cr_3Si	SB-II 6		A3B_cP8_223_c_a
***A*16**	α-S	SB-III 4		A_oF128_70_4h
***A*17**	P	SB-III 6		A_oC8_64_f.P
***A*18**	Cl_2	SB-IV 1		A_tP16_138_j
***A*19**	*Po*	SB-IV 4		A_mC12_5_3c
***A*20**	α-U	SB-VI 1		A_oC4_63_c
A_a	α-Pa	Smi 214		A_tI2_139_a.alpha-Pa
A_b	β-U	Smi 214		A_tP30_136_bf2ij
A_c	α-Np	Smi 214		A_oP8_62_2c
A_d	β-Np	Smi 214		A_tP4_129_ac
A_e	CuTi (alloy)	Smi 214	***A*20** SB-VI 1	
A_f	simple hexagonal	Smi 214		A_hP1_191_a
A_g	T-50 B	Smi 214		A_tP50_134_b2m2n
A_h	α-Po	Smi 215		A_cP1_221_a
A_i	β-Po	Smi 215		A_hR1_166_a.beta-Po
A_k	α-Se	Smi 215		A_mP64_14_16e
A_l	β-Se	Smi 215		A_mP32_14_8e
***B*1**	NaCl (rock salt)	SB-I 72		AB_cF8_225_a_b
***B*2**	CsCl	SB-I 74		AB_cP2_221_b_a
***B*3**	ZnS (zincblende)	SB-I 76		AB_cF8_216_c_a
***B*4**	ZnS (wurtzite)	SB-I 78		AB_hP4_186_b_b
***B*5**	SiC (moissanite-4H)	SB-I 80		AB_hP8_186_ab_ab
***B*6**	SiC (moissanite-6H)	SB-I 82		AB_hP12_186_a2b_a2b
***B*7**	SiC (moissanite-15R)	SB-I 83		AB_hR10_160_5a_5a
***B*8$_1$**	NiAs	Smi67 184	*B*8 SB-I 84	AB_hP4_194_c_a
***B*8$_2$**	Ni_2In	Smi67 184		AB2_hP6_194_c_ad
***B*9**	HgS (cinnabar)	SB-I 86		AB_hP6_154_a_b
***B*10**	PbO (litharge)	SB-I 89		AB_tP4_129_a_c
***B*11**	γ-CuTi	SB-I 94		AB_tP4_129_c_c
***B*12**	*BN*	SB-I 95		AB_hP4_186_b_a
***B*13**	NiS (millerite)	SB-I 740		AB_hR6_160_b_b
***B*14**	FeAs (westerveldite)	SB-II 7		AB_oP8_62_c_c.FeAs
*B*15	FeB	SB-II 7	***B*27** SB-III 12	
***B*16**	GeS	SB-II 8		AB_oP8_62_c_c.GeS
***B*17**	PtS (cooperite)	SB-II 9		AB_tP4_131_c_e
***B*18**	CuS (covellite)	SB-II 9		AB_hP12_194_df_ce
***B*19**	β'-AuCd	SB-II 11		AB_oP4_51_e_f
***B*20**	FeSi	SB-II 13		AB_cP8_198_a_a.FeSi

SB	Prototype	Source	Other SB	AFLOW
***B*21**	α-CO	SB-II 13		AB_cP8_198_a_a.alpha-CO
***B*22**	KSH	SB-III 8		AB_hR2_166_a_b.KSH
***B*23**	α-AgI	SB-III 8		A21B_cI44_229_bdh_a
***B*24**	TlF	SB-III 9		AB_oF8_69_a_b
***B*25**	NH_4Br	SB-III 10		AB4C_tP12_129_c_i_a
***B*26**	CuO (tenorite)	SB-III 11		AB_mC8_15_c_e
***B*27**	FeB	SB-III 12	*B*15 SB-II 7	AB_oP8_62_c_c.FeB
*B*28	FeSi	SB-III 13	***B*20** SB-II 13	
***B*29**	SnS	SB-III 14		AB_oP8_62_c_c.SnS
***B*30**	*MgZn*	SB-III 16		AB_oI48_44_6d_ab2cde
***B*31**	MnP	SB-III 17		AB_oP8_62_c_c.MnP
***B*32**	NaTl	SB-III 19		AB_cF16_227_a_b
***B*33**	TlI	SB-IV 6		AB_oC8_63_c_c
***B*34**	PdS	SB-V 1		AB_tP16_84_cej_k
***B*35**	CoSn	SB-VI 4		AB_hP6_191_f_ad
***B*36**	$LiOH{\cdot}H_2O$	SB-VII 4		A3BC2_mC24_12_ij_h_gi
***B*37**	TlSe	SB-VII 6		AB_tI16_140_ab_h
$\boldsymbol{B_a}$	CoU	Smi 217		AB_cI16_199_a_a
$\boldsymbol{B_b}$	ζ-AgZn	Smi 217		A2B_hP9_147_g_ad
B_c	CaSi	Smi 218	***B*33** SB-IV 6, B_f Smi 218	
$\boldsymbol{B_d}$	η-NiSi	Smi 218		AB_oP8_62_c_c.NiSi
$\boldsymbol{B_e}$	CdSb	Smi 218		AB_oP16_61_c_c
B_f	CrB	Smi 218	***B*33** SB-IV 6, B_c Smi 218	
$\boldsymbol{B_g}$	MoB	Smi 218		AB_tI16_141_e_e
$\boldsymbol{B_h}$	WC	Smi 218		AB_hP2_187_d_a
$\boldsymbol{B_i}$	TiAs/γ'-MoC	Smi 218 Per 95		AB_hP8_194_ad_f
$\boldsymbol{B_k}$	BN	Smi 218		AB_hP4_194_c_d
$\boldsymbol{B_l}$	AsS (realgar)	Smi 219		AB_mP32_14_4e_4e.realgar
B_m	TiB	Smi 219	***B*27** SB-III 12	
***C*1**	CaF_2 (fluorite)	SB-I 148		AB2_cF12_225_a_c
$\boldsymbol{C1_b}$	MgAgAs (half-Heusler)	Per 96		ABC_cF12_216_b_c_a
***C*2**	FeS_2 (pyrite)	SB-I 150		AB2_cP12_205_a_c
***C*3**	Cu_2O (cuprite)	SB-I 153		A2B_cP6_224_b_a
***C*4**	TiO_2 (rutile)	SB-I 155		A2B_tP6_136_f_a
***C*5**	TiO_2 (anatase)	SB-I 158		A2B_tI12_141_e_a
***C*6**	CdI_2 (hexagonal ω-phase)	SB-I 161		AB2_hP3_164_a_d
***C*7**	MoS_2 (molybdenite)	SB-I 164		AB2_hP6_194_c_f
***C*8**	SiO_2 (β-quartz)	SB-I 166		A2B_hP9_180_j_c
***C*9**	SiO_2 (β-cristobalite)	SB-I 169		A2B_cF24_227_c_a
***C*10**	SiO_2 (β-tridymite)	SB-I 171		A2B_hP12_194_cg_f
*C*11	CaC_2	SB-I 174	$\boldsymbol{C11_b}$ Smi 219	
$\boldsymbol{C11_a}$	CaC_2	Smi 219	*C*11 SB-I 174, 740	A2B_tI6_139_e_a
$\boldsymbol{C11_b}$	$MoSi_2$	Smi 219	*C*11 SB-I 740	AB2_tI6_139_a_e
***C*12**	$CaSi_2$	SB-I 175		AB2_hR6_166_c_2c
***C*13**	HgI_2	SB-I 177		AB2_tP6_137_a_d
***C*14**	$MgZn_2$ (hexagonal laves)	SB-I 180		AB2_hP12_194_f_ah
***C*15**	Cu_2Mg (cubic laves)	SB-I 490		A2B_cF24_227_d_a
$\boldsymbol{C15_b}$	$AuBe_5$	Per 99		AB5_cF24_216_a_ce
***C*16**	$CuAl_2$ (khatyrkite)	SB-I 491		A2B_tI12_140_h_a
***C*17**	Fe_2B	SB-I 493		AB2_tI12_121_ab_i
***C*18**	FeS_2 (marcasite)	SB-I 495		AB2_oP6_58_a_g.FeS2

SB	Prototype	Source	Other SB	AFLOW
***C*19**	α-Sm	SB-I 742		A_hR3_166_ac
***C*21**	TiO_2 (brookite)	SB-II 14		A2B_oP24_61_2c_c
***C*22**	Fe_2P	SB-II 15		A2B_hP9_150_ef_bd
***C*23**	$PbCl_2$ (cotunnite)	SB-II 16		A2B_oP12_62_2c_c.PbCl2
***C*24**	$HgBr_2$	SB-II 18		A2B_oC12_36_2a_a
***C*25**	$HgCl_2$	SB-II 19		A2B_oP12_62_2c_c.HgCl2
***C*26**	NO_2	SB-II 21	$C26_b$ SB-II 21	AB2_cI36_204_d_g
***C*26$_a$**	NO_2	SB-II 21		AB2_cI36_199_b_c
$C26_b$	NO_2	SB-II 21	***C*26** SB-II 21	
***C*27**	CdI_2	SB-III 22		AB2_hP6_186_b_ab
$C28$	$HgCl_2$	SB-III 23	***C*25** SB-II 19	
***C*29**	SrH_2	SB-III 24		A2B_oP12_62_2c_c.SrH2
***C*30**	SiO_2 (α-cristobalite)	SB-III 25		A2B_tP12_92_b_a
***C*31**	ϵ-$Zn(OH)_2$ (wülfingite)	SB-III 26		A2B2C_oP20_19_2a_2a_a
***C*32**	AlB_2 (hexagonal ω)	SB-III 28		AB2_hP3_191_a_d
***C*33**	Bi_2Te_3/Bi_2Te_2S	Sb-III 28		A2B3_hR5_166_c_ac
***C*34**	$AuTe_2$ (calaverite)	SB-III 30		AB2_mC6_12_a_i
***C*35**	$CaCl_2$ (hydrophilite)	SB-III 30		AB2_oP6_58_a_g.CaCl2
***C*36**	$MgNi_2$ (hexagonal laves)	SB-III 31		AB2_hP24_194_ef_fgh
***C*37**	Co_2Si	SB-III 32		A2B_oP12_62_2c_c.Co2Si
***C*38**	Cu_2Sb	SB-III. 33		A2B_tP6_129_ac_c
$C39$	ZrW_2	SB-III 34	***C*15** Sb-I 660	
***C*40**	$CrSi_2$	SB-III 35		AB2_hP9_180_d_j
$C41$	Fe_2W	SB-III 36	***C*14** SB-I 180	
***C*42**	SiS_2	SB-III 37		A2B_oI12_72_j_a
***C*43**	ZrO_2 (baddeleyite)	SB-IV 9		A2B_mP12_14_2e_e
***C*44**	GeS_2	SB-IV 11		AB2_oF72_43_ab_3b
***C*45**	$CuCl_2 \cdot 2H_2O$ (eriochalcite)	SB-IV 13		A2BC4D2_oP18_53_h_a_i_e
***C*46**	$AuTe_2$ (krennerite)	SB-IV 15		AB2_oP24_28_acd_2c3d
***C*47**	SeO_2 (downeyite)	SB-V 4		A2B_tP24_135_gh_h
$C48$	Cr_2Al	SB-V 5	***C*11$_b$** SB-I 147, 740; Smi 219	
***C*49**	$ZrSi_2$	SB-V 5		A2B_oC12_63_2c_c
***C*50**	α-$PdCl_2$	SB-VI 5		A2B_oP6_58_g_a
$C51$	$CoSn_2$	SB-VI 7	***C*16** SB-I 491	
***C*52**	β-TeO_2 (tellurite)	SB-VII 8		A2B_oP24_61_2c_c.TeO2
***C*53**	Br_2Sr	SB-VII 10		A2B_oP12_62_2c_c.SrBr2
***C*54**	$TiSi_2$	SB-VII 12		A2B_oF24_70_e_a
***C*$_a$**	Mg_2Ni	Smi 223		A2B_hP18_180_fi_bd
***C*$_b$**	Mg_2Cu	Smi 223		AB2_oF48_70_g_fg
***C*$_c$**	α-$ThSi_2$	Smi 223		A2B_tI12_141_e_a.ThSi2
***C*$_e$**	$PdSn_2$/$CoGe_2$	Smi 223		AB2_oC24_41_2a_2b
C_f	$RbSn_2$/$PdSn_2$	Smi 233	insufficient structural information	
***C*$_g$**	ThC_2	Smi 223		A2B_mC12_15_f_e
***C*$_h$**	Cu_2Te	Smi 223		A2B_hP6_191_h_e
***C*$_k$**	$LiZn_2$	Smi 223		AB_hP4_194_a_c
***D*0$_1$**	NH_3 (ammonia)	SB-I 230	$D1$	A3B_cP16_198_b_a
***D*0$_2$**	$CoAs_3$ (skutterudite)	SB-I 232	$D2$	A3B_cI32_204_g_c
***D*0$_3$**	BiF_3	SB-II 22		AB3_cF16_225_a_bc
***D*0$_4$**	$CrCl_3$	SB-II 23		A3B_hP24_151_3c_2a
***D*0$_5$**	BiI_3	SB-I 25		AB3_hR8_148_c_f
***D*0$_6$**	LaF_3	SB-II 27		A3B_hP24_193_ack_g

SB	Prototype	Source	Other SB	AFLOW
$\boldsymbol{D0_7}$	CrO_3	SB-II 29		AB3_oC16_20_a_bc
$\boldsymbol{D0_8}$	MoO_3	SB-II 30		AB3_oP16_62_c_3c
$\boldsymbol{D0_9}$	α-ReO_3	SB-II 31		A3B_cP4_221_d_a
$\boldsymbol{D0_{10}}$	WO_3	SB-II 32		A3B_oP16_57_a2d_d
$\boldsymbol{D0_{11}}$	Fe_3C (cementite)	SB-II 33		AB3_oP16_62_c_cd
$\boldsymbol{D0_{12}}$	FeF_3	SB-II 34		A3B_hR8_167_e_b
$\boldsymbol{D0_{13}}$	$AlCl_3$	SB-II 36		AB3_hP4_164_b_ad
$\boldsymbol{D0_{14}}$	AlF_3	SB-III 40		AB3_hR8_155_c_de
$\boldsymbol{D0_{15}}$	$AlCl_3$	SB-III 41		AB3_mC16_5_c_3c
$\boldsymbol{D0_{16}}$	NH_4I_3	SB-III 43		A3B_oP16_62_3c_c
$\boldsymbol{D0_{17}}$	BaS_3	SB-IV 18		AB3_tP8_113_a_ce
$\boldsymbol{D0_{18}}$	Na_3As	SB-V 6		AB3_hP8_194_c_bf
$\boldsymbol{D0_{19}}$	Ni_3Sn	SB-V 7		A3B_hP8_194_h_c
$\boldsymbol{D0_{20}}$	ϵ-$NiAl_3$	SB-V 8		A3B_oP16_62_cd_c
$\boldsymbol{D0_{21}}$	Cu_3P	SB-VI 7		A3B_hP24_165_bdg_f
$\boldsymbol{D0_{22}}$	$TiAl_3$	SB-VII 13		A3B_tI8_139_bd_a
$\boldsymbol{D0_{23}}$	$ZrAl_3$	SB-VII 14		A3B_tI16_139_cde_e
$\boldsymbol{D0_{24}}$	Ni_3Ti	SB-VII 14		A3B_hP16_194_gh_ac
$\boldsymbol{D0_a}$	β-$TiCu_3$	Smi 225		A3B_oP8_59_bf_a
$D0_b$	ζ-Ag_3Ga	Smi 225	$\boldsymbol{B_b}$ Smi. 217	
$\boldsymbol{D0_c}$	SiU_3	Smi 225	$D0'_c$ Gmelin 367	AB3_tI16_140_b_ah
$\boldsymbol{D0_d}$	Mn_3As	Smi 225		AB3_oC16_63_c_3c
$\boldsymbol{D0_e}$	Ni_3P	Per 105		A3B_tI32_82_3g_g
$\boldsymbol{D1_1}$	SnI_4	SB-I 234	$D11$	A4B_cP40_205_cd_c
$\boldsymbol{D1_2}$	SiF_4	SB-II 37		A4B_cI10_217_c_a
$\boldsymbol{D1_3}$	Al_4Ba	SB-VII 15		A4B_tI10_139_de_a
$\boldsymbol{D1_a}$	Ni_4Mo	Smi 225		AB4_tI10_87_a_h
$\boldsymbol{D1_b}$	Al_4U	Smi 226		A4B_oI20_74_beh_e
$\boldsymbol{D1_c}$	$PtSn_4$	Smi 226		AB4_oC20_41_a_2b
$\boldsymbol{D1_d}$	$PtPb_4$	Smi 226		A4B_tP10_125_m_a
$\boldsymbol{D1_e}$	ThB_4/UB_4	Smi 226		A4B_tP20_127_ehj_g
$\boldsymbol{D1_f}$	Mn_2B	Smi 226		AB2_oF48_70_f_fg
$\boldsymbol{D1_g}$	$B_{13}C_2$	Smi 226		A13B2_hR15_166_b2h_c
$\boldsymbol{D2_1}$	CaB_6	SB-II 37		A6B_cP7_221_f_a
$\boldsymbol{D2_2}$	$MgZn_5$	SB III 47		AB5_mC48_12_2i_ac5i2j
$\boldsymbol{D2_3}$	$NaZn_{13}$	SB VI 8		AB13_cF112_226_a_bi
$\boldsymbol{D2_a}$	$Be_{12}Ti$	Smi 227		A12B_hP13_191_cdei_a
$\boldsymbol{D2_b}$	$ThMn_{12}$	Smi 227		A12B_tI26_139_fij_a
$\boldsymbol{D2_c}$	MnU_6	Smi 227		AB6_tI28_140_a_hk
$\boldsymbol{D2_d}$	$CaCu_5$	Smi 227		AB5_hP6_191_a_cg
$\boldsymbol{D2_e}$	$BaHg_{11}$	Smi 227		AB11_cP36_221_c_agij
$\boldsymbol{D2_f}$	UB_{12}	Smi 227		A12B_cF52_225_i_a
$\boldsymbol{D2_g}$	Fe_8N	Smi 227		A8B_tI18_139_deh_a
$\boldsymbol{D2_h}$	$MnAl_6$	Per 109		A6B_oC28_63_efg_c
$\boldsymbol{D3_1}$	HgCl	SB I 237	$D31$	AB_tI8_139_e_e
$D4_1$	BH_3	SB I 239	$D41$, $\boldsymbol{A10}$ SB-I 31	
$\boldsymbol{D5_1}$	Al_2O_3 (corundum)	SB-I 240	$D51$	A2B3_hR10_167_c_e
$\boldsymbol{D5_2}$	La_2O_3	SB-I 242	$D52$	A2B3_hP5_164_d_ad
$\boldsymbol{D5_3}$	Mn_2O_3 (bixbyite)	SB-II 38		AB3C6_cI80_206_b_d_e
$D5_4$	Sb_2O_3 (senarmonite)	Gmelin 368	$\boldsymbol{D6_1}$ SB-I 245	
$\boldsymbol{D5_5}$	Mg_3P_2	SB-II 40		A3B2_cP10_224_d_b
$\boldsymbol{D5_6}$	Al_2O_3 (β-alumina)	SB-II 41		A2B3_hP60_194_3fk_cdef2k
$\boldsymbol{D5_7}$	γ-Al_2O_3/γ-Fe_2O_3 (maghemite)	SB-II 43		A2B3_cP60_212_bcd_ace

SB	Prototype	Source	Other SB	AFLOW
$D5_8$	Sb_2S_3 (stibnite)	SB-III 49		A3B2_oP20_62_3c_2c
$D5_9$	Zn_3P_2	SB-III 51		A2B3_tP40_137_cdf_3g
$D5_{10}$	Cr_3C_2 (tongbaite)	SB-III 53		A2B3_oP20_62_2c_3c
$D5_{11}$	Sb_2O_3 (valentinite)	SB-IV 20		A3B2_oP20_56_ce_e
$D5_{12}$	β-Bi_2O_3	SB-V 9		A2B3_tP20_117_i_adgh
$D5_{13}$	Ni_2Al_3	SB-V 10		A3B2_hP5_164_ad_d
$D5_a$	Si_2U_3	Smi 228		A2B3_tP10_127_g_ah
$D5_b$	Pt_2Sn_3	Smi 229		A2B3_hP10_194_f_bf
$D5_c$	Pu_2C_3	Smi 229		A3B2_cI40_220_d_c
$D5_e$	Ni_3S_2	Smi 229		A3B2_hR5_155_e_c
$D5_f$	As_2S_3 (orpiment)	Smi 229		A2B3_mP20_14_2e_3e
$D6_1$	Sm_2O_3 (senarmontite)	SB-I 245		A3B2_cF80_227_f_e
$D6_2$	Sb_2O_4	SB-III 54		A2B_cF96_227_abf_cd
$D7_1$	Al_4C_3	SB-III 56		A4B3_hR7_166_2c_ac
$D7_2$	Co_3O_4 (spinel)	SB-VI 9		A3B4_cF56_227_ad_e
$D7_3$	Th_3P_4	SB-VII 15		A4B3_cI28_220_c_a
$D7_a$	δ-Ni_3Sn_4	Smi 229		A3B4_mC14_12_ai_2i
$D7_b$	Ta_3B_4	Smi 230		A4B3_oI14_71_gh_cg
$D8_1$	Fe_3Zn_{10} (γ-brass)	SB-I 497		A3B10_cI52_229_e_fh
$D8_2$	Cu_5Zn_8 (γ-brass)	SB-I 497		A5B8_cI52_217_ce_cg
$D8_3$	Cu_9Al_4 (γ-brass)	SB-I 498		A4B9_cP52_215_ei_3efgi
$D8_4$	$Cr_{23}C_6$	SB-III 59		A6B23_cF116_225_e_acfh
$D8_5$	Fe_7W_6 (μ-phase)	SB-III 61		A7B6_hR13_166_ah_3c
$D8_6$	$Cu_{15}Si_4$	SB-III 62		A15B4_cI76_220_ae_c
$D8_7$	V_2O_5 *(shcherbinaite)*	SB-IV 22		A5B2_oP14_31_a2b_b
$D8_8$	Mn_5Si_3 (mavlyanovite)	SB-IV 24		A5B3_hP16_193_dg_g
$D8_9$	Co_9S_8	SB-IV 26		A9B8_cF68_225_af_ce
$D8_{10}$	Cr_5Al_8	SB-V 10		A8B5_hR26_160_a3bc_a3b
$D8_{11}$	Co_2Al_5	SB-VI 11		A5B2_hP28_194_ahk_ch
$D8_a$	Th_6Mn_{23}	Smi 231		A23B6_cF116_225_bd2f_e
$D8_b$	σ-CrFe (σ-phase)	Smi 231		sigma_tP30_136_bf2ij
$D8_c$	Mg_2Zn_{11} or $Mg_2Cu_6Al_5$	Smi 232		A2B11_cP39_200_f_aghij
$D8_d$	Co_2Al_9	Smi 232		A9B2_mP22_14_a4e_e
$D8_e$	$Mg_{32}(Al,Zn)_{49}$ (Bergman)	Smi 232		AB32C48_cI162_204_a_2efg_2gh
$D8_f$	Ir_3Ge_7	Smi 233		A7B3_cI40_229_df_e
$D8_g$	Mg_5Ga_2	Smi 233		A2B5_oI28_72_j_bfj
$D8_h$	W_2B_5	Smi 233		A5B2_hP14_194_abdf_f
$D8_i$	Mo_2B_5	Smi 233		A5B2_hR7_166_a2c_c
$D8_k$	Th_7S_{12}	Smi 233		A3B2_hP20_176_2h_ah
$D8_l$	Cr_5B_3	Per 117		A3B5_tI32_140_ah_cl
$D8_m$	W_5Si_3	Per 117		A3B5_tI32_140_ah_bk
$D10_1$	C_3Cr_7	Per 117		A3B7_oP40_62_cd_3c2d
$D10_2$	Fe_3Th_7	Per 117		A3B7_hP20_186_c_b2c
$E0_1$	PbFCl (matlockite)	SB-II 45		ABC_tP6_129_c_a_c
$E0_2$	AlO(OH) (diaspore)	SB-II 46		ABC2_oP16_62_c_c_2c
$E0_3$	Cd(OH)Cl	SB-III 65		ABCD_hP8_186_b_b_a_a
$E0_4$	γ-FeO(OH)	SB-III 66		AB2C2_oC20_63_c_f_2c
$E0_5$	FeOCl	SB-III 67		ABC_oP6_59_a_b_a
$E0_6$	MnO(OH) (manganite)	SB-IV 28		ABC2_mP16_14_e_e_2e
$E0_7$	FeAsS (arsenopyrite)	SB-IV 30		ABC_mP12_14_e_e_e
$E1_1$	$CuFeS_2$ (chalcopyrite)	SB-II 48		ABC2_tI16_122_a_b_d
$E1_2$	$Zn(NH_3)_2Cl_2$	SB-IV 31		A2B6C2D_oI44_74_h_ij_i_e
$E1_3$	$Mg(NH_3)_2Cl_2$ or $Cd(NH_3)_2Cl_2$	SB-IV 33		A2B8CD2_oC26_65_h_r_a_i

SB	Prototype	Source	Other SB	AFLOW
$E1_4$	$PNCl_2$	SB-VII 17		A2BC_tP32_86_2g_g_g
$E1_a$	$MgCuAl_2$	Smi 233		A2BC_oC16_63_f_c_c
$E1_b$	$AuAgTe_4$ (sylvanite)	Smi 234		ABC4_mP12_13_e_a_2g
$E2_1$	$CaTiO_3$ (perovskite)	SB-II 49	$G5/G0_5$ SB-I 300	AB3C_cP5_221_a_c_b
$E2_2$	$FeTiO_3$ (ilmenite)	Gmelin 369	$G4/G0_4$ SB-I 300	AB3C_hR10_148_c_f_c
$E2_3$	α-$LiIO_3$	SB-II 49		ABC3_hP10_182_c_b_g
$E2_4$	$(NH_4)CdCl_3$	SB-VI 13		AB3C_oP20_62_c_3c_c
$E2_5$	$(NH_4)HgCl_3$	SB-VI 15		A3BC_tP5_123_cg_a_d
$E2_6$	$KMg(H_2O)_6(Cl,Br)_3$ (bromocarnallite)	SB-VII 19		A3B6CD_tP44_85_bcg_3g_ac_e
$E3$	$CdAl_2S_4$	Per 118		A2BC4_tI14_82_bc_a_g
$E3_1$	β-Ag_2HgI_4	SB-II 50		A2BC4_tP7_111_f_a_n
$E3_2$	CaB_2O_4 I	SB-II 54		A2BC4_oP28_60_d_c_2d
$E3_3$	$FeSb_2S_4$ (berthierite)	SB-IV 35		AB4C2_oP28_62_c_4c_2c
$E3_4$	$K_2HgCl_4{\cdot}H_2O$	SB-VI 16		A4BCD2_oP32_55_ghi_f_e_gh
$E3_5$	$K_2SnCl_4{\cdot}H_2O$	SB-VII 22		A4BC2D_oP32_62_2cd_b_2c_a
$E4_1$	Fe_2TiO_5 (pseudobrookite)	SB-II 53		A2B5C_oC32_63_f_c2f_c
$E5_1$	$FeNb_2O_6$ (columbite or niobite)	SB-II 55		AB2C6_oP36_60_c_d_3d
$E6_1$	$Sr(OH)_2(H_2O)_8$	SB-II 56		A8B2C_tP11_123_r_f_a
$E6_2$	$SrO_2(H_2O)_8$	SB-II 57		A8B2C_tP11_123_r_h_a
$E8_1$	$Eu_2Ir_2O_7$ (cubic pyrochlore)	SB-II 58		A2B2C7_cF88_227_c_d_af
$E9_1$	$Ca_3Al_2O_6$	SB-II 60		A2B3C6_cP33_221_cd_ag_fh
$E9_2$	$NaBe_4SbO_7$ (swedenborgite)	SB-III 69		A4BC7D_hP26_186_ac_b_a2c_b
$E9_3$	Fe_3W_3C	SB-III 71		AB3C3_cF112_227_c_de_f
$E9_4$	Al_5C_3N	SB-III 73		A5B3C_hP18_186_2a3b_2ab_b
$E9_a$	$FeCu_2Al_7$	Smi 234		A7B2C_tP40_128_egi_h_e
$E9_b$	π-$FeMg_3Al_8Si_6$	Smi 235		A8BC3D6_hP18_189_bfh_a_g_i
$E9_c$	Al_9Mn_3Si	Smi 234		A9B3C_hP26_194_hk_h_a
$E9_d$	$AlLi_3N_2$	Smi 235		AB3C2_cI96_206_c_e_ad
$E9_e$	$CuFe_2S_3$ (cubanite)	Smi 235		AB2C3_oP24_62_c_d_cd
$F0_1$	NiSSb (ullmanite)	SB-I 269		ABC_cP12_198_a_a_a
$F0_2$	COS	SB-II 62		ABC_hR3_160_a_a_a
$F1_1$	$Hg(CN)_2$	SB-I 271	$F11$	A2BC2_tI40_122_e_d_e
$F2_1$	$K_4[Mo(CN)_8]{\cdot}2H_2O$	SB-VII 25		A8B4C4DE8F2_oP108_62_4c2d_2d_2cd_c_4c2d_d
$F4_1$	$Fe_2(CO)_9$	SB-VII 28		A9B2C9_hP40_176_hi_f_hi
$F5_1$	$NaCrS_2$ (caswellsilverite)/$NaHF_2$	SB-I 271	$F51$	ABC2_hR4_166_a_b_c
$F5_2$	KHF_2	SB-I 276	$F52$	A2BC_tI16_140_h_d_a
$F5_3$	$CuFeS_2$ (chalcopyrite)	SB-II XXII	$E1_1$ SB-II 48	
$F5_4$	NH_4ClO_2	SB-II 64		ABC2_tP8_100_b_a_c
$F5_5$	$NaNO_2$	SB-II 65		ABC2_oI8_44_a_a_c
$F5_6$	$CuSbS_2$ (chalcostibite)	SB-III 75		AB2C_oP16_62_c_2c_c
$F5_7$	$NH_4H_2PO_2$	SB-III 77		A2BC2D_oC24_67_m_a_n_g
$F5_8$	NH_4HF_2	SB-III 79		A2BC_oP16_53_eh_ab_g
$F5_9$	KCNS	SB-III 80		ABCD_oP16_57_d_c_d_d
$F5_{10}$	$KAg(CN)_2$	SB-III 82		AB2CD2_hP36_163_h_i_bf_i
$F5_{11}$	KNO_2	SB-IV 38		ABC2_mC8_8_a_a_b
$F5_{12}$	$AgNO_2$	SB-IV 41		ABC2_oI8_44_a_a_d
$F5_{13}$	KBO_2	SB-V 12		ABC2_hR24_167_e_e_2e
$F5_a$	$FeKS_2$	Smi 235		ABC2_mC16_15_e_e_f
$F6_1$	$CuFeS_2$ (chalcopyrite)	SB-I 279		ABC2_tP4_115_a_c_g
$G0_1$	$CaCO_3$ (calcite)	SB-I 292	$G1$	ABC3_hR10_167_a_b_e.CaCO3
$G0_2$	$CaCO_3$ (aragonite)	SB-I 295	$G2$	ABC3_oP20_62_c_c_cd
$G0_3$	$NaClO_3$	SB-I 297	$G3$	ABC3_cP20_198_a_a_b

SB	Prototype	Source	Other SB	AFLOW
$G0_4$	$FeTiO_3$	SB-I 300	*G*4 $\boldsymbol{E2_2}$ Gmelin 369	
$G0_5$	$CaTiO_3$ (perovskite)	SB-I 300	*G*5 $\boldsymbol{E2_1}$ SB-II 49	
$\boldsymbol{G0_6}$	$KClO_3$	SB-II 66		ABC3_mP10_11_e_e_ef
$\boldsymbol{G0_7}$	$KBrO_3$	SB-II 68		ABC3_hR5_160_a_a_b.KBrO3
$\boldsymbol{G0_8}$	NH_4NO_3 I	SB-II 69		AB_cP2_221_a_b.NH4.NO3
$\boldsymbol{G0_9}$	NH_4NO_3 II	SB-II 69		ABC3_tP10_100_b_a_bc
$\boldsymbol{G0_{10}}$	NH_4NO_3 III	SB-II 70		ABC3_oP20_62_c_c_cd.N.NH4.O
$\boldsymbol{G0_{11}}$	NH_4NO_3 IV	SB-II 71		A4B2C3_oP18_59_ef_ab_af
$\boldsymbol{G0_{12}}$	$NaHCO_3$	SB-III 86		ABCD3_mP24_14_e_e_e_3e
$\boldsymbol{G1_1}$	$MgCa(CO_3)_2$ (dolomite)	SB-I 303	*G*11	A2BCD6_hR10_148_c_a_b_f
$\boldsymbol{G2_1}$	$Pb(NO_3)_2$	SB-II 73	*G*21	A2B6C_cP36_205_c_d_a
$\boldsymbol{G2_2}$	$Nd(BrO_3)_3{\cdot}9H_2O$	SB-VII 29		A3B9CD9_hP44_186_c_3c_b_cd
$G3_1$	$Be_3Al_2(SiO_3)_6$ (beryl)	SB-I 305	*G*31 $\boldsymbol{S3_1}$ SB-II XXIV	
$\boldsymbol{G3_2}$	$Na_2(SO_3)$	SB-II 74		A2B3C_hP12_147_abd_g_d
$\boldsymbol{G5_1}$	H_3BO_3 (boric acid)	SB-III 87		AB3C3_aP28_2_2i_6i_6i
$\boldsymbol{G5_2}$	$Al(PO_3)_3$	SB-V 15		AB9C3_cI208_220_c_3e_e
$\boldsymbol{G7_1}$	$CeF(CO_3)$ (bastnäsite)	SB-II 75		ABCD3_hP36_190_h_g_af_hi
$\boldsymbol{G7_2}$	$Be_2BO_3(OH)$ (hambergite)	SB-II 78		AB2CD4_oP64_61_c_2c_c_4c
$\boldsymbol{G7_3}$	*$Na_3MgCl(CO_3)_2$ (northupite)*	SB-II 80		A2BCD3E6_cF208_227_e_c_d_f_g
$\boldsymbol{G7_4}$	$Cu_3(CO_3)_2(OH)_2$ (azurite)	SB-II 81		A2B3C2D8_mP30_14_e_ce_e_4e
$\boldsymbol{G7_5}$	*$PbCO_3{\cdot}PbCl_2$ (phosgenite)*	SB-III 89		AB2C3D2_tP16_90_c_f_ce_e
$\boldsymbol{H0_1}$	$CaSO_4$ (anhydrite)	SB-I 340	*H*1	AB4C_oC24_63_c_fg_c
$\boldsymbol{H0_2}$	$BaSO_4$ (barite)	SB-I 313	*H*2	AB4C_oP24_62_c_2cd_c
$H0_3$	$ZrSiO_4$ (zircon)	SB-I 345	*H*3 $\boldsymbol{S1_1}$ SB-II XXIV	
$\boldsymbol{H0_4}$	$CaWO_4$ (scheelite)	SB-I 347		AB4C_tI24_88_b_f_a
$\boldsymbol{H0_5}$	$KClO_4$ (high-temperature phase)	SB-II 84		ABC4_cF24_216_b_a_e
$\boldsymbol{H0_6}$	$MgWO_4$ (huanzalaite)	SB-II 85		AB4C_mP12_13_f_2g_e
$\boldsymbol{H0_7}$	BPO_4	SB-III 92		AB4C_tI12_82_c_g_a
$\boldsymbol{H0_8}$	$TlAlF_4$	SB-V 17		AB4C_tP6_123_d_eh_a
$\boldsymbol{H0_9}$	$AgMnO_4$	SB-VI 19		ABC4_mP24_14_e_e_4e
$\boldsymbol{H0_{10}}$	$KICl_4{\cdot}H_2O$	SB-VI 20		A4BCD_mP28_14_4e_e_e_e
$\boldsymbol{H1_1}$	Al_2MgO_4 (spinel)	SB-I 350	*H*11	A2BC4_cF56_227_d_a_e
$H1_2$	Mg_2SiO_4 (forsterite)	SB-I 352	*H*12 $\boldsymbol{S1_2}$	
$H1_3$	Be_2SiO_4 (phenakite)	SB-I 356	*H*13 $\boldsymbol{S1_3}$	
$\boldsymbol{H1_4}$	$KLiSO_4$	SB-I 357	*H*14	ABC4D_hP14_173_a_b_bc_b
$\boldsymbol{H1_5}$	K_2PtCl_4	SB-I 358	*H*15	A4B2C_tP7_123_j_e_a
$\boldsymbol{H1_6}$	K_2SO_4 (arcanite)	SB-II 86		A2B4C_oP28_62_2c_2cd_c
$\boldsymbol{H1_7}$	Na_2SO_4 (thenardite)	SB-II 88		A2B4C_oF56_70_g_h_a
$\boldsymbol{H1_8}$	Na_2CrO_4	SB-IV 45		AB2C4_oC28_63_c_bc_fg
$\boldsymbol{H2_1}$	Ag_3PO_4	SB-I 360	*H*21	A3B4C_cP16_218_c_e_a
$\boldsymbol{H2_2}$	KH_2PO_4	SB-I 362	*H*22	A4BC4D_tI40_122_e_b_e_a
$\boldsymbol{H2_4}$	Cu_3VS_4 (sulvanite)	SB-III 94		A3B4C_cP8_215_d_e_a
$\boldsymbol{H2_5}$	Cu_3AsS_4 (enargite)	SB-III 96		AB3C4_oP16_31_a_ab_2ab
$\boldsymbol{H2_6}$	Cu_2FeSnS_4 (stannite)	SB-III 96		A2BC4D_tI16_121_d_a_i_b
$\boldsymbol{H2_7}$	$Zn_2(OH)AsO_4$ (adamite)	SB-V 17		ABC5D2_oP36_58_g_g_3gh_eg
$\boldsymbol{H2_8}$	$BaAl_2O_4$	SB-V 19		A2BC6_hP18_182_f_b_gh
$H3_1$	$Al_2Ca_3(SiO_4)_3$ (garnet)	SB-I 363	*H*31 $\boldsymbol{S1_4}$ SB-III 150	
$\boldsymbol{H3_2}$	$AlK(SO_4)_2$ (steklite)	SB-II 90		ABC8D2_hP12_150_b_a_dg_d
$\boldsymbol{H4_1}$	$K_2CuCl_4{\cdot}2H_2O$	SB-I 366	*H*41	A4BC4D2E2_tP26_136_fg_a_j_d_e
$H4_2$	$KAl(SO_4)_2{\cdot}12H_2O$ (α-alum)	SB-I 369	*H*42 $\boldsymbol{H4_{13}}$ SB-III 108	
$\boldsymbol{H4_3}$	$BeSO_4{\cdot}4H_2O$	SB-II 91		AB8C8D_tI72_120_c_2i_2i_b
$\boldsymbol{H4_4}$	$Cu(NH_4)_2(SO_4)_2{\cdot}6H_2O$	SB-II 93		AB20C2D14E2_mP78_14_a_10e_e_7e_e
$\boldsymbol{H4_5}$	α-$NiSO_4{\cdot}6H_2O$ (retgersite)	SB-II 95		A12BC10D_tP96_92_6b_a_5b_a

SB	Prototype	Source	Other SB	AFLOW
$\boldsymbol{H4_6}$	$CaSO_4{\cdot}2H_2O$ (gypsum)	SB-II 97/IV 47		AB4C6D_mC48_15_e_2f_3f_e
$\boldsymbol{H4_7}$	$CaSO_4(H_2O)_{0.5}$ (bassanite)	SB-III 97		A2B2C9D2_mC90_5_ab2c_3c_b13c_3c
$\boldsymbol{H4_8}$	$Li_2SO_4{\cdot}H_2O$	SB-III 98		A2B2C5D_mP20_4_2a_2a_5a_a
$\boldsymbol{H4_9}$	$Pd(NH_3)_4Cl_2{\cdot}H_2O$	SB-III 100		A2BC4D_tP16_127_h_d_i_a
$\boldsymbol{H4_{10}}$	$CuSO_4{\cdot}5H_2O$ (chalcanthite)	SB-III 102		AB10C9D_aP42_2_ae_10i_9i_i
$\boldsymbol{H4_{11}}$	$Mg(ClO_4)_2{\cdot}6H_2O$	SB-III 103		A2B6CD8_oP34_31_2a_2a2b_a_4a2b
$\boldsymbol{H4_{12}}$	$NiSO_4{\cdot}7H_2O$ (morenosite)	SB-III 105		A14BC11D_oP108_19_14a_a_11a_a
$\boldsymbol{H4_{13}}$	$KAl(SO_4)_2{\cdot}12H_2O$ (α-alum)	SB-III 108		AB24CD28E2_cP224_205_a_4d_b_2c4d_c
$\boldsymbol{H4_{14}}$	$Al(NH_3CH_3)_2(SO_4)_2{\cdot}12H_2O$ (β-alum)	SB-III 111		AB2C36D2E20F2_cP252_205_a_c_6d_c_c3d_c
$\boldsymbol{H4_{15}}$	$AlNa(SO_4)_2{\cdot}12H_2O$ (γ-alum)	SB-III 112		AB24CD20E2_cP192_205_a_4d_b_c3d_c
$\boldsymbol{H4_{16}}$	$H_3PW_{12}O_{40}{\cdot}5H_2O$	SB-III 113		A5B40CD12_cP116_224_cd_e3k_a_k
$\boldsymbol{H4_{17}}$	$Ag_2SO_4{\cdot}4NH_3$	SB-III 115		A2B12C4D4E_tP46_114_d_3e_e_e_a
$\boldsymbol{H4_{18}}$	$LiClO_4{\cdot}3H_2O$	SB-III 117		AB6CD7_hP30_186_b_d_a_b2c
$\boldsymbol{H4_{19}}$	$KAuBr_4{\cdot}2H_2O$	SB-IV 50		AB4C2D_mP32_14_e_4e_2e_e
$\boldsymbol{H4_{20}}$	$(CdSO_4)_3{\cdot}8H_2O$	SB-IV 52		A3B16C20D3_mC168_15_ef_8f_10f_ef
$\boldsymbol{H4_{21}}$	$H_3PW_{12}O_{40}{\cdot}29H_2O$	SB-IV 56		A29B40CD12_cF656_227_ae2fg_e3g_b_g
$\boldsymbol{H4_{22}}$	$BaNi(CN)_4{\cdot}4H_2O$	SB-VI 23		AB4C4D4E_mC56_15_e_2f_2f_2f_a
$\boldsymbol{H4_{23}}$	$K_2Mn(SO_4)_2{\cdot}4H_2O$ (leonite)	SB-VII 32		A8B2CD15E2_mC112_12_2i3j_j_ad_g4i5j_2i
$H5_1$	Al_2SiO_5 (kyanite)	SB-II 109	$\boldsymbol{S0_1}$ SB-II XXIII	
$H5_2$	Al_2SiO_5 (sillimanite)	SB-II 112	$\boldsymbol{S0_3}$ SB-II XXIII	
$H5_3$	Al_2SiO_5 (andalusite)	SB-II 110	$\boldsymbol{S0_2}$ SB-II XXIII	
$H5_4$	$Fe(OH)_2Al_4SiO_{10}$	SB-II 113	$\boldsymbol{S0_4}$ SB-II XXIV	
$H5_5$	$Al_2SiO_4F_2$ (topaz)	SB-II 117	$\boldsymbol{S0_5}$ SB-II XXIV	
$\boldsymbol{H5_6}$	*$Na_6Mg_2SO_4(CO_3)_4$ (tychite)*	SB-II 98		A4B2C6D16E_cF232_227_e_d_f_eg_a
$\boldsymbol{H5_7}$	$Ca_5F(PO_4)_3$ (fluorapatite)	SB-II 99		A5BC12D3_hP42_176_fh_a_2hi_h
$\boldsymbol{H5_8}$	$Na_6ClF(SO_4)_2$ (sulphohalite)	SB-III 118		ABC6D8E2_cF72_225_b_a_e_f_c
$\boldsymbol{H5_9}$	*$Ca(UO_2)_2(PO_4)_2 \cdot 10\frac{1}{2}H_2O$ (autunite)*	SB-VI 24		AB2C2_tI10_139_a_d_e
$\boldsymbol{H5_{10}}$	$Ca(UO_2)_2(PO_4)_2 \cdot 6H_2O$ (meta-autunite I)	SB-VI 25		AB4C6DE_tP26_129_c_j_2ci_a_c
$H6_1$	K_2PtCl_6	SB-I 429	$H61$ $I1_1$ SB-II XXIII $\boldsymbol{J1_1}$ SB-III 121	
$\boldsymbol{H6_2}$	$K_2Sn(OH)_6$	SB-I 431	$I1_2$ SB-II XXIII	A6B2C6D_hR15_148_f_c_f_a
$\boldsymbol{H6_3}$	$K_2Pt(SCN)_6$	SB-I 433	$I1_3$ SB-II 102	A2BC6_hP9_164_d_a_i
$\boldsymbol{H6_4}$	*$Ni(NO_3)_2(NH_3)_6$*	SB-I 435	$H64$ $I1_4$ SB-II XXIII	A2B6CD6_cP60_205_c_d_a_d
$H7_1$	$(NH_4)_3AlF_6$	SB-I 437	$H71$ $I2_1$ SB-II XXIII $\boldsymbol{J2_1}$ SB-III 127	
$I1_1$	K_2PtCl_6	SB-II XXIII	$H61/H6_1$ SB-I 429 $\boldsymbol{J1_1}$ SB-III 121	
$I1_2$	$K_2Sn(OH)_6$	SB-II XXIII	$H62/H6_2$ SB-I 431	
$\boldsymbol{I1_3}$	*$SrCl_2{\cdot}(H_2O)_6$*	SB-II 102	$H63$ $H6_3$ SB-I 433	A2B6C_hP9_162_d_k_a
$I1_4$	*$Ni(NO_3)_2(NH_3)_6$*	SB-II XXIII	$H64$ $\boldsymbol{H6_4}$ SB-I 435	
$I2_1$	$(NH_4)_3AlF_6$	SB-II XXIII	$H71/H7_1$ SB-I 437 SB-II XXIII $\boldsymbol{J2_1}$ SB-III 127	
$\boldsymbol{I6_1}$	$Ni(H_2O)_6SnCl_6$	SB-II 102		A6B6CD_hR14_148_f_f_b_a
$\boldsymbol{J1_1}$	K_2PtCl_6	SB-III 121	$H61/H6_1$ SB-I 429 $I1_1$ SB-II XXIII	A6B2C_cF36_225_e_c_a
$\boldsymbol{J1_5}$	$K_2OsO_2Cl_4$	SB-III 122		A4B2C2D_tI18_139_h_d_e_a
$\boldsymbol{J1_6}$	$(NH_4)_2SiF_6$ (bararite)	SB-III 123		A6B2C_hP9_164_i_d_a
$\boldsymbol{J1_7}$	$MgCl_2 \cdot 6H_2O$ (bischofite)	SB-III 124		A2B12CD6_mC42_12_i_2i2j_a_ij
$\boldsymbol{J1_8}$	$RhCl_2(NH_3)_5Cl$	SB-III 125		A3B15C5D_oP96_62_cd_3c6d_3cd_c
$\boldsymbol{J1_9}$	$Ag[Co(NH_3)_2(NO_2)_4]$	SB-IV 60		ABC4D2E8_tP32_126_a_b_h_e_k
$\boldsymbol{J1_{10}}$	$Zn(BrO_3)_2 \cdot 6H_2O$	SB-IV 63		A2B6C6D_cP60_205_c_d_d_a

SB	Prototype	Source	Other SB	AFLOW
$\boldsymbol{J1_{11}}$	$NaSb(OH)_6$	SB-VI 26		AB6C_tP32_86_d_3g_c
$\boldsymbol{J1_{12}}$	$NaSbF_4(OH)_2$	SB-VI 27		A6BC_hP16_163_i_b_c
$\boldsymbol{J1_{13}}$	K_2GeF_6	SB-VII 34		A6BC2_hP9_164_i_a_d
$\boldsymbol{J2_1}$	$(NH_4)_3AlF_6$	SB-III 34	$H7_1/H7_1$ SB-I 437 $I2_1$ SB-II XXIII	AB30C16D3_cF200_225_a_ej_2f_bc
$\boldsymbol{J2_2}$	$CrCl_3(H_2O)_6$	SB-III 127		A3BC6_hR20_167_e_b_f
$\boldsymbol{J2_3}$	$Ca_3Al_2(OH)_{12}$	SB-III 129		A2B3C12D12_cI232_230_a_c_h_h
$\boldsymbol{J2_4}$	$K_3Co(NO_2)_6$	SB-IV 65		AB3C6D12_cF88_202_a_bc_e_h
$\boldsymbol{J2_5}$	$Cu_3[Fe(CN)_6]_2 \cdot xH_2O$	SB-VI 28		A6B9CD2E6_cF96_225_e_bf_a_c_e
$\boldsymbol{J2_6}$	Na_3AlF_6 (cryolite)	SB-VI 29		AB6C3_mP20_14_a_3e_de
$\boldsymbol{J3_1}$	$K_3TlCl_6{\cdot}2H_2O$	SB-III 131		A6B2C3D_tI168_139_egikl2m_ejn_bh2n_acf
$\boldsymbol{K0_1}$	$K_2S_2O_5$	SB-II 104	$K1_1$ SB-II XXIII	A2B5C2_mP18_11_2e_e2f_2e
$\boldsymbol{K1_1}$	KSO_3	SB-II 106	$K2_1$ SB-II XXIII	AB3C_hP30_150_ef_3g_c2d
$\boldsymbol{K1_2}$	$CaSO_3$	SB-II 107	$K2_2$ SB-II XXIII	AB3C_hP20_190_ac_i_f
$\boldsymbol{K3_1}$	Cs_3CoCl_5	SB-III 134		A5BC3_tI36_140_cl_b_ah
$\boldsymbol{K3_3}$	Tl_2AlF_5	SB-V 20		AB5C2_oC32_20_b_a2bc_c
$\boldsymbol{K3_4}$	$NH_4Pb_2Br_5$	SB-V 22		A5BC2_tI32_140_bl_a_h
$\boldsymbol{K3_5}$	$KB_5O_8 \cdot 4H_2O$ (santite)	SB-V 23		A5B8CD12_oC104_41_a2b_4b_a_6b
$\boldsymbol{K4_1}$	$(NH_4)SO_4$	SB-III 136		AB4C_mP24_14_e_4e_e.K41
$\boldsymbol{K5_1}$	$K_2S_3O_6$	SB-III 138		A2B6C3_oP44_62_2c_2c2d_3c
$\boldsymbol{K6_1}$	ZrP_2O_7	SB-III 140		A7B2C_cP40_205_bd_c_a
$\boldsymbol{K6_2}$	K_2NbF_7	SB-VII 36		A7B2C_mP40_14_7e_2e_e
$\boldsymbol{K7_1}$	$K_3W_2Cl_9$	SB-III 142		A9B3C2_hP28_176_hi_af_f
$\boldsymbol{K7_2}$	$Cs_3Tl_2Cl_9$	SB-III 145		A9B3C2_hR28_167_ef_e_c
$\boldsymbol{K7_3}$	$Cs_3As_2Cl_9$	SB-III 145		A2B9C3_hP14_150_d_eg_ad
$\boldsymbol{K7_4}$	$12CaO \cdot 7Al_2O_3$ (mayenite)	SB-IV 68		A7B12C19_cI152_220_bc_2d_ace
$\boldsymbol{K7_5}$	$Na_5Al_3F_{14}$ (chiolite)	SB-VI 30		A3B14C5_tP44_128_ac_ehi_bg
$\boldsymbol{K7_6}$	$AuCsCl_3$	SB-VI 33		AB3C_tI20_139_ab_eh_d
$\boldsymbol{L1_0}$	CuAu	SB-I 484	$L10$	AB_tP2_123_a_d
$\boldsymbol{L1_1}$	CuPt (rhombohedral)	SB-I 485	$L11$	AB_hR2_166_a_b
$\boldsymbol{L1_2}$	Cu_3Au (bogdanovite)	SB-I 496	$L12$	AB3_cP4_221_a_c
$\boldsymbol{L1_3(I)}$	CuPt (cubic)	SB-I 486	$L13$	AB_cF32_227_c_d
$\boldsymbol{L1_3(II)}$	$CdPt_3$	Hart		AB3_oC8_65_a_bf
$\boldsymbol{L1_a}$	$CuPt_3$	Smi 236		AB7_cF32_225_b_ad
$\boldsymbol{L'1_0}$	γ-Fe_4N	SB-I 487		A4B_cP5_221_bc_a
$L'1_2$	$AlFe_3C$	Smi 237	$\boldsymbol{E2_1}$ SB-II 49	
$L2_0$	PdCu	Sb-I 487	$\boldsymbol{B2}$ SB-I 72	
$\boldsymbol{L2_1}$	$AlCu_2Mn$ (heusler)	SB-I 488		AB2C_cF16_225_a_c_b
$\boldsymbol{L2_2}$	Sb_2Tl_7	SB-I 489		A2B7_cI54_229_e_afh
$\boldsymbol{L2_a}$	δ-TiCu	Smi 237		AB_tP2_123_a_d.CuTi
$\boldsymbol{L'2_0}$	C_xFe (“martensite type”)	SB-I 489		AB_tI4_139_b_a
$\boldsymbol{L'3_0}$	Fe_2N	Smi 237 Per 121		AB_hP4_194_c_a.Fe2N
$\boldsymbol{L'3_2}$	β-V_2N	Gmelin 371		AB2_hP9_162_ad_k
$\boldsymbol{L6_0}$	$CuTi_3$	Smi 237		AB3_tP4_123_a_ce
$L'6_0$	M_4N_x	Smi 237	insufficient structural information	
$\boldsymbol{S0_1}$	Al_2SiO_5 (kyanite)	SB-II 109	$H51$ $H5_1$ SB-I 427	A2B5C_aP32_2_4i_10i_2i
$\boldsymbol{S0_2}$	Al_2SiO_5 (andalusite)	SB-II 110	$H5_3$ SB-II XXIII	A2B5C_oP32_58_eg_3gh_g
$\boldsymbol{S0_3}$	Al_2SiO_5 (sillimanite)	SB-II 112	$H5_2$ SB-II XXIII	A2B5C_oP32_62_bc_3cd_c
$\boldsymbol{S0_4}$	$Fe(OH)_2Al_4Si_2O_{10}$ *(staurolite)*	SB-II 113	$H5_4$ SB-II XXIV	A4BC12D2_oC76_63_eg_c_f3gh_g
$\boldsymbol{S0_5}$	$Al_2SiO_4F_2$ (topaz)	SB-II 116	$H5_5$ SB-II XXIV	A2B2C4D_oP36_62_d_d_2cd_c
$\boldsymbol{S0_6}$	$CaTiSiO_5$ (titanite)	SB-II 117		AB5CD_mC32_15_e_e2f_e_b

SB	Prototype	Source	Other SB	AFLOW
$S0_7$	$Mg(F,OH)_2{\cdot}Mg_2SiO_4$ (norbergite)	SB-II 117		A2B3C4D_oP40_62_d_cd_2cd_c
$S0_8$	$Al_{13}(OH,F)_{18}Si_5O_{20}Cl$ (zunyite)	SB-III 147		A13BC18D20E5_cF228_216_dh_b_fh_2eh_ce
$S1_1$	$ZrSiO_4$ (zircon)	SB-II XXIV	$H3\ H0_3$ SB-I 345	A4BC_tI24_141_h_b_a
$S1_2$	Mg_2SiO_4 (forsterite/olivine)	SB-II XXIV	$H12\ H1_2$ SB-I 352	A2B4C_oP28_62_ac_2cd_c
$S1_3$	Be_2SiO_4 (phenakite)	SB-II XXIV	$H13\ H1_3$ SB-I 356	A2B4C_hR42_148_2f_4f_f
$S1_4$	$Co_3Al_2(SiO_4)_3$ (garnet)	SB-II XXIV	$H31\ H3_1$ SB-I 363	A2B3C12D3_cI160_230_a_c_h_d
$S1_5$	$Bi_4(SiO_4)_3$ (eulytine)	SB-II 122		A4B12C3_cI76_220_c_e_a
$S2_1$	$(Sc,Y)_2Si_2O_7$ (thortveitite)	SB-II 124		A7B2C2_mC22_12_aij_h_i
$S2_2$	$Zn_4Si_2O_7(OH)_2{\cdot}H_2O$ (hemimorphite)	SB-II 125		A2B5CD2_oI40_44_2c_abcde_d_e
$S2_3$	$Ca_{10}Al_4(Mg,Fe)_2Si_9O_{34}(OH)_4$	SB-II 126		A4B10C2D34E4F9_tP252_126_k_ce2k_f_h8k_k_d2k
$S3_1$	$Be_3Al_2(SiO_3)_6$ (beryl)	SB-II XXIV	$G31\ G3_1$ SB-I 305	A2B3C18D6_hP58_192_c_f_lm_l
$S3_2$	$BaSi_4O_9$	SB-II 128		AB9C4_hP28_188_e_kl_ak
$S3_3(I)$	$Na_6Ca_2Al_6Si_6O_{24}(CO_3)_2$ (crancrinite)	SB-III 150		A3BCD3E15F3_hP52_173_c_b_b_c_5c_c
$S3_3(II)$	$CaSiO_3$ (parawollastonite)	SB-IV 71		AB3C_mP60_14_3e_9e_3e
$S3_4(I)$	$Ca_{1.4}Sr_{0.3}Al_{3.8}Si_{8.3}O_{24}{\cdot}13H_2O$ (chabazite)	SB-III 151		A5B21C24D12_hR62_166_a2c_ehi_fg2h_i
$S3_4(II)$	*$Na_2Zr(SiO_3)_3{\cdot}H_2O$ (catapleiite)*	SB-V 24		A3B2C9D3E_hP36_194_g_f_hk_h_a
$S4_1$	$CaMg(SiO_3)_2$ (diopside)	SB-II 130		ABC6D2_mC40_15_e_e_3f_f.S4_1
$S4_2$	$Ca_2Mg_5Si_8O_{22}(OH)_2$ (tremolite)	SB-II 131		A2B2C5D24E8_mC82_12_h_i_agh_2i5j_2j
$S4_3$	$MgSiO_3$ (enstatite)	SB-II 134		AB3C_oP80_61_2c_6c_2c
$S4_4$	$H_2Mg_5Fe_2Si_8O_{24}$ (anthophyllite)	SB-II 137		A2B5C22D2E8_oP156_62_d_c2d_2c10d_2c_4d
$S4_5$	$H_4Mg_3Si_2O_9$ (chrysotile)	SB-II 139		AB6C11D6E4_mC112_12_e_gi2j_i5j_2i2j_2j
$S4_6$	$Be_4Si_2O_7(OH)_2$ (bertrandite)	SB-II 141		A4B7C2D2_oC60_36_2b_a3b_2a_b
$S4_7$	$BeHNaO_8Si_3$ (epididymite)	SB-III 153		ABCD8E3_oP112_62_d_2c_d_4c6d_3d
$S5_1$	$KH_2Al_3Si_3O_{12}$ (muscovite)	SB-II 143		A2BC10D2E4_mC76_15_f_e_5f_f_2f
$S5_2$	$KCa_4Si_8O_{20}F{\cdot}8H_2O$ (apophyllite)	SB-II 145		A4BC16DE28F8_tP116_128_h_a_2i_b_g3i_i
$S5_3$	$Ca_2MgSi_2O_7$ (akermanite)	SB-II 146		A2BC7D2_tP24_113_e_a_cef_e
$S5_4$	$Al_2Si_2O_5(OH)_4$ (nacrite)	SB-III 155/VII 37		A2B4C9D2_mC68_9_2a_4a_9a_2a
$S5_5$	$Mg_3(Mg_2Al)(Si_3Al)O_{10}(OH)_8$ (chlorite)	SB-III 156		A3B5C4D2_mC112_15_a3ef_5f_4f_2f
$S5_6$	$AlSi_2O_5(OH)$ (pyrophyllite)	SB-III 159		AB5CD2_mC72_15_f_5f_f_2f
$S5_7$	$Fe(Fe,Si)[(OH)_2,O]O_3$ (cronstedtite)	SB-VII 40		AB3C2D_hR7_160_a_b_2a_a
$S6_1$	$NaAlSi_2O_6{\cdot}H_2O$ (analcime)	SB-II 148		A2B2C3D12E4_tI184_142_f_f_be_3g_g
$S6_2$	$Na_4(AlSiO_4)_3Cl$ (sodalite)	SB-II 150		A3BC4D12E3_cP46_218_d_a_e_i_c
$S6_3$	$CaB_2Si_2O_8$ (danburite)	SB-II 153		A2BC8D2_oP52_62_d_c_2c3d_d
$S6_4$	$Na_4Cl(AlSi_3)_3O_{24}$ (scapolite)	SB-II 155		AB4C24D12_tI82_87_a_h_2h2i_hi
$S6_5$	$NaAlSiO_4$ (α-carnegieite)	SB-II 158		ABC4D_cP28_198_a_a_ab_a
$S6_6$	Na_2CaSiO_4	SB-II 159		AB2C4D_cP32_198_a_2a_ab_a
$S6_7$	$KAlSi_3O_8$ (sanidine)	SB-III 161		AB8C4_mC52_12_i_gi3j_2j
$S6_8$	$NaAlSi_3O_8$ (albite)	SB-III 164		ABC8D3_aP26_2_i_i_8i_3i
$S6_9$	$(Na_{0.5}Ca_{0.3}K_{0.2})_8(Al_6Si_6O_{24})(SO_4)_{1.5}$ (hauyne)	SB-III 166		A3B4C4D4E16F4G3_cP76_218_c_e_e_e_ei_e_d
$S6_{10}$	$Na_2Al_2Si_3O_{10}{\cdot}2H_2O$ (natrolite)	SB-III 168		A2B4C2D12E3_oF184_43_b_2b_b_6b_ab

6. Future work

New structure-types are constantly being investigated for inclusion into the Prototype Encyclopedia. AFLOW's symmetry (AFLOW-SYM) [60] and crystal prototype finder (AFLOW-XtalFinder) [37] software modules are actively being employed to identify and classify prototypes from literature (*e.g.*, ICSD [61]). Unique prototypes — after comparing against those in Parts 1, 2, and 3 — will be added to future installments of the Prototype Encyclopedia. The symmetries of the prototypes (*i.e.*, Pearson symbol, space group, and Wyckoff positions) will be calculated autonomously and self-consistently via AFLOW-SYM [60]. AFLOW-XtalFinder will map structures into the degrees of freedom consistent with the International Tables for Crystallography, determine the AFLOW prototype label and parameters, and create geometry files for structures in this representation with a symbolic prototype generator [37].

7. User contributions

Online functionality has been developed to analyze the prototype of a user's input structure (http://aflow.org/prototype-encyclopedia/xtal-finder.html). The structure is automatically compared to all existing prototype entries via AFLOW-XtalFinder [37]. If a matching entry is found, the relevant prototype information is returned (*e.g.*, prototype label and parameters, *Strukturbericht* designation, symmetry descriptions, and weblink to the entry). If no matching entry is found — signaling a new prototype — users can report the structure by providing the following information in an online form: **i.** geometry file (in any standardized format), **ii.** reference(s) that reported the structure, **iii.** supplemental reference(s) used to find the structure (*e.g.*, catalogs or online databases), **iv.** any notable comments, and **v.** uploader's information (to credit the contributor on that structure's entry page, if desired). Unique prototype suggestions will be collected for future enrichment of the encyclopedia.

8. Outreach and education

AFLOW Schools (http://aflow.org/aflow-school/) provide learning modules to teach the theory of and provide hands-on examples for analyzing the structure of materials; including crystallographic symmetry, prototype classification, and structural similarity. Modules are also available for studying thermodynamics (polar and non-polar materials), accessing data in AFLOW.org (APIs), calculating thermomechanical properties, modeling disordered compounds, and predicting the properties of materials with machine learning.

9. Conclusion

This article presents the newest additions to the *AFLOW Library of Crystallographic Prototypes*, which now contains 1,100 entries. This includes all of the inorganic crystal structures that were given *Strukturbericht* designations in the original volumes or by references found in the literature. We also provide a brief history of the development of the *Strukturbericht* designations and the changes to the system over time.

Declaration of competing interest

The authors declare that they have no known competing financial interests or personal relationships that could have appeared to influence the work reported in this paper.

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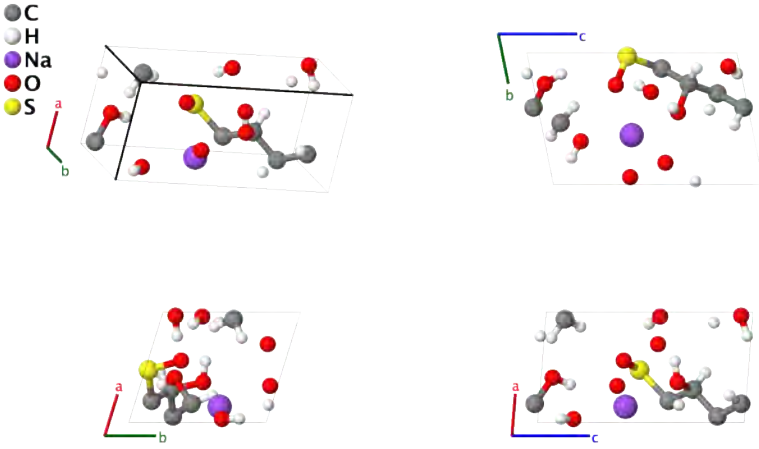
$NaC_5H_{11}O_8S$ Structure: A5B11CD8E_aP26_1_5a_11a_a_8a_a

http://aflow.org/prototype-encyclopedia/A5B11CD8E_aP26_1_5a_11a_a_8a_a

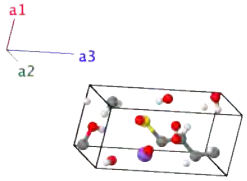


Prototype	:	$C_5H_{11}NaO_8S$
AFLOW prototype label	:	A5B11CD8E_aP26_1_5a_11a_a_8a_a
***Strukturbericht* designation**	:	None
Pearson symbol	:	aP26
Space group number	:	1
Space group symbol	:	$P1$
AFLOW prototype command	:	`aflow --proto=A5B11CD8E_aP26_1_5a_11a_a_8a_a --params=`$a, b/a, c/a, \alpha, \beta, \gamma, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}, x_{17}, y_{17}, z_{17}, x_{18}, y_{18}, z_{18}, x_{19}, y_{19}, z_{19}, x_{20}, y_{20}, z_{20}, x_{21}, y_{21}, z_{21}, x_{22}, y_{22}, z_{22}, x_{23}, y_{23}, z_{23}, x_{24}, y_{24}, z_{24}, x_{25}, y_{25}, z_{25}, x_{26}, y_{26}, z_{26}$

Triclinic primitive vectors:

$$\mathbf{a}_1 = a\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\cos\gamma\,\hat{\mathbf{x}} + b\sin\gamma\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c_x\hat{\mathbf{x}} + c_y\,\hat{\mathbf{y}} + c_z\,\hat{\mathbf{z}}$$

$$c_x = c\cos\beta$$
$$c_y = c\left(\cos\alpha - \cos\beta\cos\gamma\right)/\sin\gamma$$
$$c_z = \sqrt{c^2 - c_x^2 - c_y^2}$$



Basis vectors:

	Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$= x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	$=$	$(x_1a + y_1b\cos\gamma + z_1c_x)\,\hat{\mathbf{x}} + \left(y_1b\sin\gamma + z_1c_y\right)\hat{\mathbf{y}} + z_1c_z\,\hat{\mathbf{z}}$	$(1a)$	C I

$$\mathbf{B}_2 = x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3 = (x_2a + y_2b\cos\gamma + z_2c_x)\,\hat{\mathbf{x}} + \left(y_2b\sin\gamma + z_2c_y\right)\hat{\mathbf{y}} + z_2c_z\,\hat{\mathbf{z}} \quad (1a) \quad \text{C II}$$

$$\mathbf{B}_3 = x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3 = (x_3a + y_3b\cos\gamma + z_3c_x)\,\hat{\mathbf{x}} + \left(y_3b\sin\gamma + z_3c_y\right)\hat{\mathbf{y}} + z_3c_z\,\hat{\mathbf{z}} \quad (1a) \quad \text{C III}$$

$$\mathbf{B}_4 = x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3 = (x_4a + y_4b\cos\gamma + z_4c_x)\,\hat{\mathbf{x}} + \left(y_4b\sin\gamma + z_4c_y\right)\hat{\mathbf{y}} + z_4c_z\,\hat{\mathbf{z}} \quad (1a) \quad \text{C IV}$$

$$\mathbf{B}_5 = x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3 = (x_5a + y_5b\cos\gamma + z_5c_x)\,\hat{\mathbf{x}} + \left(y_5b\sin\gamma + z_5c_y\right)\hat{\mathbf{y}} + z_5c_z\,\hat{\mathbf{z}} \quad (1a) \quad \text{C V}$$

$$\mathbf{B}_6 = x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 = (x_6a + y_6b\cos\gamma + z_6c_x)\,\hat{\mathbf{x}} + \left(y_6b\sin\gamma + z_6c_y\right)\hat{\mathbf{y}} + z_6c_z\,\hat{\mathbf{z}} \quad (1a) \quad \text{H I}$$

$$\mathbf{B}_7 = x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 = (x_7a + y_7b\cos\gamma + z_7c_x)\,\hat{\mathbf{x}} + \left(y_7b\sin\gamma + z_7c_y\right)\hat{\mathbf{y}} + z_7c_z\,\hat{\mathbf{z}} \quad (1a) \quad \text{H II}$$

$$\mathbf{B}_8 = x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 = (x_8a + y_8b\cos\gamma + z_8c_x)\,\hat{\mathbf{x}} + \left(y_8b\sin\gamma + z_8c_y\right)\hat{\mathbf{y}} + z_8c_z\,\hat{\mathbf{z}} \quad (1a) \quad \text{H III}$$

$$\mathbf{B}_9 = x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 = (x_9a + y_9b\cos\gamma + z_9c_x)\,\hat{\mathbf{x}} + \left(y_9b\sin\gamma + z_9c_y\right)\hat{\mathbf{y}} + z_9c_z\,\hat{\mathbf{z}} \quad (1a) \quad \text{H IV}$$

$$\mathbf{B}_{10} = x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 = (x_{10}a + y_{10}b\cos\gamma + z_{10}c_x)\,\hat{\mathbf{x}} + \left(y_{10}b\sin\gamma + z_{10}c_y\right)\hat{\mathbf{y}} + z_{10}c_z\,\hat{\mathbf{z}} \quad (1a) \quad \text{H V}$$

$$\mathbf{B}_{11} = x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 = (x_{11}a + y_{11}b\cos\gamma + z_{11}c_x)\,\hat{\mathbf{x}} + \left(y_{11}b\sin\gamma + z_{11}c_y\right)\hat{\mathbf{y}} + z_{11}c_z\,\hat{\mathbf{z}} \quad (1a) \quad \text{H VI}$$

$$\mathbf{B}_{12} = x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 = (x_{12}a + y_{12}b\cos\gamma + z_{12}c_x)\,\hat{\mathbf{x}} + \left(y_{12}b\sin\gamma + z_{12}c_y\right)\hat{\mathbf{y}} + z_{12}c_z\,\hat{\mathbf{z}} \quad (1a) \quad \text{H VII}$$

$$\mathbf{B}_{13} = x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3 = (x_{13}a + y_{13}b\cos\gamma + z_{13}c_x)\,\hat{\mathbf{x}} + \left(y_{13}b\sin\gamma + z_{13}c_y\right)\hat{\mathbf{y}} + z_{13}c_z\,\hat{\mathbf{z}} \quad (1a) \quad \text{H VIII}$$

$$\mathbf{B}_{14} = x_{14}\,\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3 = (x_{14}a + y_{14}b\cos\gamma + z_{14}c_x)\,\hat{\mathbf{x}} + \left(y_{14}b\sin\gamma + z_{14}c_y\right)\hat{\mathbf{y}} + z_{14}c_z\,\hat{\mathbf{z}} \quad (1a) \quad \text{H IX}$$

$$\mathbf{B}_{15} = x_{15}\,\mathbf{a}_1 + y_{15}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 = (x_{15}a + y_{15}b\cos\gamma + z_{15}c_x)\,\hat{\mathbf{x}} + \left(y_{15}b\sin\gamma + z_{15}c_y\right)\hat{\mathbf{y}} + z_{15}c_z\,\hat{\mathbf{z}} \quad (1a) \quad \text{H X}$$

$$\mathbf{B}_{16} = x_{16}\,\mathbf{a}_1 + y_{16}\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 = (x_{16}a + y_{16}b\cos\gamma + z_{16}c_x)\,\hat{\mathbf{x}} + \left(y_{16}b\sin\gamma + z_{16}c_y\right)\hat{\mathbf{y}} + z_{16}c_z\,\hat{\mathbf{z}} \quad (1a) \quad \text{H XI}$$

$$\mathbf{B}_{17} = x_{17}\,\mathbf{a}_1 + y_{17}\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 = (x_{17}a + y_{17}b\cos\gamma + z_{17}c_x)\,\hat{\mathbf{x}} + \left(y_{17}b\sin\gamma + z_{17}c_y\right)\hat{\mathbf{y}} + z_{17}c_z\,\hat{\mathbf{z}} \quad (1a) \quad \text{Na}$$

$$\mathbf{B}_{18} = x_{18}\,\mathbf{a}_1 + y_{18}\,\mathbf{a}_2 + z_{18}\,\mathbf{a}_3 = (x_{18}a + y_{18}b\cos\gamma + z_{18}c_x)\,\hat{\mathbf{x}} + \left(y_{18}b\sin\gamma + z_{18}c_y\right)\hat{\mathbf{y}} + z_{18}c_z\,\hat{\mathbf{z}} \quad (1a) \quad \text{O I}$$

$$\mathbf{B}_{19} = x_{19}\,\mathbf{a}_1 + y_{19}\,\mathbf{a}_2 + z_{19}\,\mathbf{a}_3 = (x_{19}a + y_{19}b\cos\gamma + z_{19}c_x)\,\hat{\mathbf{x}} + \left(y_{19}b\sin\gamma + z_{19}c_y\right)\hat{\mathbf{y}} + z_{19}c_z\,\hat{\mathbf{z}} \quad (1a) \quad \text{O II}$$

$$\mathbf{B}_{20} = x_{20}\,\mathbf{a}_1 + y_{20}\,\mathbf{a}_2 + z_{20}\,\mathbf{a}_3 = (x_{20}a + y_{20}b\cos\gamma + z_{20}c_x)\,\hat{\mathbf{x}} + \left(y_{20}b\sin\gamma + z_{20}c_y\right)\hat{\mathbf{y}} + z_{20}c_z\,\hat{\mathbf{z}} \quad (1a) \quad \text{O III}$$

$$\mathbf{B}_{21} = x_{21}\,\mathbf{a}_1 + y_{21}\,\mathbf{a}_2 + z_{21}\,\mathbf{a}_3 = (x_{21}a + y_{21}b\cos\gamma + z_{21}c_x)\,\hat{\mathbf{x}} + \left(y_{21}b\sin\gamma + z_{21}c_y\right)\hat{\mathbf{y}} + z_{21}c_z\,\hat{\mathbf{z}} \quad (1a) \quad \text{O IV}$$

$$\mathbf{B}_{22} = x_{22}\,\mathbf{a}_1 + y_{22}\,\mathbf{a}_2 + z_{22}\,\mathbf{a}_3 = (x_{22}a + y_{22}b\cos\gamma + z_{22}c_x)\,\hat{\mathbf{x}} + \left(y_{22}b\sin\gamma + z_{22}c_y\right)\hat{\mathbf{y}} + z_{22}c_z\,\hat{\mathbf{z}} \quad (1a) \quad \text{O V}$$

$$\mathbf{B}_{23} = x_{23}\,\mathbf{a}_1 + y_{23}\,\mathbf{a}_2 + z_{23}\,\mathbf{a}_3 = (x_{23}a + y_{23}b\cos\gamma + z_{23}c_x)\,\hat{\mathbf{x}} + \left(y_{23}b\sin\gamma + z_{23}c_y\right)\hat{\mathbf{y}} + z_{23}c_z\,\hat{\mathbf{z}} \quad (1a) \quad \text{O VI}$$

$$\mathbf{B}_{24} = x_{24}\,\mathbf{a}_1 + y_{24}\,\mathbf{a}_2 + z_{24}\,\mathbf{a}_3 = (x_{24}a + y_{24}b\cos\gamma + z_{24}c_x)\,\hat{\mathbf{x}} + \left(y_{24}b\sin\gamma + z_{24}c_y\right)\hat{\mathbf{y}} + z_{24}c_z\,\hat{\mathbf{z}}$$ (1*a*) O VII

$$\mathbf{B}_{25} = x_{25}\,\mathbf{a}_1 + y_{25}\,\mathbf{a}_2 + z_{25}\,\mathbf{a}_3 = (x_{25}a + y_{25}b\cos\gamma + z_{25}c_x)\,\hat{\mathbf{x}} + \left(y_{25}b\sin\gamma + z_{25}c_y\right)\hat{\mathbf{y}} + z_{25}c_z\,\hat{\mathbf{z}}$$ (1*a*) O VIII

$$\mathbf{B}_{26} = x_{26}\,\mathbf{a}_1 + y_{26}\,\mathbf{a}_2 + z_{26}\,\mathbf{a}_3 = (x_{26}a + y_{26}b\cos\gamma + z_{26}c_x)\,\hat{\mathbf{x}} + \left(y_{26}b\sin\gamma + z_{26}c_y\right)\hat{\mathbf{y}} + z_{26}c_z\,\hat{\mathbf{z}}$$ (1*a*) S

References:

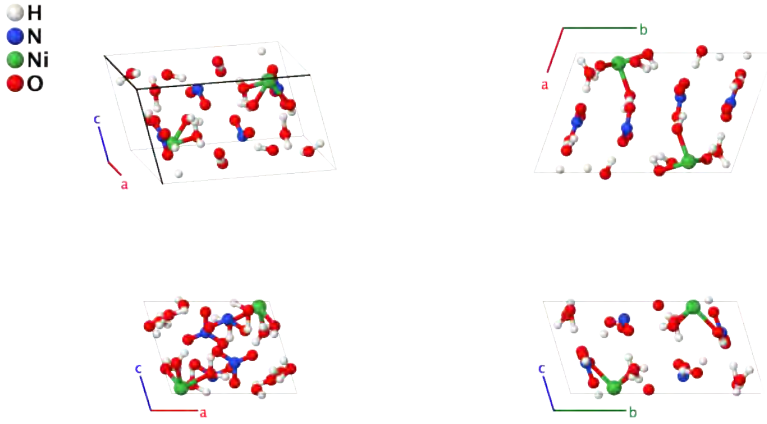
- A. H. Haines and D. L. Hughes, *Crystal structure of sodium (1S)-D-lyxit-1-yl sulfonate*, Acta Crystallogr. E **72**, 628–631 (2016), doi:10.1107/S2056989016005375.

Geometry files:

- CIF: pp. 1499
- POSCAR: pp. 1499

$Ni(NO_3)_2(H_2O)_6$ Structure: A12B2CD12_aP54_2_12i_2i_i_12i

http://aflow.org/prototype-encyclopedia/A12B2CD12_aP54_2_12i_2i_i_12i



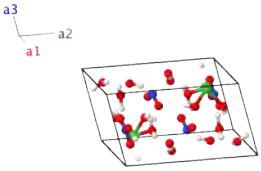
Prototype	:	$H_{12}N_2NiO_{12}$
AFLOW prototype label	:	A12B2CD12_aP54_2_12i_2i_i_12i
***Strukturbericht* designation**	:	None
Pearson symbol	:	aP54
Space group number	:	2
Space group symbol	:	$P\bar{1}$
AFLOW prototype command	:	`aflow --proto=A12B2CD12_aP54_2_12i_2i_i_12i --params=`$a, b/a, c/a, \alpha, \beta, \gamma, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}, x_{17}, y_{17}, z_{17}, x_{18}, y_{18}, z_{18}, x_{19}, y_{19}, z_{19}, x_{20}, y_{20}, z_{20}, x_{21}, y_{21}, z_{21}, x_{22}, y_{22}, z_{22}, x_{23}, y_{23}, z_{23}, x_{24}, y_{24}, z_{24}, x_{25}, y_{25}, z_{25}, x_{26}, y_{26}, z_{26}, x_{27}, y_{27}, z_{27}$

- This is the accepted structure for $Ni(NO_3)_2(H_2O)_6$, replacing the $H6_4$ structure. The positions of the hydrogen atoms are approximate.

Triclinic primitive vectors:

$$\mathbf{a}_1 = a\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\cos\gamma\,\hat{\mathbf{x}} + b\sin\gamma\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c_x\hat{\mathbf{x}} + c_y\,\hat{\mathbf{y}} + c_z\,\hat{\mathbf{z}}$$

$$c_x = c\cos\beta$$
$$c_y = c\,(\cos\alpha - \cos\beta\cos\gamma)/\sin\gamma$$
$$c_z = \sqrt{c^2 - c_x^2 - c_y^2}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$$
\begin{array}{lcccccc}
\mathbf{B}_1 & = & x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3 & = & \begin{array}{l}(x_1a + y_1b\cos\gamma + z_1c_x)\;\hat{\mathbf{x}} + \\ \left(y_1b\sin\gamma + z_1c_y\right)\;\hat{\mathbf{y}} + z_1c_z\;\hat{\mathbf{z}}\end{array} & (2i) & \text{H I} \\
\mathbf{B}_2 & = & -x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3 & = & \begin{array}{l}(-x_1a - y_1b\cos\gamma - z_1c_x)\;\hat{\mathbf{x}} + \\ \left(-y_1b\sin\gamma - z_1c_y\right)\;\hat{\mathbf{y}} - z_1c_z\;\hat{\mathbf{z}}\end{array} & (2i) & \text{H I} \\
\mathbf{B}_3 & = & x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3 & = & \begin{array}{l}(x_2a + y_2b\cos\gamma + z_2c_x)\;\hat{\mathbf{x}} + \\ \left(y_2b\sin\gamma + z_2c_y\right)\;\hat{\mathbf{y}} + z_2c_z\;\hat{\mathbf{z}}\end{array} & (2i) & \text{H II} \\
\mathbf{B}_4 & = & -x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3 & = & \begin{array}{l}(-x_2a - y_2b\cos\gamma - z_2c_x)\;\hat{\mathbf{x}} + \\ \left(-y_2b\sin\gamma - z_2c_y\right)\;\hat{\mathbf{y}} - z_2c_z\;\hat{\mathbf{z}}\end{array} & (2i) & \text{H II} \\
\mathbf{B}_5 & = & x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3 & = & \begin{array}{l}(x_3a + y_3b\cos\gamma + z_3c_x)\;\hat{\mathbf{x}} + \\ \left(y_3b\sin\gamma + z_3c_y\right)\;\hat{\mathbf{y}} + z_3c_z\;\hat{\mathbf{z}}\end{array} & (2i) & \text{H III} \\
\mathbf{B}_6 & = & -x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3 & = & \begin{array}{l}(-x_3a - y_3b\cos\gamma - z_3c_x)\;\hat{\mathbf{x}} + \\ \left(-y_3b\sin\gamma - z_3c_y\right)\;\hat{\mathbf{y}} - z_3c_z\;\hat{\mathbf{z}}\end{array} & (2i) & \text{H III} \\
\mathbf{B}_7 & = & x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3 & = & \begin{array}{l}(x_4a + y_4b\cos\gamma + z_4c_x)\;\hat{\mathbf{x}} + \\ \left(y_4b\sin\gamma + z_4c_y\right)\;\hat{\mathbf{y}} + z_4c_z\;\hat{\mathbf{z}}\end{array} & (2i) & \text{H IV} \\
\mathbf{B}_8 & = & -x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3 & = & \begin{array}{l}(-x_4a - y_4b\cos\gamma - z_4c_x)\;\hat{\mathbf{x}} + \\ \left(-y_4b\sin\gamma - z_4c_y\right)\;\hat{\mathbf{y}} - z_4c_z\;\hat{\mathbf{z}}\end{array} & (2i) & \text{H IV} \\
\mathbf{B}_9 & = & x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3 & = & \begin{array}{l}(x_5a + y_5b\cos\gamma + z_5c_x)\;\hat{\mathbf{x}} + \\ \left(y_5b\sin\gamma + z_5c_y\right)\;\hat{\mathbf{y}} + z_5c_z\;\hat{\mathbf{z}}\end{array} & (2i) & \text{H V} \\
\mathbf{B}_{10} & = & -x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3 & = & \begin{array}{l}(-x_5a - y_5b\cos\gamma - z_5c_x)\;\hat{\mathbf{x}} + \\ \left(-y_5b\sin\gamma - z_5c_y\right)\;\hat{\mathbf{y}} - z_5c_z\;\hat{\mathbf{z}}\end{array} & (2i) & \text{H V} \\
\mathbf{B}_{11} & = & x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 & = & \begin{array}{l}(x_6a + y_6b\cos\gamma + z_6c_x)\;\hat{\mathbf{x}} + \\ \left(y_6b\sin\gamma + z_6c_y\right)\;\hat{\mathbf{y}} + z_6c_z\;\hat{\mathbf{z}}\end{array} & (2i) & \text{H VI} \\
\mathbf{B}_{12} & = & -x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3 & = & \begin{array}{l}(-x_6a - y_6b\cos\gamma - z_6c_x)\;\hat{\mathbf{x}} + \\ \left(-y_6b\sin\gamma - z_6c_y\right)\;\hat{\mathbf{y}} - z_6c_z\;\hat{\mathbf{z}}\end{array} & (2i) & \text{H VI} \\
\mathbf{B}_{13} & = & x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 & = & \begin{array}{l}(x_7a + y_7b\cos\gamma + z_7c_x)\;\hat{\mathbf{x}} + \\ \left(y_7b\sin\gamma + z_7c_y\right)\;\hat{\mathbf{y}} + z_7c_z\;\hat{\mathbf{z}}\end{array} & (2i) & \text{H VII} \\
\mathbf{B}_{14} & = & -x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3 & = & \begin{array}{l}(-x_7a - y_7b\cos\gamma - z_7c_x)\;\hat{\mathbf{x}} + \\ \left(-y_7b\sin\gamma - z_7c_y\right)\;\hat{\mathbf{y}} - z_7c_z\;\hat{\mathbf{z}}\end{array} & (2i) & \text{H VII} \\
\mathbf{B}_{15} & = & x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 & = & \begin{array}{l}(x_8a + y_8b\cos\gamma + z_8c_x)\;\hat{\mathbf{x}} + \\ \left(y_8b\sin\gamma + z_8c_y\right)\;\hat{\mathbf{y}} + z_8c_z\;\hat{\mathbf{z}}\end{array} & (2i) & \text{H VIII} \\
\mathbf{B}_{16} & = & -x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3 & = & \begin{array}{l}(-x_8a - y_8b\cos\gamma - z_8c_x)\;\hat{\mathbf{x}} + \\ \left(-y_8b\sin\gamma - z_8c_y\right)\;\hat{\mathbf{y}} - z_8c_z\;\hat{\mathbf{z}}\end{array} & (2i) & \text{H VIII} \\
\mathbf{B}_{17} & = & x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 & = & \begin{array}{l}(x_9a + y_9b\cos\gamma + z_9c_x)\;\hat{\mathbf{x}} + \\ \left(y_9b\sin\gamma + z_9c_y\right)\;\hat{\mathbf{y}} + z_9c_z\;\hat{\mathbf{z}}\end{array} & (2i) & \text{H IX} \\
\mathbf{B}_{18} & = & -x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3 & = & \begin{array}{l}(-x_9a - y_9b\cos\gamma - z_9c_x)\;\hat{\mathbf{x}} + \\ \left(-y_9b\sin\gamma - z_9c_y\right)\;\hat{\mathbf{y}} - z_9c_z\;\hat{\mathbf{z}}\end{array} & (2i) & \text{H IX} \\
\mathbf{B}_{19} & = & x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 & = & \begin{array}{l}(x_{10}a + y_{10}b\cos\gamma + z_{10}c_x)\;\hat{\mathbf{x}} + \\ \left(y_{10}b\sin\gamma + z_{10}c_y\right)\;\hat{\mathbf{y}} + z_{10}c_z\;\hat{\mathbf{z}}\end{array} & (2i) & \text{H X} \\
\mathbf{B}_{20} & = & -x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 & = & \begin{array}{l}(-x_{10}a - y_{10}b\cos\gamma - z_{10}c_x)\;\hat{\mathbf{x}} + \\ \left(-y_{10}b\sin\gamma - z_{10}c_y\right)\;\hat{\mathbf{y}} - z_{10}c_z\;\hat{\mathbf{z}}\end{array} & (2i) & \text{H X} \\
\mathbf{B}_{21} & = & x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 & = & \begin{array}{l}(x_{11}a + y_{11}b\cos\gamma + z_{11}c_x)\;\hat{\mathbf{x}} + \\ \left(y_{11}b\sin\gamma + z_{11}c_y\right)\;\hat{\mathbf{y}} + z_{11}c_z\;\hat{\mathbf{z}}\end{array} & (2i) & \text{H XI} \\
\mathbf{B}_{22} & = & -x_{11}\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3 & = & \begin{array}{l}(-x_{11}a - y_{11}b\cos\gamma - z_{11}c_x)\;\hat{\mathbf{x}} + \\ \left(-y_{11}b\sin\gamma - z_{11}c_y\right)\;\hat{\mathbf{y}} - z_{11}c_z\;\hat{\mathbf{z}}\end{array} & (2i) & \text{H XI}
\end{array}
$$

$$\mathbf{B}_{23} = x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 = \begin{array}{l}(x_{12}a + y_{12}b\cos\gamma + z_{12}c_x)\,\hat{\mathbf{x}} + \\ \left(y_{12}b\sin\gamma + z_{12}c_y\right)\hat{\mathbf{y}} + z_{12}c_z\,\hat{\mathbf{z}}\end{array} \quad (2i) \quad \text{H XII}$$

$$\mathbf{B}_{24} = -x_{12}\,\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3 = \begin{array}{l}(-x_{12}a - y_{12}b\cos\gamma - z_{12}c_x)\,\hat{\mathbf{x}} + \\ \left(-y_{12}b\sin\gamma - z_{12}c_y\right)\hat{\mathbf{y}} - z_{12}c_z\,\hat{\mathbf{z}}\end{array} \quad (2i) \quad \text{H XII}$$

$$\mathbf{B}_{25} = x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3 = \begin{array}{l}(x_{13}a + y_{13}b\cos\gamma + z_{13}c_x)\,\hat{\mathbf{x}} + \\ \left(y_{13}b\sin\gamma + z_{13}c_y\right)\hat{\mathbf{y}} + z_{13}c_z\,\hat{\mathbf{z}}\end{array} \quad (2i) \quad \text{N I}$$

$$\mathbf{B}_{26} = -x_{13}\,\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 - z_{13}\,\mathbf{a}_3 = \begin{array}{l}(-x_{13}a - y_{13}b\cos\gamma - z_{13}c_x)\,\hat{\mathbf{x}} + \\ \left(-y_{13}b\sin\gamma - z_{13}c_y\right)\hat{\mathbf{y}} - z_{13}c_z\,\hat{\mathbf{z}}\end{array} \quad (2i) \quad \text{N I}$$

$$\mathbf{B}_{27} = x_{14}\,\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3 = \begin{array}{l}(x_{14}a + y_{14}b\cos\gamma + z_{14}c_x)\,\hat{\mathbf{x}} + \\ \left(y_{14}b\sin\gamma + z_{14}c_y\right)\hat{\mathbf{y}} + z_{14}c_z\,\hat{\mathbf{z}}\end{array} \quad (2i) \quad \text{N II}$$

$$\mathbf{B}_{28} = -x_{14}\,\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 - z_{14}\,\mathbf{a}_3 = \begin{array}{l}(-x_{14}a - y_{14}b\cos\gamma - z_{14}c_x)\,\hat{\mathbf{x}} + \\ \left(-y_{14}b\sin\gamma - z_{14}c_y\right)\hat{\mathbf{y}} - z_{14}c_z\,\hat{\mathbf{z}}\end{array} \quad (2i) \quad \text{N II}$$

$$\mathbf{B}_{29} = x_{15}\,\mathbf{a}_1 + y_{15}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 = \begin{array}{l}(x_{15}a + y_{15}b\cos\gamma + z_{15}c_x)\,\hat{\mathbf{x}} + \\ \left(y_{15}b\sin\gamma + z_{15}c_y\right)\hat{\mathbf{y}} + z_{15}c_z\,\hat{\mathbf{z}}\end{array} \quad (2i) \quad \text{Ni}$$

$$\mathbf{B}_{30} = -x_{15}\,\mathbf{a}_1 - y_{15}\,\mathbf{a}_2 - z_{15}\,\mathbf{a}_3 = \begin{array}{l}(-x_{15}a - y_{15}b\cos\gamma - z_{15}c_x)\,\hat{\mathbf{x}} + \\ \left(-y_{15}b\sin\gamma - z_{15}c_y\right)\hat{\mathbf{y}} - z_{15}c_z\,\hat{\mathbf{z}}\end{array} \quad (2i) \quad \text{Ni}$$

$$\mathbf{B}_{31} = x_{16}\,\mathbf{a}_1 + y_{16}\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 = \begin{array}{l}(x_{16}a + y_{16}b\cos\gamma + z_{16}c_x)\,\hat{\mathbf{x}} + \\ \left(y_{16}b\sin\gamma + z_{16}c_y\right)\hat{\mathbf{y}} + z_{16}c_z\,\hat{\mathbf{z}}\end{array} \quad (2i) \quad \text{O I}$$

$$\mathbf{B}_{32} = -x_{16}\,\mathbf{a}_1 - y_{16}\,\mathbf{a}_2 - z_{16}\,\mathbf{a}_3 = \begin{array}{l}(-x_{16}a - y_{16}b\cos\gamma - z_{16}c_x)\,\hat{\mathbf{x}} + \\ \left(-y_{16}b\sin\gamma - z_{16}c_y\right)\hat{\mathbf{y}} - z_{16}c_z\,\hat{\mathbf{z}}\end{array} \quad (2i) \quad \text{O I}$$

$$\mathbf{B}_{33} = x_{17}\,\mathbf{a}_1 + y_{17}\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 = \begin{array}{l}(x_{17}a + y_{17}b\cos\gamma + z_{17}c_x)\,\hat{\mathbf{x}} + \\ \left(y_{17}b\sin\gamma + z_{17}c_y\right)\hat{\mathbf{y}} + z_{17}c_z\,\hat{\mathbf{z}}\end{array} \quad (2i) \quad \text{O II}$$

$$\mathbf{B}_{34} = -x_{17}\,\mathbf{a}_1 - y_{17}\,\mathbf{a}_2 - z_{17}\,\mathbf{a}_3 = \begin{array}{l}(-x_{17}a - y_{17}b\cos\gamma - z_{17}c_x)\,\hat{\mathbf{x}} + \\ \left(-y_{17}b\sin\gamma - z_{17}c_y\right)\hat{\mathbf{y}} - z_{17}c_z\,\hat{\mathbf{z}}\end{array} \quad (2i) \quad \text{O II}$$

$$\mathbf{B}_{35} = x_{18}\,\mathbf{a}_1 + y_{18}\,\mathbf{a}_2 + z_{18}\,\mathbf{a}_3 = \begin{array}{l}(x_{18}a + y_{18}b\cos\gamma + z_{18}c_x)\,\hat{\mathbf{x}} + \\ \left(y_{18}b\sin\gamma + z_{18}c_y\right)\hat{\mathbf{y}} + z_{18}c_z\,\hat{\mathbf{z}}\end{array} \quad (2i) \quad \text{O III}$$

$$\mathbf{B}_{36} = -x_{18}\,\mathbf{a}_1 - y_{18}\,\mathbf{a}_2 - z_{18}\,\mathbf{a}_3 = \begin{array}{l}(-x_{18}a - y_{18}b\cos\gamma - z_{18}c_x)\,\hat{\mathbf{x}} + \\ \left(-y_{18}b\sin\gamma - z_{18}c_y\right)\hat{\mathbf{y}} - z_{18}c_z\,\hat{\mathbf{z}}\end{array} \quad (2i) \quad \text{O III}$$

$$\mathbf{B}_{37} = x_{19}\,\mathbf{a}_1 + y_{19}\,\mathbf{a}_2 + z_{19}\,\mathbf{a}_3 = \begin{array}{l}(x_{19}a + y_{19}b\cos\gamma + z_{19}c_x)\,\hat{\mathbf{x}} + \\ \left(y_{19}b\sin\gamma + z_{19}c_y\right)\hat{\mathbf{y}} + z_{19}c_z\,\hat{\mathbf{z}}\end{array} \quad (2i) \quad \text{O IV}$$

$$\mathbf{B}_{38} = -x_{19}\,\mathbf{a}_1 - y_{19}\,\mathbf{a}_2 - z_{19}\,\mathbf{a}_3 = \begin{array}{l}(-x_{19}a - y_{19}b\cos\gamma - z_{19}c_x)\,\hat{\mathbf{x}} + \\ \left(-y_{19}b\sin\gamma - z_{19}c_y\right)\hat{\mathbf{y}} - z_{19}c_z\,\hat{\mathbf{z}}\end{array} \quad (2i) \quad \text{O IV}$$

$$\mathbf{B}_{39} = x_{20}\,\mathbf{a}_1 + y_{20}\,\mathbf{a}_2 + z_{20}\,\mathbf{a}_3 = \begin{array}{l}(x_{20}a + y_{20}b\cos\gamma + z_{20}c_x)\,\hat{\mathbf{x}} + \\ \left(y_{20}b\sin\gamma + z_{20}c_y\right)\hat{\mathbf{y}} + z_{20}c_z\,\hat{\mathbf{z}}\end{array} \quad (2i) \quad \text{O V}$$

$$\mathbf{B}_{40} = -x_{20}\,\mathbf{a}_1 - y_{20}\,\mathbf{a}_2 - z_{20}\,\mathbf{a}_3 = \begin{array}{l}(-x_{20}a - y_{20}b\cos\gamma - z_{20}c_x)\,\hat{\mathbf{x}} + \\ \left(-y_{20}b\sin\gamma - z_{20}c_y\right)\hat{\mathbf{y}} - z_{20}c_z\,\hat{\mathbf{z}}\end{array} \quad (2i) \quad \text{O V}$$

$$\mathbf{B}_{41} = x_{21}\,\mathbf{a}_1 + y_{21}\,\mathbf{a}_2 + z_{21}\,\mathbf{a}_3 = \begin{array}{l}(x_{21}a + y_{21}b\cos\gamma + z_{21}c_x)\,\hat{\mathbf{x}} + \\ \left(y_{21}b\sin\gamma + z_{21}c_y\right)\hat{\mathbf{y}} + z_{21}c_z\,\hat{\mathbf{z}}\end{array} \quad (2i) \quad \text{O VI}$$

$$\mathbf{B}_{42} = -x_{21}\,\mathbf{a}_1 - y_{21}\,\mathbf{a}_2 - z_{21}\,\mathbf{a}_3 = \begin{array}{l}(-x_{21}a - y_{21}b\cos\gamma - z_{21}c_x)\,\hat{\mathbf{x}} + \\ \left(-y_{21}b\sin\gamma - z_{21}c_y\right)\hat{\mathbf{y}} - z_{21}c_z\,\hat{\mathbf{z}}\end{array} \quad (2i) \quad \text{O VI}$$

$$\mathbf{B}_{43} = x_{22}\,\mathbf{a}_1 + y_{22}\,\mathbf{a}_2 + z_{22}\,\mathbf{a}_3 = \begin{array}{l}(x_{22}a + y_{22}b\cos\gamma + z_{22}c_x)\,\hat{\mathbf{x}} + \\ \left(y_{22}b\sin\gamma + z_{22}c_y\right)\hat{\mathbf{y}} + z_{22}c_z\,\hat{\mathbf{z}}\end{array} \quad (2i) \quad \text{O VII}$$

$$\mathbf{B}_{44} = -x_{22}\,\mathbf{a}_1 - y_{22}\,\mathbf{a}_2 - z_{22}\,\mathbf{a}_3 = \begin{array}{l}(-x_{22}a - y_{22}b\cos\gamma - z_{22}c_x)\,\hat{\mathbf{x}} + \\ \left(-y_{22}b\sin\gamma - z_{22}c_y\right)\hat{\mathbf{y}} - z_{22}c_z\,\hat{\mathbf{z}}\end{array} \quad (2i) \quad \text{O VII}$$

$\mathbf{B}_{45}$	=	$x_{23}\,\mathbf{a}_1 + y_{23}\,\mathbf{a}_2 + z_{23}\,\mathbf{a}_3$	=	$(x_{23}a + y_{23}b\cos\gamma + z_{23}c_x)\,\hat{\mathbf{x}} + \left(y_{23}b\sin\gamma + z_{23}c_y\right)\hat{\mathbf{y}} + z_{23}c_z\,\hat{\mathbf{z}}$	(2*i*)	O VIII
$\mathbf{B}_{46}$	=	$-x_{23}\,\mathbf{a}_1 - y_{23}\,\mathbf{a}_2 - z_{23}\,\mathbf{a}_3$	=	$(-x_{23}a - y_{23}b\cos\gamma - z_{23}c_x)\,\hat{\mathbf{x}} + \left(-y_{23}b\sin\gamma - z_{23}c_y\right)\hat{\mathbf{y}} - z_{23}c_z\,\hat{\mathbf{z}}$	(2*i*)	O VIII
$\mathbf{B}_{47}$	=	$x_{24}\,\mathbf{a}_1 + y_{24}\,\mathbf{a}_2 + z_{24}\,\mathbf{a}_3$	=	$(x_{24}a + y_{24}b\cos\gamma + z_{24}c_x)\,\hat{\mathbf{x}} + \left(y_{24}b\sin\gamma + z_{24}c_y\right)\hat{\mathbf{y}} + z_{24}c_z\,\hat{\mathbf{z}}$	(2*i*)	O IX
$\mathbf{B}_{48}$	=	$-x_{24}\,\mathbf{a}_1 - y_{24}\,\mathbf{a}_2 - z_{24}\,\mathbf{a}_3$	=	$(-x_{24}a - y_{24}b\cos\gamma - z_{24}c_x)\,\hat{\mathbf{x}} + \left(-y_{24}b\sin\gamma - z_{24}c_y\right)\hat{\mathbf{y}} - z_{24}c_z\,\hat{\mathbf{z}}$	(2*i*)	O IX
$\mathbf{B}_{49}$	=	$x_{25}\,\mathbf{a}_1 + y_{25}\,\mathbf{a}_2 + z_{25}\,\mathbf{a}_3$	=	$(x_{25}a + y_{25}b\cos\gamma + z_{25}c_x)\,\hat{\mathbf{x}} + \left(y_{25}b\sin\gamma + z_{25}c_y\right)\hat{\mathbf{y}} + z_{25}c_z\,\hat{\mathbf{z}}$	(2*i*)	O X
$\mathbf{B}_{50}$	=	$-x_{25}\,\mathbf{a}_1 - y_{25}\,\mathbf{a}_2 - z_{25}\,\mathbf{a}_3$	=	$(-x_{25}a - y_{25}b\cos\gamma - z_{25}c_x)\,\hat{\mathbf{x}} + \left(-y_{25}b\sin\gamma - z_{25}c_y\right)\hat{\mathbf{y}} - z_{25}c_z\,\hat{\mathbf{z}}$	(2*i*)	O X
$\mathbf{B}_{51}$	=	$x_{26}\,\mathbf{a}_1 + y_{26}\,\mathbf{a}_2 + z_{26}\,\mathbf{a}_3$	=	$(x_{26}a + y_{26}b\cos\gamma + z_{26}c_x)\,\hat{\mathbf{x}} + \left(y_{26}b\sin\gamma + z_{26}c_y\right)\hat{\mathbf{y}} + z_{26}c_z\,\hat{\mathbf{z}}$	(2*i*)	O XI
$\mathbf{B}_{52}$	=	$-x_{26}\,\mathbf{a}_1 - y_{26}\,\mathbf{a}_2 - z_{26}\,\mathbf{a}_3$	=	$(-x_{26}a - y_{26}b\cos\gamma - z_{26}c_x)\,\hat{\mathbf{x}} + \left(-y_{26}b\sin\gamma - z_{26}c_y\right)\hat{\mathbf{y}} - z_{26}c_z\,\hat{\mathbf{z}}$	(2*i*)	O XI
$\mathbf{B}_{53}$	=	$x_{27}\,\mathbf{a}_1 + y_{27}\,\mathbf{a}_2 + z_{27}\,\mathbf{a}_3$	=	$(x_{27}a + y_{27}b\cos\gamma + z_{27}c_x)\,\hat{\mathbf{x}} + \left(y_{27}b\sin\gamma + z_{27}c_y\right)\hat{\mathbf{y}} + z_{27}c_z\,\hat{\mathbf{z}}$	(2*i*)	O XII
$\mathbf{B}_{54}$	=	$-x_{27}\,\mathbf{a}_1 - y_{27}\,\mathbf{a}_2 - z_{27}\,\mathbf{a}_3$	=	$(-x_{27}a - y_{27}b\cos\gamma - z_{27}c_x)\,\hat{\mathbf{x}} + \left(-y_{27}b\sin\gamma - z_{27}c_y\right)\hat{\mathbf{y}} - z_{27}c_z\,\hat{\mathbf{z}}$	(2*i*)	O XII

References:

- F. Bigoli, A. Braibanti, A. Tiripicchio, and M. T. Camellini, *The crystal structures of nitrates of divalent hexaaquocations. III. Hexaaquonickel nitrate*, Acta Crystallogr. Sect. B Struct. Sci. **27**, 1427–1434 (1971), doi:10.1107/S0567740871004084.

Found in:

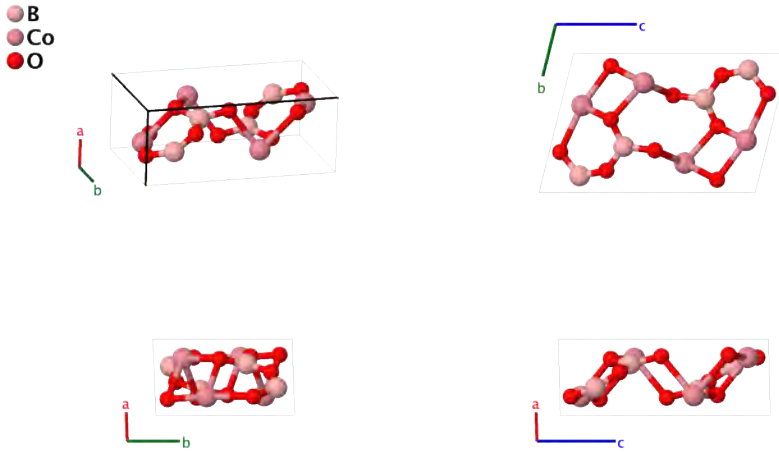
- J. Breternitz, L. J. Farrugia, A. Godula-Jopek, S. Saremi-Yarahmadi, I. E. Malka, T. K. A. Hoang, and D. H. Gregory, *Reaction of [Ni(H$_2$O)$_6$](NO$_3$)$_2$ with gaseous NH$_3$; crystal growth* via *in-situ solvation*, J. Cryst. Growth **412**, 1–6 (2015), doi:10.1016/j.jcrysgro.2014.11.021.

Geometry files:

- CIF: pp. 1499
- POSCAR: pp. 1500

$Co_2B_2O_5$ Structure: A2B2C5_aP18_2_2i_2i_5i

http://aflow.org/prototype-encyclopedia/A2B2C5_aP18_2_2i_2i_5i

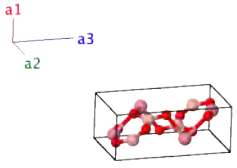


Prototype	:	$B_2Co_2O_5$
AFLOW prototype label	:	A2B2C5_aP18_2_2i_2i_5i
***Strukturbericht* designation**	:	None
Pearson symbol	:	aP18
Space group number	:	2
Space group symbol	:	$P\bar{1}$
AFLOW prototype command	:	`aflow --proto=A2B2C5_aP18_2_2i_2i_5i --params=`$a, b/a, c/a, \alpha, \beta, \gamma, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9$

Triclinic primitive vectors:

$$\mathbf{a}_1 = a\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\cos\gamma\,\hat{\mathbf{x}} + b\sin\gamma\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c_x\hat{\mathbf{x}} + c_y\,\hat{\mathbf{y}} + c_z\,\hat{\mathbf{z}}$$

$$c_x = c\,\cos\beta$$
$$c_y = c\,(\cos\alpha - \cos\beta\cos\gamma)/\sin\gamma$$
$$c_z = \sqrt{c^2 - c_x^2 - c_y^2}$$



Basis vectors:

	Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$= x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	$=$	$(x_1a + y_1b\cos\gamma + z_1c_x)\,\hat{\mathbf{x}} + \left(y_1b\sin\gamma + z_1c_y\right)\,\hat{\mathbf{y}} + z_1c_z\,\hat{\mathbf{z}}$	$(2i)$	B I
$\mathbf{B}_2$	$= -x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	$=$	$(-x_1a - y_1b\cos\gamma - z_1c_x)\,\hat{\mathbf{x}} + \left(-y_1b\sin\gamma - z_1c_y\right)\,\hat{\mathbf{y}} - z_1c_z\,\hat{\mathbf{z}}$	$(2i)$	B I
$\mathbf{B}_3$	$= x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	$=$	$(x_2a + y_2b\cos\gamma + z_2c_x)\,\hat{\mathbf{x}} + \left(y_2b\sin\gamma + z_2c_y\right)\,\hat{\mathbf{y}} + z_2c_z\,\hat{\mathbf{z}}$	$(2i)$	B II

$\mathbf{B}_4$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2a - y_2b\cos\gamma - z_2c_x)\,\hat{\mathbf{x}} + \left(-y_2b\sin\gamma - z_2c_y\right)\,\hat{\mathbf{y}} - z_2c_z\,\hat{\mathbf{z}}$	$(2i)$	B II
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3a + y_3b\cos\gamma + z_3c_x)\,\hat{\mathbf{x}} + \left(y_3b\sin\gamma + z_3c_y\right)\,\hat{\mathbf{y}} + z_3c_z\,\hat{\mathbf{z}}$	$(2i)$	Co I
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3a - y_3b\cos\gamma - z_3c_x)\,\hat{\mathbf{x}} + \left(-y_3b\sin\gamma - z_3c_y\right)\,\hat{\mathbf{y}} - z_3c_z\,\hat{\mathbf{z}}$	$(2i)$	Co I
$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4a + y_4b\cos\gamma + z_4c_x)\,\hat{\mathbf{x}} + \left(y_4b\sin\gamma + z_4c_y\right)\,\hat{\mathbf{y}} + z_4c_z\,\hat{\mathbf{z}}$	$(2i)$	Co II
$\mathbf{B}_8$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4a - y_4b\cos\gamma - z_4c_x)\,\hat{\mathbf{x}} + \left(-y_4b\sin\gamma - z_4c_y\right)\,\hat{\mathbf{y}} - z_4c_z\,\hat{\mathbf{z}}$	$(2i)$	Co II
$\mathbf{B}_9$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5a + y_5b\cos\gamma + z_5c_x)\,\hat{\mathbf{x}} + \left(y_5b\sin\gamma + z_5c_y\right)\,\hat{\mathbf{y}} + z_5c_z\,\hat{\mathbf{z}}$	$(2i)$	O I
$\mathbf{B}_{10}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5a - y_5b\cos\gamma - z_5c_x)\,\hat{\mathbf{x}} + \left(-y_5b\sin\gamma - z_5c_y\right)\,\hat{\mathbf{y}} - z_5c_z\,\hat{\mathbf{z}}$	$(2i)$	O I
$\mathbf{B}_{11}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6a + y_6b\cos\gamma + z_6c_x)\,\hat{\mathbf{x}} + \left(y_6b\sin\gamma + z_6c_y\right)\,\hat{\mathbf{y}} + z_6c_z\,\hat{\mathbf{z}}$	$(2i)$	O II
$\mathbf{B}_{12}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6a - y_6b\cos\gamma - z_6c_x)\,\hat{\mathbf{x}} + \left(-y_6b\sin\gamma - z_6c_y\right)\,\hat{\mathbf{y}} - z_6c_z\,\hat{\mathbf{z}}$	$(2i)$	O II
$\mathbf{B}_{13}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7a + y_7b\cos\gamma + z_7c_x)\,\hat{\mathbf{x}} + \left(y_7b\sin\gamma + z_7c_y\right)\,\hat{\mathbf{y}} + z_7c_z\,\hat{\mathbf{z}}$	$(2i)$	O III
$\mathbf{B}_{14}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7a - y_7b\cos\gamma - z_7c_x)\,\hat{\mathbf{x}} + \left(-y_7b\sin\gamma - z_7c_y\right)\,\hat{\mathbf{y}} - z_7c_z\,\hat{\mathbf{z}}$	$(2i)$	O III
$\mathbf{B}_{15}$	=	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8a + y_8b\cos\gamma + z_8c_x)\,\hat{\mathbf{x}} + \left(y_8b\sin\gamma + z_8c_y\right)\,\hat{\mathbf{y}} + z_8c_z\,\hat{\mathbf{z}}$	$(2i)$	O IV
$\mathbf{B}_{16}$	=	$-x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8a - y_8b\cos\gamma - z_8c_x)\,\hat{\mathbf{x}} + \left(-y_8b\sin\gamma - z_8c_y\right)\,\hat{\mathbf{y}} - z_8c_z\,\hat{\mathbf{z}}$	$(2i)$	O IV
$\mathbf{B}_{17}$	=	$x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$(x_9a + y_9b\cos\gamma + z_9c_x)\,\hat{\mathbf{x}} + \left(y_9b\sin\gamma + z_9c_y\right)\,\hat{\mathbf{y}} + z_9c_z\,\hat{\mathbf{z}}$	$(2i)$	O V
$\mathbf{B}_{18}$	=	$-x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$(-x_9a - y_9b\cos\gamma - z_9c_x)\,\hat{\mathbf{x}} + \left(-y_9b\sin\gamma - z_9c_y\right)\,\hat{\mathbf{y}} - z_9c_z\,\hat{\mathbf{z}}$	$(2i)$	O V

References:

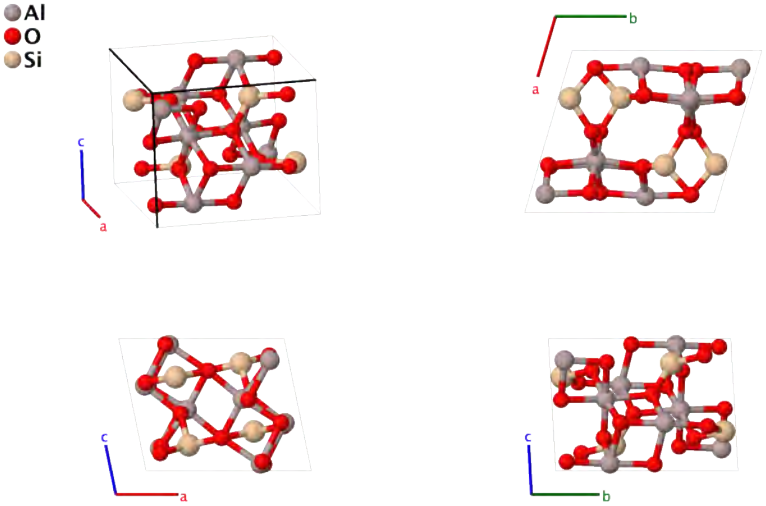
- S. V. Berger, *The Crystal Structure of Cobaltpyroborate*, Acta Chem. Scand. **4**, 1054–1065 (1950), doi:10.3891/acta.chem.scand.04-1054.

Geometry files:

- CIF: pp. 1500
- POSCAR: pp. 1501

Kyanite (Al_2SiO_5, $S0_1$) Structure: A2B5C_aP32_2_4i_10i_2i

http://aflow.org/prototype-encyclopedia/A2B5C_aP32_2_4i_10i_2i



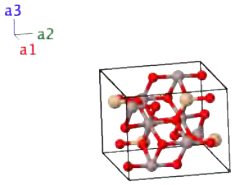
Prototype	:	Al_2O_5Si
AFLOW prototype label	:	A2B5C_aP32_2_4i_10i_2i
***Strukturbericht* designation**	:	$S0_1$
Pearson symbol	:	aP32
Space group number	:	2
Space group symbol	:	$P\bar{1}$
AFLOW prototype command	:	`aflow --proto=A2B5C_aP32_2_4i_10i_2i --params=`$a, b/a, c/a, \alpha, \beta, \gamma, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}$

- Three crystal polymorphs of Al_2SiO_5 have been characterized: kyanite ($S0_1$), space group $P\bar{1}$ #2, andalusite ($S0_2$), space group *Pnnm* #58, and sillimanite ($S0_3$), space group *Pnma* #62. All are characterized by chains of edge-sharing SiO_6 tetrahedra and Al octahedra.
- (Ewald, 1931) originally gave Kyanite the $H5_1$ *Strukturbericht* symbol, but this was changed to $S0_1$ in (Hermann, 1937).
- We use the ambient pressure data of (Yang, 1997).

Triclinic primitive vectors:

$$\mathbf{a}_1 = a\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\cos\gamma\,\hat{\mathbf{x}} + b\sin\gamma\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c_x\hat{\mathbf{x}} + c_y\,\hat{\mathbf{y}} + c_z\,\hat{\mathbf{z}}$$

$$c_x = c\,\cos\beta$$
$$c_y = c\,(\cos\alpha - \cos\beta\cos\gamma)/\sin\gamma$$
$$c_z = \sqrt{c^2 - c_x^2 - c_y^2}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	$(x_1a + y_1b\cos\gamma + z_1c_x)\,\hat{\mathbf{x}} + \left(y_1b\sin\gamma + z_1c_y\right)\hat{\mathbf{y}} + z_1c_z\,\hat{\mathbf{z}}$	$(2i)$	Al I
$\mathbf{B}_2$	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	$(-x_1a - y_1b\cos\gamma - z_1c_x)\,\hat{\mathbf{x}} + \left(-y_1b\sin\gamma - z_1c_y\right)\hat{\mathbf{y}} - z_1c_z\,\hat{\mathbf{z}}$	$(2i)$	Al I
$\mathbf{B}_3$	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	$(x_2a + y_2b\cos\gamma + z_2c_x)\,\hat{\mathbf{x}} + \left(y_2b\sin\gamma + z_2c_y\right)\hat{\mathbf{y}} + z_2c_z\,\hat{\mathbf{z}}$	$(2i)$	Al II
$\mathbf{B}_4$	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	$(-x_2a - y_2b\cos\gamma - z_2c_x)\,\hat{\mathbf{x}} + \left(-y_2b\sin\gamma - z_2c_y\right)\hat{\mathbf{y}} - z_2c_z\,\hat{\mathbf{z}}$	$(2i)$	Al II
$\mathbf{B}_5$	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$(x_3a + y_3b\cos\gamma + z_3c_x)\,\hat{\mathbf{x}} + \left(y_3b\sin\gamma + z_3c_y\right)\hat{\mathbf{y}} + z_3c_z\,\hat{\mathbf{z}}$	$(2i)$	Al III
$\mathbf{B}_6$	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	$(-x_3a - y_3b\cos\gamma - z_3c_x)\,\hat{\mathbf{x}} + \left(-y_3b\sin\gamma - z_3c_y\right)\hat{\mathbf{y}} - z_3c_z\,\hat{\mathbf{z}}$	$(2i)$	Al III
$\mathbf{B}_7$	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$(x_4a + y_4b\cos\gamma + z_4c_x)\,\hat{\mathbf{x}} + \left(y_4b\sin\gamma + z_4c_y\right)\hat{\mathbf{y}} + z_4c_z\,\hat{\mathbf{z}}$	$(2i)$	Al IV
$\mathbf{B}_8$	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	$(-x_4a - y_4b\cos\gamma - z_4c_x)\,\hat{\mathbf{x}} + \left(-y_4b\sin\gamma - z_4c_y\right)\hat{\mathbf{y}} - z_4c_z\,\hat{\mathbf{z}}$	$(2i)$	Al IV
$\mathbf{B}_9$	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	$(x_5a + y_5b\cos\gamma + z_5c_x)\,\hat{\mathbf{x}} + \left(y_5b\sin\gamma + z_5c_y\right)\hat{\mathbf{y}} + z_5c_z\,\hat{\mathbf{z}}$	$(2i)$	O I
$\mathbf{B}_{10}$	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	$(-x_5a - y_5b\cos\gamma - z_5c_x)\,\hat{\mathbf{x}} + \left(-y_5b\sin\gamma - z_5c_y\right)\hat{\mathbf{y}} - z_5c_z\,\hat{\mathbf{z}}$	$(2i)$	O I
$\mathbf{B}_{11}$	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	$(x_6a + y_6b\cos\gamma + z_6c_x)\,\hat{\mathbf{x}} + \left(y_6b\sin\gamma + z_6c_y\right)\hat{\mathbf{y}} + z_6c_z\,\hat{\mathbf{z}}$	$(2i)$	O II
$\mathbf{B}_{12}$	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	$(-x_6a - y_6b\cos\gamma - z_6c_x)\,\hat{\mathbf{x}} + \left(-y_6b\sin\gamma - z_6c_y\right)\hat{\mathbf{y}} - z_6c_z\,\hat{\mathbf{z}}$	$(2i)$	O II
$\mathbf{B}_{13}$	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	$(x_7a + y_7b\cos\gamma + z_7c_x)\,\hat{\mathbf{x}} + \left(y_7b\sin\gamma + z_7c_y\right)\hat{\mathbf{y}} + z_7c_z\,\hat{\mathbf{z}}$	$(2i)$	O III
$\mathbf{B}_{14}$	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	$(-x_7a - y_7b\cos\gamma - z_7c_x)\,\hat{\mathbf{x}} + \left(-y_7b\sin\gamma - z_7c_y\right)\hat{\mathbf{y}} - z_7c_z\,\hat{\mathbf{z}}$	$(2i)$	O III
$\mathbf{B}_{15}$	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	$(x_8a + y_8b\cos\gamma + z_8c_x)\,\hat{\mathbf{x}} + \left(y_8b\sin\gamma + z_8c_y\right)\hat{\mathbf{y}} + z_8c_z\,\hat{\mathbf{z}}$	$(2i)$	O IV
$\mathbf{B}_{16}$	$-x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	$(-x_8a - y_8b\cos\gamma - z_8c_x)\,\hat{\mathbf{x}} + \left(-y_8b\sin\gamma - z_8c_y\right)\hat{\mathbf{y}} - z_8c_z\,\hat{\mathbf{z}}$	$(2i)$	O IV

$$\mathbf{B}_{17} = x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 = (x_9a + y_9b\cos\gamma + z_9c_x)\,\hat{\mathbf{x}} + \left(y_9b\sin\gamma + z_9c_y\right)\hat{\mathbf{y}} + z_9c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{O V}$$

$$\mathbf{B}_{18} = -x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3 = (-x_9a - y_9b\cos\gamma - z_9c_x)\,\hat{\mathbf{x}} + \left(-y_9b\sin\gamma - z_9c_y\right)\hat{\mathbf{y}} - z_9c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{O V}$$

$$\mathbf{B}_{19} = x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 = (x_{10}a + y_{10}b\cos\gamma + z_{10}c_x)\,\hat{\mathbf{x}} + \left(y_{10}b\sin\gamma + z_{10}c_y\right)\hat{\mathbf{y}} + z_{10}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{O VI}$$

$$\mathbf{B}_{20} = -x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 = (-x_{10}a - y_{10}b\cos\gamma - z_{10}c_x)\,\hat{\mathbf{x}} + \left(-y_{10}b\sin\gamma - z_{10}c_y\right)\hat{\mathbf{y}} - z_{10}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{O VI}$$

$$\mathbf{B}_{21} = x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 = (x_{11}a + y_{11}b\cos\gamma + z_{11}c_x)\,\hat{\mathbf{x}} + \left(y_{11}b\sin\gamma + z_{11}c_y\right)\hat{\mathbf{y}} + z_{11}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{O VII}$$

$$\mathbf{B}_{22} = -x_{11}\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3 = (-x_{11}a - y_{11}b\cos\gamma - z_{11}c_x)\,\hat{\mathbf{x}} + \left(-y_{11}b\sin\gamma - z_{11}c_y\right)\hat{\mathbf{y}} - z_{11}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{O VII}$$

$$\mathbf{B}_{23} = x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 = (x_{12}a + y_{12}b\cos\gamma + z_{12}c_x)\,\hat{\mathbf{x}} + \left(y_{12}b\sin\gamma + z_{12}c_y\right)\hat{\mathbf{y}} + z_{12}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{O VIII}$$

$$\mathbf{B}_{24} = -x_{12}\,\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3 = (-x_{12}a - y_{12}b\cos\gamma - z_{12}c_x)\,\hat{\mathbf{x}} + \left(-y_{12}b\sin\gamma - z_{12}c_y\right)\hat{\mathbf{y}} - z_{12}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{O VIII}$$

$$\mathbf{B}_{25} = x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3 = (x_{13}a + y_{13}b\cos\gamma + z_{13}c_x)\,\hat{\mathbf{x}} + \left(y_{13}b\sin\gamma + z_{13}c_y\right)\hat{\mathbf{y}} + z_{13}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{O IX}$$

$$\mathbf{B}_{26} = -x_{13}\,\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 - z_{13}\,\mathbf{a}_3 = (-x_{13}a - y_{13}b\cos\gamma - z_{13}c_x)\,\hat{\mathbf{x}} + \left(-y_{13}b\sin\gamma - z_{13}c_y\right)\hat{\mathbf{y}} - z_{13}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{O IX}$$

$$\mathbf{B}_{27} = x_{14}\,\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3 = (x_{14}a + y_{14}b\cos\gamma + z_{14}c_x)\,\hat{\mathbf{x}} + \left(y_{14}b\sin\gamma + z_{14}c_y\right)\hat{\mathbf{y}} + z_{14}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{O X}$$

$$\mathbf{B}_{28} = -x_{14}\,\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 - z_{14}\,\mathbf{a}_3 = (-x_{14}a - y_{14}b\cos\gamma - z_{14}c_x)\,\hat{\mathbf{x}} + \left(-y_{14}b\sin\gamma - z_{14}c_y\right)\hat{\mathbf{y}} - z_{14}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{O X}$$

$$\mathbf{B}_{29} = x_{15}\,\mathbf{a}_1 + y_{15}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 = (x_{15}a + y_{15}b\cos\gamma + z_{15}c_x)\,\hat{\mathbf{x}} + \left(y_{15}b\sin\gamma + z_{15}c_y\right)\hat{\mathbf{y}} + z_{15}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{Si I}$$

$$\mathbf{B}_{30} = -x_{15}\,\mathbf{a}_1 - y_{15}\,\mathbf{a}_2 - z_{15}\,\mathbf{a}_3 = (-x_{15}a - y_{15}b\cos\gamma - z_{15}c_x)\,\hat{\mathbf{x}} + \left(-y_{15}b\sin\gamma - z_{15}c_y\right)\hat{\mathbf{y}} - z_{15}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{Si I}$$

$$\mathbf{B}_{31} = x_{16}\,\mathbf{a}_1 + y_{16}\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 = (x_{16}a + y_{16}b\cos\gamma + z_{16}c_x)\,\hat{\mathbf{x}} + \left(y_{16}b\sin\gamma + z_{16}c_y\right)\hat{\mathbf{y}} + z_{16}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{Si II}$$

$$\mathbf{B}_{32} = -x_{16}\,\mathbf{a}_1 - y_{16}\,\mathbf{a}_2 - z_{16}\,\mathbf{a}_3 = (-x_{16}a - y_{16}b\cos\gamma - z_{16}c_x)\,\hat{\mathbf{x}} + \left(-y_{16}b\sin\gamma - z_{16}c_y\right)\hat{\mathbf{y}} - z_{16}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{Si II}$$

References:

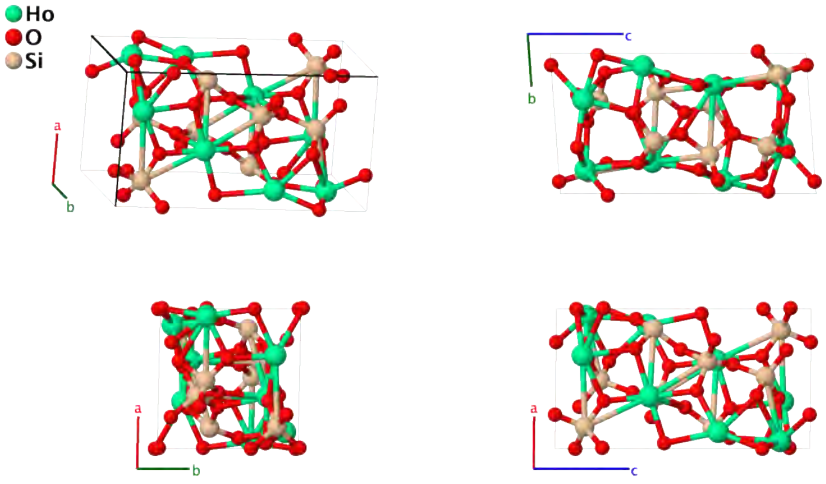
- H. Yang, R. T. Downs, L. W. Finger, R. M. Hazen, and C. T. Prewitt, *Compressibility and crystal structure of kyanite, Al_2SiO_5, at high pressure*, Am. Mineral. **82**, 467–474 (1997), doi:10.2138/am-1997-5-604.
- P. P. Ewald and C. Hermann, eds., *Strukturbericht 1913-1928* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1931).
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1501
- POSCAR: pp. 1501

α-$Ho_2Si_2O_7$ Structure: A2B7C2_aP44_2_4i_14i_4i

http://aflow.org/prototype-encyclopedia/A2B7C2_aP44_2_4i_14i_4i

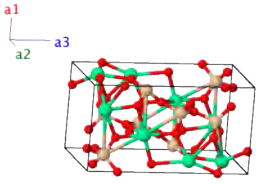


Prototype	:	$Ho_2O_7Si_2$
AFLOW prototype label	:	A2B7C2_aP44_2_4i_14i_4i
***Strukturbericht* designation**	:	None
Pearson symbol	:	aP44
Space group number	:	2
Space group symbol	:	$P\bar{1}$
AFLOW prototype command	:	`aflow --proto=A2B7C2_aP44_2_4i_14i_4i --params=`$a, b/a, c/a, \alpha, \beta, \gamma, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}, x_{17}, y_{17}, z_{17}, x_{18}, y_{18}, z_{18}, x_{19}, y_{19}, z_{19}, x_{20}, y_{20}, z_{20}, x_{21}, y_{21}, z_{21}, x_{22}, y_{22}, z_{22}$

Triclinic primitive vectors:

$$\mathbf{a}_1 = a\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\cos\gamma\,\hat{\mathbf{x}} + b\sin\gamma\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c_x\hat{\mathbf{x}} + c_y\,\hat{\mathbf{y}} + c_z\,\hat{\mathbf{z}}$$

$$c_x = c\cos\beta$$
$$c_y = c\,(\cos\alpha - \cos\beta\cos\gamma)/\sin\gamma$$
$$c_z = \sqrt{c^2 - c_x^2 - c_y^2}$$



Basis vectors:

	Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$ =	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1a + y_1b\cos\gamma + z_1c_x)\,\hat{\mathbf{x}} + \left(y_1b\sin\gamma + z_1c_y\right)\hat{\mathbf{y}} + z_1c_z\,\hat{\mathbf{z}}$	$(2i)$	Ho I
$\mathbf{B}_2$ =	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$(-x_1a - y_1b\cos\gamma - z_1c_x)\,\hat{\mathbf{x}} + \left(-y_1b\sin\gamma - z_1c_y\right)\hat{\mathbf{y}} - z_1c_z\,\hat{\mathbf{z}}$	$(2i)$	Ho I

$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2a + y_2b\cos\gamma + z_2c_x)\,\hat{\mathbf{x}} + \left(y_2b\sin\gamma + z_2c_y\right)\hat{\mathbf{y}} + z_2c_z\,\hat{\mathbf{z}}$	(2*i*)	Ho II
$\mathbf{B}_4$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2a - y_2b\cos\gamma - z_2c_x)\,\hat{\mathbf{x}} + \left(-y_2b\sin\gamma - z_2c_y\right)\hat{\mathbf{y}} - z_2c_z\,\hat{\mathbf{z}}$	(2*i*)	Ho II
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3a + y_3b\cos\gamma + z_3c_x)\,\hat{\mathbf{x}} + \left(y_3b\sin\gamma + z_3c_y\right)\hat{\mathbf{y}} + z_3c_z\,\hat{\mathbf{z}}$	(2*i*)	Ho III
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3a - y_3b\cos\gamma - z_3c_x)\,\hat{\mathbf{x}} + \left(-y_3b\sin\gamma - z_3c_y\right)\hat{\mathbf{y}} - z_3c_z\,\hat{\mathbf{z}}$	(2*i*)	Ho III
$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4a + y_4b\cos\gamma + z_4c_x)\,\hat{\mathbf{x}} + \left(y_4b\sin\gamma + z_4c_y\right)\hat{\mathbf{y}} + z_4c_z\,\hat{\mathbf{z}}$	(2*i*)	Ho IV
$\mathbf{B}_8$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4a - y_4b\cos\gamma - z_4c_x)\,\hat{\mathbf{x}} + \left(-y_4b\sin\gamma - z_4c_y\right)\hat{\mathbf{y}} - z_4c_z\,\hat{\mathbf{z}}$	(2*i*)	Ho IV
$\mathbf{B}_9$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5a + y_5b\cos\gamma + z_5c_x)\,\hat{\mathbf{x}} + \left(y_5b\sin\gamma + z_5c_y\right)\hat{\mathbf{y}} + z_5c_z\,\hat{\mathbf{z}}$	(2*i*)	O I
$\mathbf{B}_{10}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5a - y_5b\cos\gamma - z_5c_x)\,\hat{\mathbf{x}} + \left(-y_5b\sin\gamma - z_5c_y\right)\hat{\mathbf{y}} - z_5c_z\,\hat{\mathbf{z}}$	(2*i*)	O I
$\mathbf{B}_{11}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6a + y_6b\cos\gamma + z_6c_x)\,\hat{\mathbf{x}} + \left(y_6b\sin\gamma + z_6c_y\right)\hat{\mathbf{y}} + z_6c_z\,\hat{\mathbf{z}}$	(2*i*)	O II
$\mathbf{B}_{12}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6a - y_6b\cos\gamma - z_6c_x)\,\hat{\mathbf{x}} + \left(-y_6b\sin\gamma - z_6c_y\right)\hat{\mathbf{y}} - z_6c_z\,\hat{\mathbf{z}}$	(2*i*)	O II
$\mathbf{B}_{13}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7a + y_7b\cos\gamma + z_7c_x)\,\hat{\mathbf{x}} + \left(y_7b\sin\gamma + z_7c_y\right)\hat{\mathbf{y}} + z_7c_z\,\hat{\mathbf{z}}$	(2*i*)	O III
$\mathbf{B}_{14}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7a - y_7b\cos\gamma - z_7c_x)\,\hat{\mathbf{x}} + \left(-y_7b\sin\gamma - z_7c_y\right)\hat{\mathbf{y}} - z_7c_z\,\hat{\mathbf{z}}$	(2*i*)	O III
$\mathbf{B}_{15}$	=	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8a + y_8b\cos\gamma + z_8c_x)\,\hat{\mathbf{x}} + \left(y_8b\sin\gamma + z_8c_y\right)\hat{\mathbf{y}} + z_8c_z\,\hat{\mathbf{z}}$	(2*i*)	O IV
$\mathbf{B}_{16}$	=	$-x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8a - y_8b\cos\gamma - z_8c_x)\,\hat{\mathbf{x}} + \left(-y_8b\sin\gamma - z_8c_y\right)\hat{\mathbf{y}} - z_8c_z\,\hat{\mathbf{z}}$	(2*i*)	O IV
$\mathbf{B}_{17}$	=	$x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$(x_9a + y_9b\cos\gamma + z_9c_x)\,\hat{\mathbf{x}} + \left(y_9b\sin\gamma + z_9c_y\right)\hat{\mathbf{y}} + z_9c_z\,\hat{\mathbf{z}}$	(2*i*)	O V
$\mathbf{B}_{18}$	=	$-x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$(-x_9a - y_9b\cos\gamma - z_9c_x)\,\hat{\mathbf{x}} + \left(-y_9b\sin\gamma - z_9c_y\right)\hat{\mathbf{y}} - z_9c_z\,\hat{\mathbf{z}}$	(2*i*)	O V
$\mathbf{B}_{19}$	=	$x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$(x_{10}a + y_{10}b\cos\gamma + z_{10}c_x)\,\hat{\mathbf{x}} + \left(y_{10}b\sin\gamma + z_{10}c_y\right)\hat{\mathbf{y}} + z_{10}c_z\,\hat{\mathbf{z}}$	(2*i*)	O VI
$\mathbf{B}_{20}$	=	$-x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3$	=	$(-x_{10}a - y_{10}b\cos\gamma - z_{10}c_x)\,\hat{\mathbf{x}} + \left(-y_{10}b\sin\gamma - z_{10}c_y\right)\hat{\mathbf{y}} - z_{10}c_z\,\hat{\mathbf{z}}$	(2*i*)	O VI
$\mathbf{B}_{21}$	=	$x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$(x_{11}a + y_{11}b\cos\gamma + z_{11}c_x)\,\hat{\mathbf{x}} + \left(y_{11}b\sin\gamma + z_{11}c_y\right)\hat{\mathbf{y}} + z_{11}c_z\,\hat{\mathbf{z}}$	(2*i*)	O VII
$\mathbf{B}_{22}$	=	$-x_{11}\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3$	=	$(-x_{11}a - y_{11}b\cos\gamma - z_{11}c_x)\,\hat{\mathbf{x}} + \left(-y_{11}b\sin\gamma - z_{11}c_y\right)\hat{\mathbf{y}} - z_{11}c_z\,\hat{\mathbf{z}}$	(2*i*)	O VII
$\mathbf{B}_{23}$	=	$x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3$	=	$(x_{12}a + y_{12}b\cos\gamma + z_{12}c_x)\,\hat{\mathbf{x}} + \left(y_{12}b\sin\gamma + z_{12}c_y\right)\hat{\mathbf{y}} + z_{12}c_z\,\hat{\mathbf{z}}$	(2*i*)	O VIII
$\mathbf{B}_{24}$	=	$-x_{12}\,\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3$	=	$(-x_{12}a - y_{12}b\cos\gamma - z_{12}c_x)\,\hat{\mathbf{x}} + \left(-y_{12}b\sin\gamma - z_{12}c_y\right)\hat{\mathbf{y}} - z_{12}c_z\,\hat{\mathbf{z}}$	(2*i*)	O VIII

$$\mathbf{B}_{25} = x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3 = (x_{13}a + y_{13}b\cos\gamma + z_{13}c_x)\,\hat{\mathbf{x}} + \left(y_{13}b\sin\gamma + z_{13}c_y\right)\hat{\mathbf{y}} + z_{13}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{O IX}$$

$$\mathbf{B}_{26} = -x_{13}\,\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 - z_{13}\,\mathbf{a}_3 = (-x_{13}a - y_{13}b\cos\gamma - z_{13}c_x)\,\hat{\mathbf{x}} + \left(-y_{13}b\sin\gamma - z_{13}c_y\right)\hat{\mathbf{y}} - z_{13}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{O IX}$$

$$\mathbf{B}_{27} = x_{14}\,\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3 = (x_{14}a + y_{14}b\cos\gamma + z_{14}c_x)\,\hat{\mathbf{x}} + \left(y_{14}b\sin\gamma + z_{14}c_y\right)\hat{\mathbf{y}} + z_{14}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{O X}$$

$$\mathbf{B}_{28} = -x_{14}\,\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 - z_{14}\,\mathbf{a}_3 = (-x_{14}a - y_{14}b\cos\gamma - z_{14}c_x)\,\hat{\mathbf{x}} + \left(-y_{14}b\sin\gamma - z_{14}c_y\right)\hat{\mathbf{y}} - z_{14}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{O X}$$

$$\mathbf{B}_{29} = x_{15}\,\mathbf{a}_1 + y_{15}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 = (x_{15}a + y_{15}b\cos\gamma + z_{15}c_x)\,\hat{\mathbf{x}} + \left(y_{15}b\sin\gamma + z_{15}c_y\right)\hat{\mathbf{y}} + z_{15}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{O XI}$$

$$\mathbf{B}_{30} = -x_{15}\,\mathbf{a}_1 - y_{15}\,\mathbf{a}_2 - z_{15}\,\mathbf{a}_3 = (-x_{15}a - y_{15}b\cos\gamma - z_{15}c_x)\,\hat{\mathbf{x}} + \left(-y_{15}b\sin\gamma - z_{15}c_y\right)\hat{\mathbf{y}} - z_{15}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{O XI}$$

$$\mathbf{B}_{31} = x_{16}\,\mathbf{a}_1 + y_{16}\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 = (x_{16}a + y_{16}b\cos\gamma + z_{16}c_x)\,\hat{\mathbf{x}} + \left(y_{16}b\sin\gamma + z_{16}c_y\right)\hat{\mathbf{y}} + z_{16}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{O XII}$$

$$\mathbf{B}_{32} = -x_{16}\,\mathbf{a}_1 - y_{16}\,\mathbf{a}_2 - z_{16}\,\mathbf{a}_3 = (-x_{16}a - y_{16}b\cos\gamma - z_{16}c_x)\,\hat{\mathbf{x}} + \left(-y_{16}b\sin\gamma - z_{16}c_y\right)\hat{\mathbf{y}} - z_{16}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{O XII}$$

$$\mathbf{B}_{33} = x_{17}\,\mathbf{a}_1 + y_{17}\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 = (x_{17}a + y_{17}b\cos\gamma + z_{17}c_x)\,\hat{\mathbf{x}} + \left(y_{17}b\sin\gamma + z_{17}c_y\right)\hat{\mathbf{y}} + z_{17}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{O XIII}$$

$$\mathbf{B}_{34} = -x_{17}\,\mathbf{a}_1 - y_{17}\,\mathbf{a}_2 - z_{17}\,\mathbf{a}_3 = (-x_{17}a - y_{17}b\cos\gamma - z_{17}c_x)\,\hat{\mathbf{x}} + \left(-y_{17}b\sin\gamma - z_{17}c_y\right)\hat{\mathbf{y}} - z_{17}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{O XIII}$$

$$\mathbf{B}_{35} = x_{18}\,\mathbf{a}_1 + y_{18}\,\mathbf{a}_2 + z_{18}\,\mathbf{a}_3 = (x_{18}a + y_{18}b\cos\gamma + z_{18}c_x)\,\hat{\mathbf{x}} + \left(y_{18}b\sin\gamma + z_{18}c_y\right)\hat{\mathbf{y}} + z_{18}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{O XIV}$$

$$\mathbf{B}_{36} = -x_{18}\,\mathbf{a}_1 - y_{18}\,\mathbf{a}_2 - z_{18}\,\mathbf{a}_3 = (-x_{18}a - y_{18}b\cos\gamma - z_{18}c_x)\,\hat{\mathbf{x}} + \left(-y_{18}b\sin\gamma - z_{18}c_y\right)\hat{\mathbf{y}} - z_{18}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{O XIV}$$

$$\mathbf{B}_{37} = x_{19}\,\mathbf{a}_1 + y_{19}\,\mathbf{a}_2 + z_{19}\,\mathbf{a}_3 = (x_{19}a + y_{19}b\cos\gamma + z_{19}c_x)\,\hat{\mathbf{x}} + \left(y_{19}b\sin\gamma + z_{19}c_y\right)\hat{\mathbf{y}} + z_{19}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{Si I}$$

$$\mathbf{B}_{38} = -x_{19}\,\mathbf{a}_1 - y_{19}\,\mathbf{a}_2 - z_{19}\,\mathbf{a}_3 = (-x_{19}a - y_{19}b\cos\gamma - z_{19}c_x)\,\hat{\mathbf{x}} + \left(-y_{19}b\sin\gamma - z_{19}c_y\right)\hat{\mathbf{y}} - z_{19}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{Si I}$$

$$\mathbf{B}_{39} = x_{20}\,\mathbf{a}_1 + y_{20}\,\mathbf{a}_2 + z_{20}\,\mathbf{a}_3 = (x_{20}a + y_{20}b\cos\gamma + z_{20}c_x)\,\hat{\mathbf{x}} + \left(y_{20}b\sin\gamma + z_{20}c_y\right)\hat{\mathbf{y}} + z_{20}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{Si II}$$

$$\mathbf{B}_{40} = -x_{20}\,\mathbf{a}_1 - y_{20}\,\mathbf{a}_2 - z_{20}\,\mathbf{a}_3 = (-x_{20}a - y_{20}b\cos\gamma - z_{20}c_x)\,\hat{\mathbf{x}} + \left(-y_{20}b\sin\gamma - z_{20}c_y\right)\hat{\mathbf{y}} - z_{20}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{Si II}$$

$$\mathbf{B}_{41} = x_{21}\,\mathbf{a}_1 + y_{21}\,\mathbf{a}_2 + z_{21}\,\mathbf{a}_3 = (x_{21}a + y_{21}b\cos\gamma + z_{21}c_x)\,\hat{\mathbf{x}} + \left(y_{21}b\sin\gamma + z_{21}c_y\right)\hat{\mathbf{y}} + z_{21}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{Si III}$$

$$\mathbf{B}_{42} = -x_{21}\,\mathbf{a}_1 - y_{21}\,\mathbf{a}_2 - z_{21}\,\mathbf{a}_3 = (-x_{21}a - y_{21}b\cos\gamma - z_{21}c_x)\,\hat{\mathbf{x}} + \left(-y_{21}b\sin\gamma - z_{21}c_y\right)\hat{\mathbf{y}} - z_{21}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{Si III}$$

$$\mathbf{B}_{43} = x_{22}\,\mathbf{a}_1 + y_{22}\,\mathbf{a}_2 + z_{22}\,\mathbf{a}_3 = (x_{22}a + y_{22}b\cos\gamma + z_{22}c_x)\,\hat{\mathbf{x}} + \left(y_{22}b\sin\gamma + z_{22}c_y\right)\hat{\mathbf{y}} + z_{22}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{Si IV}$$

$$\mathbf{B}_{44} = -x_{22}\,\mathbf{a}_1 - y_{22}\,\mathbf{a}_2 - z_{22}\,\mathbf{a}_3 = (-x_{22}a - y_{22}b\cos\gamma - z_{22}c_x)\,\hat{\mathbf{x}} + \left(-y_{22}b\sin\gamma - z_{22}c_y\right)\hat{\mathbf{y}} - z_{22}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{Si IV}$$

References:

- J. Felsche, *A new silicate structure containing linear [Si_3O_{10}] groups*, Naturwissenschaften **59**, 35–36 (1972), doi:10.1007/BF00594623.

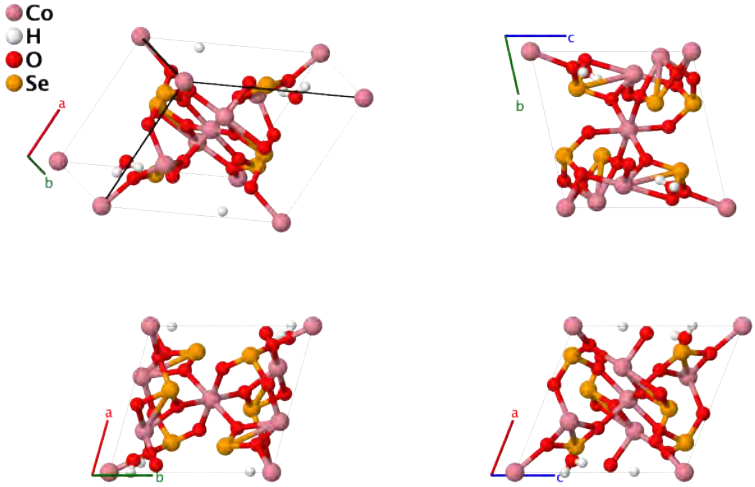
Found in:

- A. I. Becerro and A. Escudero, *Revision of the crystallographic data of polymorphic $Y_2Si_2O_7$ and Y_2SiO_5 compounds*, Phase Transit. **77**, 1093–1102 (2004), doi:10.1080/01411590412331282814.

Geometry files:
- CIF: pp. 1501
- POSCAR: pp. 1502

$Co_3(SeO_3)_3{\cdot}H_2O$ Structure: A3B2C10D3_aP36_2_ah2i_2i_10i_3i

http://aflow.org/prototype-encyclopedia/A3B2C10D3_aP36_2_ah2i_2i_10i_3i



Prototype	:	$Co_3H_2O_{10}Se_3$
AFLOW prototype label	:	A3B2C10D3_aP36_2_ah2i_2i_10i_3i
***Strukturbericht* designation**	:	None
Pearson symbol	:	aP36
Space group number	:	2
Space group symbol	:	$P\bar{1}$
AFLOW prototype command	:	`aflow --proto=A3B2C10D3_aP36_2_ah2i_2i_10i_3i --params=`$a, b/a, c/a, \alpha, \beta, \gamma, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}, x_{17}, y_{17}, z_{17}, x_{18}, y_{18}, z_{18}, x_{19}, y_{19}, z_{19}$

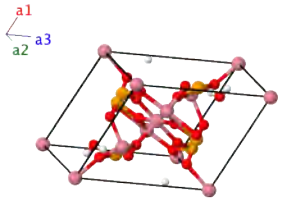
Other compounds with this structure

- $Mn_3(SeO_3)_3{\cdot}H_2O$ and $Ni_3(SeO_3)_3{\cdot}H_2O$

Triclinic primitive vectors:

$$\mathbf{a}_1 = a\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\cos\gamma\,\hat{\mathbf{x}} + b\sin\gamma\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c_x\hat{\mathbf{x}} + c_y\,\hat{\mathbf{y}} + c_z\,\hat{\mathbf{z}}$$

$$c_x = c\cos\beta$$
$$c_y = c\,(\cos\alpha - \cos\beta\cos\gamma)/\sin\gamma$$
$$c_z = \sqrt{c^2 - c_x^2 - c_y^2}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(1*a*)	Co I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(a + b\cos\gamma + c_x)\,\hat{\mathbf{x}} + \frac{1}{2}\left(b\sin\gamma + c_y\right)\hat{\mathbf{y}} + \frac{1}{2}c_z\,\hat{\mathbf{z}}$	(1*h*)	Co II
$\mathbf{B}_3$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3a + y_3b\cos\gamma + z_3c_x)\,\hat{\mathbf{x}} + \left(y_3b\sin\gamma + z_3c_y\right)\hat{\mathbf{y}} + z_3c_z\,\hat{\mathbf{z}}$	(2*i*)	Co III
$\mathbf{B}_4$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3a - y_3b\cos\gamma - z_3c_x)\,\hat{\mathbf{x}} + \left(-y_3b\sin\gamma - z_3c_y\right)\hat{\mathbf{y}} - z_3c_z\,\hat{\mathbf{z}}$	(2*i*)	Co III
$\mathbf{B}_5$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4a + y_4b\cos\gamma + z_4c_x)\,\hat{\mathbf{x}} + \left(y_4b\sin\gamma + z_4c_y\right)\hat{\mathbf{y}} + z_4c_z\,\hat{\mathbf{z}}$	(2*i*)	Co IV
$\mathbf{B}_6$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4a - y_4b\cos\gamma - z_4c_x)\,\hat{\mathbf{x}} + \left(-y_4b\sin\gamma - z_4c_y\right)\hat{\mathbf{y}} - z_4c_z\,\hat{\mathbf{z}}$	(2*i*)	Co IV
$\mathbf{B}_7$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5a + y_5b\cos\gamma + z_5c_x)\,\hat{\mathbf{x}} + \left(y_5b\sin\gamma + z_5c_y\right)\hat{\mathbf{y}} + z_5c_z\,\hat{\mathbf{z}}$	(2*i*)	H I
$\mathbf{B}_8$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5a - y_5b\cos\gamma - z_5c_x)\,\hat{\mathbf{x}} + \left(-y_5b\sin\gamma - z_5c_y\right)\hat{\mathbf{y}} - z_5c_z\,\hat{\mathbf{z}}$	(2*i*)	H I
$\mathbf{B}_9$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6a + y_6b\cos\gamma + z_6c_x)\,\hat{\mathbf{x}} + \left(y_6b\sin\gamma + z_6c_y\right)\hat{\mathbf{y}} + z_6c_z\,\hat{\mathbf{z}}$	(2*i*)	H II
$\mathbf{B}_{10}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6a - y_6b\cos\gamma - z_6c_x)\,\hat{\mathbf{x}} + \left(-y_6b\sin\gamma - z_6c_y\right)\hat{\mathbf{y}} - z_6c_z\,\hat{\mathbf{z}}$	(2*i*)	H II
$\mathbf{B}_{11}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7a + y_7b\cos\gamma + z_7c_x)\,\hat{\mathbf{x}} + \left(y_7b\sin\gamma + z_7c_y\right)\hat{\mathbf{y}} + z_7c_z\,\hat{\mathbf{z}}$	(2*i*)	O I
$\mathbf{B}_{12}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7a - y_7b\cos\gamma - z_7c_x)\,\hat{\mathbf{x}} + \left(-y_7b\sin\gamma - z_7c_y\right)\hat{\mathbf{y}} - z_7c_z\,\hat{\mathbf{z}}$	(2*i*)	O I
$\mathbf{B}_{13}$	=	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8a + y_8b\cos\gamma + z_8c_x)\,\hat{\mathbf{x}} + \left(y_8b\sin\gamma + z_8c_y\right)\hat{\mathbf{y}} + z_8c_z\,\hat{\mathbf{z}}$	(2*i*)	O II
$\mathbf{B}_{14}$	=	$-x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8a - y_8b\cos\gamma - z_8c_x)\,\hat{\mathbf{x}} + \left(-y_8b\sin\gamma - z_8c_y\right)\hat{\mathbf{y}} - z_8c_z\,\hat{\mathbf{z}}$	(2*i*)	O II
$\mathbf{B}_{15}$	=	$x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$(x_9a + y_9b\cos\gamma + z_9c_x)\,\hat{\mathbf{x}} + \left(y_9b\sin\gamma + z_9c_y\right)\hat{\mathbf{y}} + z_9c_z\,\hat{\mathbf{z}}$	(2*i*)	O III
$\mathbf{B}_{16}$	=	$-x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$(-x_9a - y_9b\cos\gamma - z_9c_x)\,\hat{\mathbf{x}} + \left(-y_9b\sin\gamma - z_9c_y\right)\hat{\mathbf{y}} - z_9c_z\,\hat{\mathbf{z}}$	(2*i*)	O III
$\mathbf{B}_{17}$	=	$x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$(x_{10}a + y_{10}b\cos\gamma + z_{10}c_x)\,\hat{\mathbf{x}} + \left(y_{10}b\sin\gamma + z_{10}c_y\right)\hat{\mathbf{y}} + z_{10}c_z\,\hat{\mathbf{z}}$	(2*i*)	O IV
$\mathbf{B}_{18}$	=	$-x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3$	=	$(-x_{10}a - y_{10}b\cos\gamma - z_{10}c_x)\,\hat{\mathbf{x}} + \left(-y_{10}b\sin\gamma - z_{10}c_y\right)\hat{\mathbf{y}} - z_{10}c_z\,\hat{\mathbf{z}}$	(2*i*)	O IV
$\mathbf{B}_{19}$	=	$x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$(x_{11}a + y_{11}b\cos\gamma + z_{11}c_x)\,\hat{\mathbf{x}} + \left(y_{11}b\sin\gamma + z_{11}c_y\right)\hat{\mathbf{y}} + z_{11}c_z\,\hat{\mathbf{z}}$	(2*i*)	O V
$\mathbf{B}_{20}$	=	$-x_{11}\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3$	=	$(-x_{11}a - y_{11}b\cos\gamma - z_{11}c_x)\,\hat{\mathbf{x}} + \left(-y_{11}b\sin\gamma - z_{11}c_y\right)\hat{\mathbf{y}} - z_{11}c_z\,\hat{\mathbf{z}}$	(2*i*)	O V
$\mathbf{B}_{21}$	=	$x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3$	=	$(x_{12}a + y_{12}b\cos\gamma + z_{12}c_x)\,\hat{\mathbf{x}} + \left(y_{12}b\sin\gamma + z_{12}c_y\right)\hat{\mathbf{y}} + z_{12}c_z\,\hat{\mathbf{z}}$	(2*i*)	O VI
$\mathbf{B}_{22}$	=	$-x_{12}\,\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3$	=	$(-x_{12}a - y_{12}b\cos\gamma - z_{12}c_x)\,\hat{\mathbf{x}} + \left(-y_{12}b\sin\gamma - z_{12}c_y\right)\hat{\mathbf{y}} - z_{12}c_z\,\hat{\mathbf{z}}$	(2*i*)	O VI

$$\mathbf{B}_{23} = x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3 = (x_{13}a + y_{13}b\cos\gamma + z_{13}c_x)\,\hat{\mathbf{x}} + \left(y_{13}b\sin\gamma + z_{13}c_y\right)\hat{\mathbf{y}} + z_{13}c_z\,\hat{\mathbf{z}}$$ (2*i*) O VII

$$\mathbf{B}_{24} = -x_{13}\,\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 - z_{13}\,\mathbf{a}_3 = (-x_{13}a - y_{13}b\cos\gamma - z_{13}c_x)\,\hat{\mathbf{x}} + \left(-y_{13}b\sin\gamma - z_{13}c_y\right)\hat{\mathbf{y}} - z_{13}c_z\,\hat{\mathbf{z}}$$ (2*i*) O VII

$$\mathbf{B}_{25} = x_{14}\,\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3 = (x_{14}a + y_{14}b\cos\gamma + z_{14}c_x)\,\hat{\mathbf{x}} + \left(y_{14}b\sin\gamma + z_{14}c_y\right)\hat{\mathbf{y}} + z_{14}c_z\,\hat{\mathbf{z}}$$ (2*i*) O VIII

$$\mathbf{B}_{26} = -x_{14}\,\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 - z_{14}\,\mathbf{a}_3 = (-x_{14}a - y_{14}b\cos\gamma - z_{14}c_x)\,\hat{\mathbf{x}} + \left(-y_{14}b\sin\gamma - z_{14}c_y\right)\hat{\mathbf{y}} - z_{14}c_z\,\hat{\mathbf{z}}$$ (2*i*) O VIII

$$\mathbf{B}_{27} = x_{15}\,\mathbf{a}_1 + y_{15}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 = (x_{15}a + y_{15}b\cos\gamma + z_{15}c_x)\,\hat{\mathbf{x}} + \left(y_{15}b\sin\gamma + z_{15}c_y\right)\hat{\mathbf{y}} + z_{15}c_z\,\hat{\mathbf{z}}$$ (2*i*) O IX

$$\mathbf{B}_{28} = -x_{15}\,\mathbf{a}_1 - y_{15}\,\mathbf{a}_2 - z_{15}\,\mathbf{a}_3 = (-x_{15}a - y_{15}b\cos\gamma - z_{15}c_x)\,\hat{\mathbf{x}} + \left(-y_{15}b\sin\gamma - z_{15}c_y\right)\hat{\mathbf{y}} - z_{15}c_z\,\hat{\mathbf{z}}$$ (2*i*) O IX

$$\mathbf{B}_{29} = x_{16}\,\mathbf{a}_1 + y_{16}\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 = (x_{16}a + y_{16}b\cos\gamma + z_{16}c_x)\,\hat{\mathbf{x}} + \left(y_{16}b\sin\gamma + z_{16}c_y\right)\hat{\mathbf{y}} + z_{16}c_z\,\hat{\mathbf{z}}$$ (2*i*) O X

$$\mathbf{B}_{30} = -x_{16}\,\mathbf{a}_1 - y_{16}\,\mathbf{a}_2 - z_{16}\,\mathbf{a}_3 = (-x_{16}a - y_{16}b\cos\gamma - z_{16}c_x)\,\hat{\mathbf{x}} + \left(-y_{16}b\sin\gamma - z_{16}c_y\right)\hat{\mathbf{y}} - z_{16}c_z\,\hat{\mathbf{z}}$$ (2*i*) O X

$$\mathbf{B}_{31} = x_{17}\,\mathbf{a}_1 + y_{17}\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 = (x_{17}a + y_{17}b\cos\gamma + z_{17}c_x)\,\hat{\mathbf{x}} + \left(y_{17}b\sin\gamma + z_{17}c_y\right)\hat{\mathbf{y}} + z_{17}c_z\,\hat{\mathbf{z}}$$ (2*i*) Se I

$$\mathbf{B}_{32} = -x_{17}\,\mathbf{a}_1 - y_{17}\,\mathbf{a}_2 - z_{17}\,\mathbf{a}_3 = (-x_{17}a - y_{17}b\cos\gamma - z_{17}c_x)\,\hat{\mathbf{x}} + \left(-y_{17}b\sin\gamma - z_{17}c_y\right)\hat{\mathbf{y}} - z_{17}c_z\,\hat{\mathbf{z}}$$ (2*i*) Se I

$$\mathbf{B}_{33} = x_{18}\,\mathbf{a}_1 + y_{18}\,\mathbf{a}_2 + z_{18}\,\mathbf{a}_3 = (x_{18}a + y_{18}b\cos\gamma + z_{18}c_x)\,\hat{\mathbf{x}} + \left(y_{18}b\sin\gamma + z_{18}c_y\right)\hat{\mathbf{y}} + z_{18}c_z\,\hat{\mathbf{z}}$$ (2*i*) Se II

$$\mathbf{B}_{34} = -x_{18}\,\mathbf{a}_1 - y_{18}\,\mathbf{a}_2 - z_{18}\,\mathbf{a}_3 = (-x_{18}a - y_{18}b\cos\gamma - z_{18}c_x)\,\hat{\mathbf{x}} + \left(-y_{18}b\sin\gamma - z_{18}c_y\right)\hat{\mathbf{y}} - z_{18}c_z\,\hat{\mathbf{z}}$$ (2*i*) Se II

$$\mathbf{B}_{35} = x_{19}\,\mathbf{a}_1 + y_{19}\,\mathbf{a}_2 + z_{19}\,\mathbf{a}_3 = (x_{19}a + y_{19}b\cos\gamma + z_{19}c_x)\,\hat{\mathbf{x}} + \left(y_{19}b\sin\gamma + z_{19}c_y\right)\hat{\mathbf{y}} + z_{19}c_z\,\hat{\mathbf{z}}$$ (2*i*) Se III

$$\mathbf{B}_{36} = -x_{19}\,\mathbf{a}_1 - y_{19}\,\mathbf{a}_2 - z_{19}\,\mathbf{a}_3 = (-x_{19}a - y_{19}b\cos\gamma - z_{19}c_x)\,\hat{\mathbf{x}} + \left(-y_{19}b\sin\gamma - z_{19}c_y\right)\hat{\mathbf{y}} - z_{19}c_z\,\hat{\mathbf{z}}$$ (2*i*) Se III

References:

- M. Wildner, *Crystal structures of $Co_3(SeO_3)_3{\cdot}H_2O$ and $Ni_3(SeO_3)_3{\cdot}H_2O$, two new isotypic compounds*, Monatshefte für Chemie - Chemical Monthly **122**, 585–594 (1991), doi:10.1007/BF00811457.

Found in:

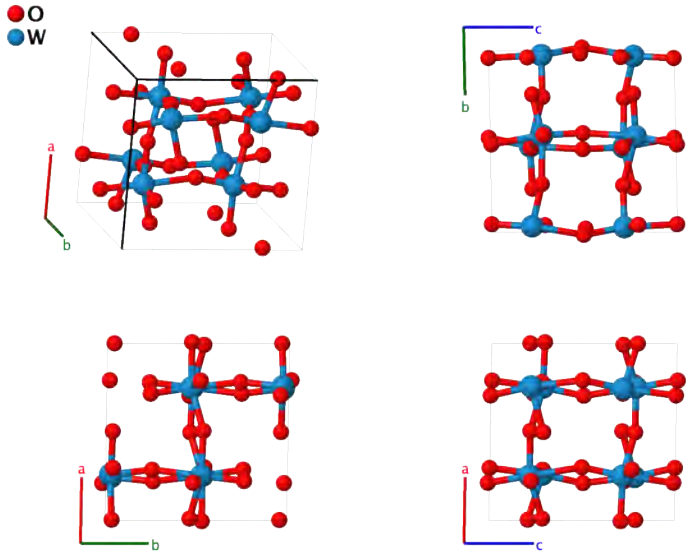
- K. M. Taddei, L. D. Sanjeewa, J. Xing, Q. Zhang, D. Parker, A. Podlesnyak, C. dela Cruz, and A. S. Sefat, *Tunable magnetic order in low-symmetry SeO_3 ligand linked $TM_3(SeO_3)_3H_2O$ (TM = Mn, Co and Ni) compounds*, Phys. Rev. Mater. **4**, 024410 (2020), doi:10.1103/PhysRevMaterials.4.024410.

Geometry files:

- CIF: pp. 1502
- POSCAR: pp. 1503

δ-WO_3 Structure: A3B_aP32_2_12i_4i

http://aflow.org/prototype-encyclopedia/A3B_aP32_2_12i_4i



Prototype	:	O_3W
AFLOW prototype label	:	A3B_aP32_2_12i_4i
***Strukturbericht* designation**	:	None
Pearson symbol	:	aP32
Space group number	:	2
Space group symbol	:	$P\bar{1}$
AFLOW prototype command	:	`aflow --proto=A3B_aP32_2_12i_4i --params=`$a, b/a, c/a, \alpha, \beta, \gamma, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}$

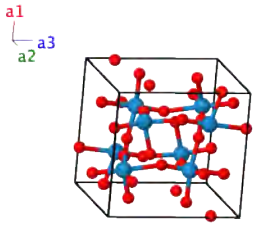
- All stable phases of WO_3 are distortions of the cubic α-ReO_3 ($D0_9$) phase. Based on (Woodward, 1997 and Vogt, 1999), the known stable phases and their approximate temperature ranges are:
  - α-WO_3 (1010-1170K) (Vogt, 1999)
  - β-WO_3 (600-1170K) (Vogt, 1999)
  - γ-WO_3 (290-600K) (Vogt, 1999)
  - δ-WO_3 (230-290K) (Diehl, 1978), this structure
  - ϵ-WO_3 (below 23K) (Woodward, 1997)

  Woodward's caption says "The transition temperatures display large hysteresis effects and universal agreement is not found in the literature."
- In addition, several other structures have been proposed and/or found:
  - The original $D0_{10}$ structure (Bräkken, 1931), (Hermann, 1937) superseded by δ-WO_3
  - Original β-WO_3 (Salje, 1977)
  - Hexagonal WO_3 (Gerand, 1979) (metastable)

Triclinic primitive vectors:

$$\begin{aligned}
\mathbf{a}_1 &= a\hat{\mathbf{x}} \\
\mathbf{a}_2 &= b\cos\gamma\,\hat{\mathbf{x}} + b\sin\gamma\,\hat{\mathbf{y}} \\
\mathbf{a}_3 &= c_x\hat{\mathbf{x}} + c_y\,\hat{\mathbf{y}} + c_z\,\hat{\mathbf{z}}
\end{aligned}$$

$$\begin{aligned}
c_x &= c\,\cos\beta \\
c_y &= c\,(\cos\alpha - \cos\beta\cos\gamma)/\sin\gamma \\
c_z &= \sqrt{c^2 - c_x^2 - c_y^2}
\end{aligned}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	$(x_1a + y_1b\cos\gamma + z_1c_x)\,\hat{\mathbf{x}} + \left(y_1b\sin\gamma + z_1c_y\right)\hat{\mathbf{y}} + z_1c_z\,\hat{\mathbf{z}}$	$(2i)$	O I
$\mathbf{B}_2$	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	$(-x_1a - y_1b\cos\gamma - z_1c_x)\,\hat{\mathbf{x}} + \left(-y_1b\sin\gamma - z_1c_y\right)\hat{\mathbf{y}} - z_1c_z\,\hat{\mathbf{z}}$	$(2i)$	O I
$\mathbf{B}_3$	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	$(x_2a + y_2b\cos\gamma + z_2c_x)\,\hat{\mathbf{x}} + \left(y_2b\sin\gamma + z_2c_y\right)\hat{\mathbf{y}} + z_2c_z\,\hat{\mathbf{z}}$	$(2i)$	O II
$\mathbf{B}_4$	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	$(-x_2a - y_2b\cos\gamma - z_2c_x)\,\hat{\mathbf{x}} + \left(-y_2b\sin\gamma - z_2c_y\right)\hat{\mathbf{y}} - z_2c_z\,\hat{\mathbf{z}}$	$(2i)$	O II
$\mathbf{B}_5$	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$(x_3a + y_3b\cos\gamma + z_3c_x)\,\hat{\mathbf{x}} + \left(y_3b\sin\gamma + z_3c_y\right)\hat{\mathbf{y}} + z_3c_z\,\hat{\mathbf{z}}$	$(2i)$	O III
$\mathbf{B}_6$	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	$(-x_3a - y_3b\cos\gamma - z_3c_x)\,\hat{\mathbf{x}} + \left(-y_3b\sin\gamma - z_3c_y\right)\hat{\mathbf{y}} - z_3c_z\,\hat{\mathbf{z}}$	$(2i)$	O III
$\mathbf{B}_7$	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$(x_4a + y_4b\cos\gamma + z_4c_x)\,\hat{\mathbf{x}} + \left(y_4b\sin\gamma + z_4c_y\right)\hat{\mathbf{y}} + z_4c_z\,\hat{\mathbf{z}}$	$(2i)$	O IV
$\mathbf{B}_8$	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	$(-x_4a - y_4b\cos\gamma - z_4c_x)\,\hat{\mathbf{x}} + \left(-y_4b\sin\gamma - z_4c_y\right)\hat{\mathbf{y}} - z_4c_z\,\hat{\mathbf{z}}$	$(2i)$	O IV
$\mathbf{B}_9$	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	$(x_5a + y_5b\cos\gamma + z_5c_x)\,\hat{\mathbf{x}} + \left(y_5b\sin\gamma + z_5c_y\right)\hat{\mathbf{y}} + z_5c_z\,\hat{\mathbf{z}}$	$(2i)$	O V
$\mathbf{B}_{10}$	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	$(-x_5a - y_5b\cos\gamma - z_5c_x)\,\hat{\mathbf{x}} + \left(-y_5b\sin\gamma - z_5c_y\right)\hat{\mathbf{y}} - z_5c_z\,\hat{\mathbf{z}}$	$(2i)$	O V
$\mathbf{B}_{11}$	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	$(x_6a + y_6b\cos\gamma + z_6c_x)\,\hat{\mathbf{x}} + \left(y_6b\sin\gamma + z_6c_y\right)\hat{\mathbf{y}} + z_6c_z\,\hat{\mathbf{z}}$	$(2i)$	O VI
$\mathbf{B}_{12}$	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	$(-x_6a - y_6b\cos\gamma - z_6c_x)\,\hat{\mathbf{x}} + \left(-y_6b\sin\gamma - z_6c_y\right)\hat{\mathbf{y}} - z_6c_z\,\hat{\mathbf{z}}$	$(2i)$	O VI
$\mathbf{B}_{13}$	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	$(x_7a + y_7b\cos\gamma + z_7c_x)\,\hat{\mathbf{x}} + \left(y_7b\sin\gamma + z_7c_y\right)\hat{\mathbf{y}} + z_7c_z\,\hat{\mathbf{z}}$	$(2i)$	O VII
$\mathbf{B}_{14}$	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	$(-x_7a - y_7b\cos\gamma - z_7c_x)\,\hat{\mathbf{x}} + \left(-y_7b\sin\gamma - z_7c_y\right)\hat{\mathbf{y}} - z_7c_z\,\hat{\mathbf{z}}$	$(2i)$	O VII
$\mathbf{B}_{15}$	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	$(x_8a + y_8b\cos\gamma + z_8c_x)\,\hat{\mathbf{x}} + \left(y_8b\sin\gamma + z_8c_y\right)\hat{\mathbf{y}} + z_8c_z\,\hat{\mathbf{z}}$	$(2i)$	O VIII
$\mathbf{B}_{16}$	$-x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	$(-x_8a - y_8b\cos\gamma - z_8c_x)\,\hat{\mathbf{x}} + \left(-y_8b\sin\gamma - z_8c_y\right)\hat{\mathbf{y}} - z_8c_z\,\hat{\mathbf{z}}$	$(2i)$	O VIII

$$\mathbf{B}_{17} = x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 = (x_9a + y_9b\cos\gamma + z_9c_x)\,\hat{\mathbf{x}} + \left(y_9b\sin\gamma + z_9c_y\right)\hat{\mathbf{y}} + z_9c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{O IX}$$

$$\mathbf{B}_{18} = -x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3 = (-x_9a - y_9b\cos\gamma - z_9c_x)\,\hat{\mathbf{x}} + \left(-y_9b\sin\gamma - z_9c_y\right)\hat{\mathbf{y}} - z_9c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{O IX}$$

$$\mathbf{B}_{19} = x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 = (x_{10}a + y_{10}b\cos\gamma + z_{10}c_x)\,\hat{\mathbf{x}} + \left(y_{10}b\sin\gamma + z_{10}c_y\right)\hat{\mathbf{y}} + z_{10}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{O X}$$

$$\mathbf{B}_{20} = -x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 = (-x_{10}a - y_{10}b\cos\gamma - z_{10}c_x)\,\hat{\mathbf{x}} + \left(-y_{10}b\sin\gamma - z_{10}c_y\right)\hat{\mathbf{y}} - z_{10}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{O X}$$

$$\mathbf{B}_{21} = x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 = (x_{11}a + y_{11}b\cos\gamma + z_{11}c_x)\,\hat{\mathbf{x}} + \left(y_{11}b\sin\gamma + z_{11}c_y\right)\hat{\mathbf{y}} + z_{11}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{O XI}$$

$$\mathbf{B}_{22} = -x_{11}\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3 = (-x_{11}a - y_{11}b\cos\gamma - z_{11}c_x)\,\hat{\mathbf{x}} + \left(-y_{11}b\sin\gamma - z_{11}c_y\right)\hat{\mathbf{y}} - z_{11}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{O XI}$$

$$\mathbf{B}_{23} = x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 = (x_{12}a + y_{12}b\cos\gamma + z_{12}c_x)\,\hat{\mathbf{x}} + \left(y_{12}b\sin\gamma + z_{12}c_y\right)\hat{\mathbf{y}} + z_{12}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{O XII}$$

$$\mathbf{B}_{24} = -x_{12}\,\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3 = (-x_{12}a - y_{12}b\cos\gamma - z_{12}c_x)\,\hat{\mathbf{x}} + \left(-y_{12}b\sin\gamma - z_{12}c_y\right)\hat{\mathbf{y}} - z_{12}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{O XII}$$

$$\mathbf{B}_{25} = x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3 = (x_{13}a + y_{13}b\cos\gamma + z_{13}c_x)\,\hat{\mathbf{x}} + \left(y_{13}b\sin\gamma + z_{13}c_y\right)\hat{\mathbf{y}} + z_{13}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{W I}$$

$$\mathbf{B}_{26} = -x_{13}\,\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 - z_{13}\,\mathbf{a}_3 = (-x_{13}a - y_{13}b\cos\gamma - z_{13}c_x)\,\hat{\mathbf{x}} + \left(-y_{13}b\sin\gamma - z_{13}c_y\right)\hat{\mathbf{y}} - z_{13}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{W I}$$

$$\mathbf{B}_{27} = x_{14}\,\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3 = (x_{14}a + y_{14}b\cos\gamma + z_{14}c_x)\,\hat{\mathbf{x}} + \left(y_{14}b\sin\gamma + z_{14}c_y\right)\hat{\mathbf{y}} + z_{14}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{W II}$$

$$\mathbf{B}_{28} = -x_{14}\,\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 - z_{14}\,\mathbf{a}_3 = (-x_{14}a - y_{14}b\cos\gamma - z_{14}c_x)\,\hat{\mathbf{x}} + \left(-y_{14}b\sin\gamma - z_{14}c_y\right)\hat{\mathbf{y}} - z_{14}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{W II}$$

$$\mathbf{B}_{29} = x_{15}\,\mathbf{a}_1 + y_{15}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 = (x_{15}a + y_{15}b\cos\gamma + z_{15}c_x)\,\hat{\mathbf{x}} + \left(y_{15}b\sin\gamma + z_{15}c_y\right)\hat{\mathbf{y}} + z_{15}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{W III}$$

$$\mathbf{B}_{30} = -x_{15}\,\mathbf{a}_1 - y_{15}\,\mathbf{a}_2 - z_{15}\,\mathbf{a}_3 = (-x_{15}a - y_{15}b\cos\gamma - z_{15}c_x)\,\hat{\mathbf{x}} + \left(-y_{15}b\sin\gamma - z_{15}c_y\right)\hat{\mathbf{y}} - z_{15}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{W III}$$

$$\mathbf{B}_{31} = x_{16}\,\mathbf{a}_1 + y_{16}\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 = (x_{16}a + y_{16}b\cos\gamma + z_{16}c_x)\,\hat{\mathbf{x}} + \left(y_{16}b\sin\gamma + z_{16}c_y\right)\hat{\mathbf{y}} + z_{16}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{W IV}$$

$$\mathbf{B}_{32} = -x_{16}\,\mathbf{a}_1 - y_{16}\,\mathbf{a}_2 - z_{16}\,\mathbf{a}_3 = (-x_{16}a - y_{16}b\cos\gamma - z_{16}c_x)\,\hat{\mathbf{x}} + \left(-y_{16}b\sin\gamma - z_{16}c_y\right)\hat{\mathbf{y}} - z_{16}c_z\,\hat{\mathbf{z}} \quad (2i) \quad \text{W IV}$$

References:

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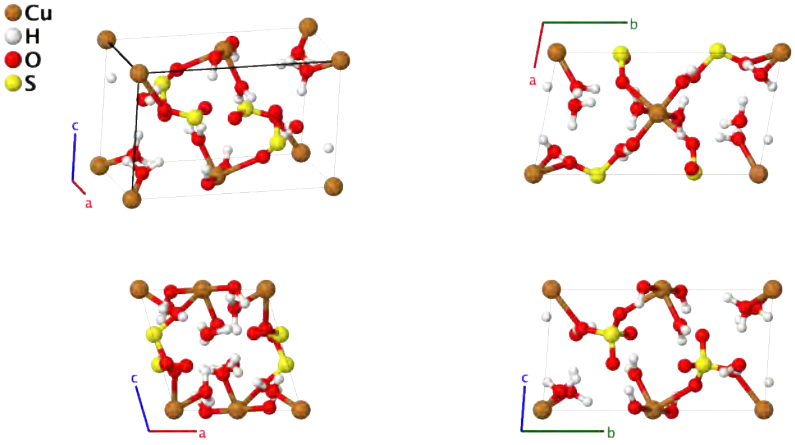
- B. Gerand, G. Nowogrocki, J. Guenot, and M. Figlarz, *Structural study of a new hexagonal form of tungsten trioxide*, J. Solid State Chem. **29**, 429–434 (1979), doi:10.1016/0022-4596(79)90199-3.

Geometry files:
- CIF: pp. 1503
- POSCAR: pp. 1503

Chalcanthite ($CuSO_4{\cdot}5H_2O$, $H4_{10}$) Structure: AB10C9D_aP42_2_ae_10i_9i_i

http://aflow.org/prototype-encyclopedia/AB10C9D_aP42_2_ae_10i_9i_i



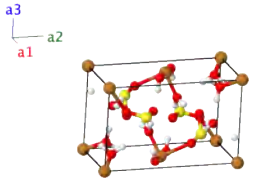
Prototype	:	$CuH_{10}O_9S$
AFLOW prototype label	:	AB10C9D_aP42_2_ae_10i_9i_i
***Strukturbericht* designation**	:	$H4_{10}$
Pearson symbol	:	aP42
Space group number	:	2
Space group symbol	:	$P\bar{1}$
AFLOW prototype command	:	`aflow --proto=AB10C9D_aP42_2_ae_10i_9i_i --params=`$a, b/a, c/a, \alpha, \beta, \gamma, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}, x_{17}, y_{17}, z_{17}, x_{18}, y_{18}, z_{18}, x_{19}, y_{19}, z_{19}, x_{20}, y_{20}, z_{20}, x_{21}, y_{21}, z_{21}, x_{22}, y_{22}, z_{22}$

- We reference (Bacon, 1975) in order to include the positions of the hydrogen atoms.

Triclinic primitive vectors:

$$\mathbf{a}_1 = a\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\cos\gamma\,\hat{\mathbf{x}} + b\sin\gamma\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c_x\hat{\mathbf{x}} + c_y\hat{\mathbf{y}} + c_z\hat{\mathbf{z}}$$

$$c_x = c\cos\beta$$
$$c_y = c\left(\cos\alpha - \cos\beta\cos\gamma\right)/\sin\gamma$$
$$c_z = \sqrt{c^2 - c_x^2 - c_y^2}$$



Basis vectors:

	Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$ =	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(1a)$	Cu I

$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}\,(a + b\cos\gamma)\;\hat{\mathbf{x}} + \frac{1}{2} b \sin\gamma\,\hat{\mathbf{y}}$	(1*e*)	Cu II
$\mathbf{B}_3$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + y_3 b\cos\gamma + z_3 c_x)\;\hat{\mathbf{x}} + \left(y_3 b\sin\gamma + z_3 c_y\right)\hat{\mathbf{y}} + z_3 c_z\,\hat{\mathbf{z}}$	(2*i*)	H I
$\mathbf{B}_4$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - y_3 b\cos\gamma - z_3 c_x)\;\hat{\mathbf{x}} + \left(-y_3 b\sin\gamma - z_3 c_y\right)\hat{\mathbf{y}} - z_3 c_z\,\hat{\mathbf{z}}$	(2*i*)	H I
$\mathbf{B}_5$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + y_4 b\cos\gamma + z_4 c_x)\;\hat{\mathbf{x}} + \left(y_4 b\sin\gamma + z_4 c_y\right)\hat{\mathbf{y}} + z_4 c_z\,\hat{\mathbf{z}}$	(2*i*)	H II
$\mathbf{B}_6$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - y_4 b\cos\gamma - z_4 c_x)\;\hat{\mathbf{x}} + \left(-y_4 b\sin\gamma - z_4 c_y\right)\hat{\mathbf{y}} - z_4 c_z\,\hat{\mathbf{z}}$	(2*i*)	H II
$\mathbf{B}_7$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + y_5 b\cos\gamma + z_5 c_x)\;\hat{\mathbf{x}} + \left(y_5 b\sin\gamma + z_5 c_y\right)\hat{\mathbf{y}} + z_5 c_z\,\hat{\mathbf{z}}$	(2*i*)	H III
$\mathbf{B}_8$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - y_5 b\cos\gamma - z_5 c_x)\;\hat{\mathbf{x}} + \left(-y_5 b\sin\gamma - z_5 c_y\right)\hat{\mathbf{y}} - z_5 c_z\,\hat{\mathbf{z}}$	(2*i*)	H III
$\mathbf{B}_9$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + y_6 b\cos\gamma + z_6 c_x)\;\hat{\mathbf{x}} + \left(y_6 b\sin\gamma + z_6 c_y\right)\hat{\mathbf{y}} + z_6 c_z\,\hat{\mathbf{z}}$	(2*i*)	H IV
$\mathbf{B}_{10}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6 a - y_6 b\cos\gamma - z_6 c_x)\;\hat{\mathbf{x}} + \left(-y_6 b\sin\gamma - z_6 c_y\right)\hat{\mathbf{y}} - z_6 c_z\,\hat{\mathbf{z}}$	(2*i*)	H IV
$\mathbf{B}_{11}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + y_7 b\cos\gamma + z_7 c_x)\;\hat{\mathbf{x}} + \left(y_7 b\sin\gamma + z_7 c_y\right)\hat{\mathbf{y}} + z_7 c_z\,\hat{\mathbf{z}}$	(2*i*)	H V
$\mathbf{B}_{12}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7 a - y_7 b\cos\gamma - z_7 c_x)\;\hat{\mathbf{x}} + \left(-y_7 b\sin\gamma - z_7 c_y\right)\hat{\mathbf{y}} - z_7 c_z\,\hat{\mathbf{z}}$	(2*i*)	H V
$\mathbf{B}_{13}$	=	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8 a + y_8 b\cos\gamma + z_8 c_x)\;\hat{\mathbf{x}} + \left(y_8 b\sin\gamma + z_8 c_y\right)\hat{\mathbf{y}} + z_8 c_z\,\hat{\mathbf{z}}$	(2*i*)	H VI
$\mathbf{B}_{14}$	=	$-x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8 a - y_8 b\cos\gamma - z_8 c_x)\;\hat{\mathbf{x}} + \left(-y_8 b\sin\gamma - z_8 c_y\right)\hat{\mathbf{y}} - z_8 c_z\,\hat{\mathbf{z}}$	(2*i*)	H VI
$\mathbf{B}_{15}$	=	$x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$(x_9 a + y_9 b\cos\gamma + z_9 c_x)\;\hat{\mathbf{x}} + \left(y_9 b\sin\gamma + z_9 c_y\right)\hat{\mathbf{y}} + z_9 c_z\,\hat{\mathbf{z}}$	(2*i*)	H VII
$\mathbf{B}_{16}$	=	$-x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$(-x_9 a - y_9 b\cos\gamma - z_9 c_x)\;\hat{\mathbf{x}} + \left(-y_9 b\sin\gamma - z_9 c_y\right)\hat{\mathbf{y}} - z_9 c_z\,\hat{\mathbf{z}}$	(2*i*)	H VII
$\mathbf{B}_{17}$	=	$x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$(x_{10} a + y_{10} b\cos\gamma + z_{10} c_x)\;\hat{\mathbf{x}} + \left(y_{10} b\sin\gamma + z_{10} c_y\right)\hat{\mathbf{y}} + z_{10} c_z\,\hat{\mathbf{z}}$	(2*i*)	H VIII
$\mathbf{B}_{18}$	=	$-x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3$	=	$(-x_{10} a - y_{10} b\cos\gamma - z_{10} c_x)\;\hat{\mathbf{x}} + \left(-y_{10} b\sin\gamma - z_{10} c_y\right)\hat{\mathbf{y}} - z_{10} c_z\,\hat{\mathbf{z}}$	(2*i*)	H VIII
$\mathbf{B}_{19}$	=	$x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$(x_{11} a + y_{11} b\cos\gamma + z_{11} c_x)\;\hat{\mathbf{x}} + \left(y_{11} b\sin\gamma + z_{11} c_y\right)\hat{\mathbf{y}} + z_{11} c_z\,\hat{\mathbf{z}}$	(2*i*)	H IX
$\mathbf{B}_{20}$	=	$-x_{11}\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3$	=	$(-x_{11} a - y_{11} b\cos\gamma - z_{11} c_x)\;\hat{\mathbf{x}} + \left(-y_{11} b\sin\gamma - z_{11} c_y\right)\hat{\mathbf{y}} - z_{11} c_z\,\hat{\mathbf{z}}$	(2*i*)	H IX
$\mathbf{B}_{21}$	=	$x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3$	=	$(x_{12} a + y_{12} b\cos\gamma + z_{12} c_x)\;\hat{\mathbf{x}} + \left(y_{12} b\sin\gamma + z_{12} c_y\right)\hat{\mathbf{y}} + z_{12} c_z\,\hat{\mathbf{z}}$	(2*i*)	H X
$\mathbf{B}_{22}$	=	$-x_{12}\,\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3$	=	$(-x_{12} a - y_{12} b\cos\gamma - z_{12} c_x)\;\hat{\mathbf{x}} + \left(-y_{12} b\sin\gamma - z_{12} c_y\right)\hat{\mathbf{y}} - z_{12} c_z\,\hat{\mathbf{z}}$	(2*i*)	H X
$\mathbf{B}_{23}$	=	$x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3$	=	$(x_{13} a + y_{13} b\cos\gamma + z_{13} c_x)\;\hat{\mathbf{x}} + \left(y_{13} b\sin\gamma + z_{13} c_y\right)\hat{\mathbf{y}} + z_{13} c_z\,\hat{\mathbf{z}}$	(2*i*)	O I
$\mathbf{B}_{24}$	=	$-x_{13}\,\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 - z_{13}\,\mathbf{a}_3$	=	$(-x_{13} a - y_{13} b\cos\gamma - z_{13} c_x)\;\hat{\mathbf{x}} + \left(-y_{13} b\sin\gamma - z_{13} c_y\right)\hat{\mathbf{y}} - z_{13} c_z\,\hat{\mathbf{z}}$	(2*i*)	O I

$\mathbf{B}_{25}$	=	$x_{14}\,\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3$	=	$(x_{14}a + y_{14}b\cos\gamma + z_{14}c_x)\,\hat{\mathbf{x}} + \left(y_{14}b\sin\gamma + z_{14}c_y\right)\hat{\mathbf{y}} + z_{14}c_z\,\hat{\mathbf{z}}$	(2*i*)	O II
$\mathbf{B}_{26}$	=	$-x_{14}\,\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 - z_{14}\,\mathbf{a}_3$	=	$(-x_{14}a - y_{14}b\cos\gamma - z_{14}c_x)\,\hat{\mathbf{x}} + \left(-y_{14}b\sin\gamma - z_{14}c_y\right)\hat{\mathbf{y}} - z_{14}c_z\,\hat{\mathbf{z}}$	(2*i*)	O II
$\mathbf{B}_{27}$	=	$x_{15}\,\mathbf{a}_1 + y_{15}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3$	=	$(x_{15}a + y_{15}b\cos\gamma + z_{15}c_x)\,\hat{\mathbf{x}} + \left(y_{15}b\sin\gamma + z_{15}c_y\right)\hat{\mathbf{y}} + z_{15}c_z\,\hat{\mathbf{z}}$	(2*i*)	O III
$\mathbf{B}_{28}$	=	$-x_{15}\,\mathbf{a}_1 - y_{15}\,\mathbf{a}_2 - z_{15}\,\mathbf{a}_3$	=	$(-x_{15}a - y_{15}b\cos\gamma - z_{15}c_x)\,\hat{\mathbf{x}} + \left(-y_{15}b\sin\gamma - z_{15}c_y\right)\hat{\mathbf{y}} - z_{15}c_z\,\hat{\mathbf{z}}$	(2*i*)	O III
$\mathbf{B}_{29}$	=	$x_{16}\,\mathbf{a}_1 + y_{16}\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3$	=	$(x_{16}a + y_{16}b\cos\gamma + z_{16}c_x)\,\hat{\mathbf{x}} + \left(y_{16}b\sin\gamma + z_{16}c_y\right)\hat{\mathbf{y}} + z_{16}c_z\,\hat{\mathbf{z}}$	(2*i*)	O IV
$\mathbf{B}_{30}$	=	$-x_{16}\,\mathbf{a}_1 - y_{16}\,\mathbf{a}_2 - z_{16}\,\mathbf{a}_3$	=	$(-x_{16}a - y_{16}b\cos\gamma - z_{16}c_x)\,\hat{\mathbf{x}} + \left(-y_{16}b\sin\gamma - z_{16}c_y\right)\hat{\mathbf{y}} - z_{16}c_z\,\hat{\mathbf{z}}$	(2*i*)	O IV
$\mathbf{B}_{31}$	=	$x_{17}\,\mathbf{a}_1 + y_{17}\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3$	=	$(x_{17}a + y_{17}b\cos\gamma + z_{17}c_x)\,\hat{\mathbf{x}} + \left(y_{17}b\sin\gamma + z_{17}c_y\right)\hat{\mathbf{y}} + z_{17}c_z\,\hat{\mathbf{z}}$	(2*i*)	O V
$\mathbf{B}_{32}$	=	$-x_{17}\,\mathbf{a}_1 - y_{17}\,\mathbf{a}_2 - z_{17}\,\mathbf{a}_3$	=	$(-x_{17}a - y_{17}b\cos\gamma - z_{17}c_x)\,\hat{\mathbf{x}} + \left(-y_{17}b\sin\gamma - z_{17}c_y\right)\hat{\mathbf{y}} - z_{17}c_z\,\hat{\mathbf{z}}$	(2*i*)	O V
$\mathbf{B}_{33}$	=	$x_{18}\,\mathbf{a}_1 + y_{18}\,\mathbf{a}_2 + z_{18}\,\mathbf{a}_3$	=	$(x_{18}a + y_{18}b\cos\gamma + z_{18}c_x)\,\hat{\mathbf{x}} + \left(y_{18}b\sin\gamma + z_{18}c_y\right)\hat{\mathbf{y}} + z_{18}c_z\,\hat{\mathbf{z}}$	(2*i*)	O VI
$\mathbf{B}_{34}$	=	$-x_{18}\,\mathbf{a}_1 - y_{18}\,\mathbf{a}_2 - z_{18}\,\mathbf{a}_3$	=	$(-x_{18}a - y_{18}b\cos\gamma - z_{18}c_x)\,\hat{\mathbf{x}} + \left(-y_{18}b\sin\gamma - z_{18}c_y\right)\hat{\mathbf{y}} - z_{18}c_z\,\hat{\mathbf{z}}$	(2*i*)	O VI
$\mathbf{B}_{35}$	=	$x_{19}\,\mathbf{a}_1 + y_{19}\,\mathbf{a}_2 + z_{19}\,\mathbf{a}_3$	=	$(x_{19}a + y_{19}b\cos\gamma + z_{19}c_x)\,\hat{\mathbf{x}} + \left(y_{19}b\sin\gamma + z_{19}c_y\right)\hat{\mathbf{y}} + z_{19}c_z\,\hat{\mathbf{z}}$	(2*i*)	O VII
$\mathbf{B}_{36}$	=	$-x_{19}\,\mathbf{a}_1 - y_{19}\,\mathbf{a}_2 - z_{19}\,\mathbf{a}_3$	=	$(-x_{19}a - y_{19}b\cos\gamma - z_{19}c_x)\,\hat{\mathbf{x}} + \left(-y_{19}b\sin\gamma - z_{19}c_y\right)\hat{\mathbf{y}} - z_{19}c_z\,\hat{\mathbf{z}}$	(2*i*)	O VII
$\mathbf{B}_{37}$	=	$x_{20}\,\mathbf{a}_1 + y_{20}\,\mathbf{a}_2 + z_{20}\,\mathbf{a}_3$	=	$(x_{20}a + y_{20}b\cos\gamma + z_{20}c_x)\,\hat{\mathbf{x}} + \left(y_{20}b\sin\gamma + z_{20}c_y\right)\hat{\mathbf{y}} + z_{20}c_z\,\hat{\mathbf{z}}$	(2*i*)	O VIII
$\mathbf{B}_{38}$	=	$-x_{20}\,\mathbf{a}_1 - y_{20}\,\mathbf{a}_2 - z_{20}\,\mathbf{a}_3$	=	$(-x_{20}a - y_{20}b\cos\gamma - z_{20}c_x)\,\hat{\mathbf{x}} + \left(-y_{20}b\sin\gamma - z_{20}c_y\right)\hat{\mathbf{y}} - z_{20}c_z\,\hat{\mathbf{z}}$	(2*i*)	O VIII
$\mathbf{B}_{39}$	=	$x_{21}\,\mathbf{a}_1 + y_{21}\,\mathbf{a}_2 + z_{21}\,\mathbf{a}_3$	=	$(x_{21}a + y_{21}b\cos\gamma + z_{21}c_x)\,\hat{\mathbf{x}} + \left(y_{21}b\sin\gamma + z_{21}c_y\right)\hat{\mathbf{y}} + z_{21}c_z\,\hat{\mathbf{z}}$	(2*i*)	O IX
$\mathbf{B}_{40}$	=	$-x_{21}\,\mathbf{a}_1 - y_{21}\,\mathbf{a}_2 - z_{21}\,\mathbf{a}_3$	=	$(-x_{21}a - y_{21}b\cos\gamma - z_{21}c_x)\,\hat{\mathbf{x}} + \left(-y_{21}b\sin\gamma - z_{21}c_y\right)\hat{\mathbf{y}} - z_{21}c_z\,\hat{\mathbf{z}}$	(2*i*)	O IX
$\mathbf{B}_{41}$	=	$x_{22}\,\mathbf{a}_1 + y_{22}\,\mathbf{a}_2 + z_{22}\,\mathbf{a}_3$	=	$(x_{22}a + y_{22}b\cos\gamma + z_{22}c_x)\,\hat{\mathbf{x}} + \left(y_{22}b\sin\gamma + z_{22}c_y\right)\hat{\mathbf{y}} + z_{22}c_z\,\hat{\mathbf{z}}$	(2*i*)	S
$\mathbf{B}_{42}$	=	$-x_{22}\,\mathbf{a}_1 - y_{22}\,\mathbf{a}_2 - z_{22}\,\mathbf{a}_3$	=	$(-x_{22}a - y_{22}b\cos\gamma - z_{22}c_x)\,\hat{\mathbf{x}} + \left(-y_{22}b\sin\gamma - z_{22}c_y\right)\hat{\mathbf{y}} - z_{22}c_z\,\hat{\mathbf{z}}$	(2*i*)	S

References:

- G. E. Bacon and D. H. Titterton, *Neutron-diffraction studies of $CuSO_4{\cdot}5H_2O$ and $CuSO_4{\cdot}5D_2O$*, Zeitschrift für Kristallographie - Crystalline Materials **141**, 330–341 (1975), doi:10.1524/zkri.1975.141.16.330.

Found in:

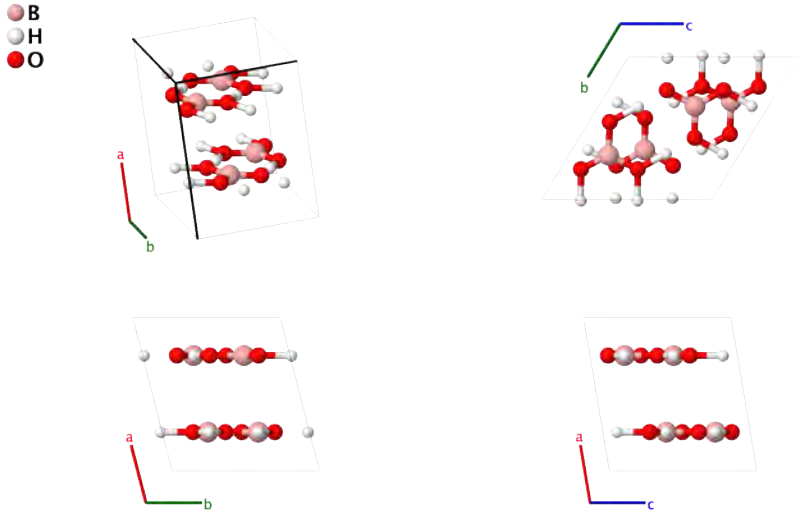
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1504
- POSCAR: pp. 1504

Boric Acid (H_3BO_3, $G5_1$) Structure: AB3C3_aP28_2_2i_6i_6i

http://aflow.org/prototype-encyclopedia/AB3C3_aP28_2_2i_6i_6i

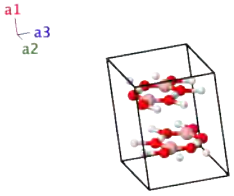


Prototype	:	BH_3O_3
AFLOW prototype label	:	AB3C3_aP28_2_2i_6i_6i
***Strukturbericht* designation**	:	$G5_1$
Pearson symbol	:	aP28
Space group number	:	2
Space group symbol	:	$P\bar{1}$
AFLOW prototype command	:	`aflow --proto=AB3C3_aP28_2_2i_6i_6i --params=`$a, b/a, c/a, \alpha, \beta, \gamma, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}$

Triclinic primitive vectors:

$$\mathbf{a}_1 = a\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\cos\gamma\,\hat{\mathbf{x}} + b\sin\gamma\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c_x\hat{\mathbf{x}} + c_y\,\hat{\mathbf{y}} + c_z\,\hat{\mathbf{z}}$$

$$c_x = c\cos\beta$$
$$c_y = c\left(\cos\alpha - \cos\beta\cos\gamma\right)/\sin\gamma$$
$$c_z = \sqrt{c^2 - c_x^2 - c_y^2}$$



Basis vectors:

	Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$= x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	$=$	$(x_1a + y_1b\cos\gamma + z_1c_x)\,\hat{\mathbf{x}} + \left(y_1b\sin\gamma + z_1c_y\right)\,\hat{\mathbf{y}} + z_1c_z\,\hat{\mathbf{z}}$	$(2i)$	B I

$$
\begin{array}{lcccccc}
\mathbf{B}_2 & = & -x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3 & = & \begin{array}{l}(-x_1a - y_1b\cos\gamma - z_1c_x)\;\hat{\mathbf{x}} + \\ \left(-y_1b\sin\gamma - z_1c_y\right)\hat{\mathbf{y}} - z_1c_z\,\hat{\mathbf{z}}\end{array} & (2i) & \text{B I} \\
\mathbf{B}_3 & = & x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3 & = & \begin{array}{l}(x_2a + y_2b\cos\gamma + z_2c_x)\;\hat{\mathbf{x}} + \\ \left(y_2b\sin\gamma + z_2c_y\right)\hat{\mathbf{y}} + z_2c_z\,\hat{\mathbf{z}}\end{array} & (2i) & \text{B II} \\
\mathbf{B}_4 & = & -x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3 & = & \begin{array}{l}(-x_2a - y_2b\cos\gamma - z_2c_x)\;\hat{\mathbf{x}} + \\ \left(-y_2b\sin\gamma - z_2c_y\right)\hat{\mathbf{y}} - z_2c_z\,\hat{\mathbf{z}}\end{array} & (2i) & \text{B II} \\
\mathbf{B}_5 & = & x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3 & = & \begin{array}{l}(x_3a + y_3b\cos\gamma + z_3c_x)\;\hat{\mathbf{x}} + \\ \left(y_3b\sin\gamma + z_3c_y\right)\hat{\mathbf{y}} + z_3c_z\,\hat{\mathbf{z}}\end{array} & (2i) & \text{H I} \\
\mathbf{B}_6 & = & -x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3 & = & \begin{array}{l}(-x_3a - y_3b\cos\gamma - z_3c_x)\;\hat{\mathbf{x}} + \\ \left(-y_3b\sin\gamma - z_3c_y\right)\hat{\mathbf{y}} - z_3c_z\,\hat{\mathbf{z}}\end{array} & (2i) & \text{H I} \\
\mathbf{B}_7 & = & x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3 & = & \begin{array}{l}(x_4a + y_4b\cos\gamma + z_4c_x)\;\hat{\mathbf{x}} + \\ \left(y_4b\sin\gamma + z_4c_y\right)\hat{\mathbf{y}} + z_4c_z\,\hat{\mathbf{z}}\end{array} & (2i) & \text{H II} \\
\mathbf{B}_8 & = & -x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3 & = & \begin{array}{l}(-x_4a - y_4b\cos\gamma - z_4c_x)\;\hat{\mathbf{x}} + \\ \left(-y_4b\sin\gamma - z_4c_y\right)\hat{\mathbf{y}} - z_4c_z\,\hat{\mathbf{z}}\end{array} & (2i) & \text{H II} \\
\mathbf{B}_9 & = & x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3 & = & \begin{array}{l}(x_5a + y_5b\cos\gamma + z_5c_x)\;\hat{\mathbf{x}} + \\ \left(y_5b\sin\gamma + z_5c_y\right)\hat{\mathbf{y}} + z_5c_z\,\hat{\mathbf{z}}\end{array} & (2i) & \text{H III} \\
\mathbf{B}_{10} & = & -x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3 & = & \begin{array}{l}(-x_5a - y_5b\cos\gamma - z_5c_x)\;\hat{\mathbf{x}} + \\ \left(-y_5b\sin\gamma - z_5c_y\right)\hat{\mathbf{y}} - z_5c_z\,\hat{\mathbf{z}}\end{array} & (2i) & \text{H III} \\
\mathbf{B}_{11} & = & x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 & = & \begin{array}{l}(x_6a + y_6b\cos\gamma + z_6c_x)\;\hat{\mathbf{x}} + \\ \left(y_6b\sin\gamma + z_6c_y\right)\hat{\mathbf{y}} + z_6c_z\,\hat{\mathbf{z}}\end{array} & (2i) & \text{H IV} \\
\mathbf{B}_{12} & = & -x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3 & = & \begin{array}{l}(-x_6a - y_6b\cos\gamma - z_6c_x)\;\hat{\mathbf{x}} + \\ \left(-y_6b\sin\gamma - z_6c_y\right)\hat{\mathbf{y}} - z_6c_z\,\hat{\mathbf{z}}\end{array} & (2i) & \text{H IV} \\
\mathbf{B}_{13} & = & x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 & = & \begin{array}{l}(x_7a + y_7b\cos\gamma + z_7c_x)\;\hat{\mathbf{x}} + \\ \left(y_7b\sin\gamma + z_7c_y\right)\hat{\mathbf{y}} + z_7c_z\,\hat{\mathbf{z}}\end{array} & (2i) & \text{H V} \\
\mathbf{B}_{14} & = & -x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3 & = & \begin{array}{l}(-x_7a - y_7b\cos\gamma - z_7c_x)\;\hat{\mathbf{x}} + \\ \left(-y_7b\sin\gamma - z_7c_y\right)\hat{\mathbf{y}} - z_7c_z\,\hat{\mathbf{z}}\end{array} & (2i) & \text{H V} \\
\mathbf{B}_{15} & = & x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 & = & \begin{array}{l}(x_8a + y_8b\cos\gamma + z_8c_x)\;\hat{\mathbf{x}} + \\ \left(y_8b\sin\gamma + z_8c_y\right)\hat{\mathbf{y}} + z_8c_z\,\hat{\mathbf{z}}\end{array} & (2i) & \text{H VI} \\
\mathbf{B}_{16} & = & -x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3 & = & \begin{array}{l}(-x_8a - y_8b\cos\gamma - z_8c_x)\;\hat{\mathbf{x}} + \\ \left(-y_8b\sin\gamma - z_8c_y\right)\hat{\mathbf{y}} - z_8c_z\,\hat{\mathbf{z}}\end{array} & (2i) & \text{H VI} \\
\mathbf{B}_{17} & = & x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 & = & \begin{array}{l}(x_9a + y_9b\cos\gamma + z_9c_x)\;\hat{\mathbf{x}} + \\ \left(y_9b\sin\gamma + z_9c_y\right)\hat{\mathbf{y}} + z_9c_z\,\hat{\mathbf{z}}\end{array} & (2i) & \text{O I} \\
\mathbf{B}_{18} & = & -x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3 & = & \begin{array}{l}(-x_9a - y_9b\cos\gamma - z_9c_x)\;\hat{\mathbf{x}} + \\ \left(-y_9b\sin\gamma - z_9c_y\right)\hat{\mathbf{y}} - z_9c_z\,\hat{\mathbf{z}}\end{array} & (2i) & \text{O I} \\
\mathbf{B}_{19} & = & x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 & = & \begin{array}{l}(x_{10}a + y_{10}b\cos\gamma + z_{10}c_x)\;\hat{\mathbf{x}} + \\ \left(y_{10}b\sin\gamma + z_{10}c_y\right)\hat{\mathbf{y}} + z_{10}c_z\,\hat{\mathbf{z}}\end{array} & (2i) & \text{O II} \\
\mathbf{B}_{20} & = & -x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 & = & \begin{array}{l}(-x_{10}a - y_{10}b\cos\gamma - z_{10}c_x)\;\hat{\mathbf{x}} + \\ \left(-y_{10}b\sin\gamma - z_{10}c_y\right)\hat{\mathbf{y}} - z_{10}c_z\,\hat{\mathbf{z}}\end{array} & (2i) & \text{O II} \\
\mathbf{B}_{21} & = & x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 & = & \begin{array}{l}(x_{11}a + y_{11}b\cos\gamma + z_{11}c_x)\;\hat{\mathbf{x}} + \\ \left(y_{11}b\sin\gamma + z_{11}c_y\right)\hat{\mathbf{y}} + z_{11}c_z\,\hat{\mathbf{z}}\end{array} & (2i) & \text{O III} \\
\mathbf{B}_{22} & = & -x_{11}\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3 & = & \begin{array}{l}(-x_{11}a - y_{11}b\cos\gamma - z_{11}c_x)\;\hat{\mathbf{x}} + \\ \left(-y_{11}b\sin\gamma - z_{11}c_y\right)\hat{\mathbf{y}} - z_{11}c_z\,\hat{\mathbf{z}}\end{array} & (2i) & \text{O III} \\
\mathbf{B}_{23} & = & x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 & = & \begin{array}{l}(x_{12}a + y_{12}b\cos\gamma + z_{12}c_x)\;\hat{\mathbf{x}} + \\ \left(y_{12}b\sin\gamma + z_{12}c_y\right)\hat{\mathbf{y}} + z_{12}c_z\,\hat{\mathbf{z}}\end{array} & (2i) & \text{O IV}
\end{array}
$$

$\mathbf{B}_{24}$	=	$-x_{12}\,\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3$	=	$(-x_{12}a - y_{12}b\cos\gamma - z_{12}c_x)\,\hat{\mathbf{x}} + \left(-y_{12}b\sin\gamma - z_{12}c_y\right)\,\hat{\mathbf{y}} - z_{12}c_z\,\hat{\mathbf{z}}$	(2*i*)	O IV
$\mathbf{B}_{25}$	=	$x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3$	=	$(x_{13}a + y_{13}b\cos\gamma + z_{13}c_x)\,\hat{\mathbf{x}} + \left(y_{13}b\sin\gamma + z_{13}c_y\right)\,\hat{\mathbf{y}} + z_{13}c_z\,\hat{\mathbf{z}}$	(2*i*)	O V
$\mathbf{B}_{26}$	=	$-x_{13}\,\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 - z_{13}\,\mathbf{a}_3$	=	$(-x_{13}a - y_{13}b\cos\gamma - z_{13}c_x)\,\hat{\mathbf{x}} + \left(-y_{13}b\sin\gamma - z_{13}c_y\right)\,\hat{\mathbf{y}} - z_{13}c_z\,\hat{\mathbf{z}}$	(2*i*)	O V
$\mathbf{B}_{27}$	=	$x_{14}\,\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3$	=	$(x_{14}a + y_{14}b\cos\gamma + z_{14}c_x)\,\hat{\mathbf{x}} + \left(y_{14}b\sin\gamma + z_{14}c_y\right)\,\hat{\mathbf{y}} + z_{14}c_z\,\hat{\mathbf{z}}$	(2*i*)	O VI
$\mathbf{B}_{28}$	=	$-x_{14}\,\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 - z_{14}\,\mathbf{a}_3$	=	$(-x_{14}a - y_{14}b\cos\gamma - z_{14}c_x)\,\hat{\mathbf{x}} + \left(-y_{14}b\sin\gamma - z_{14}c_y\right)\,\hat{\mathbf{y}} - z_{14}c_z\,\hat{\mathbf{z}}$	(2*i*)	O VI

References:

- W. H. Zachariasen, *The Crystal Lattice of Boric Acid, BO_3H_3*, Zeitschrift für Kristallographie - Crystalline Materials **88**, 150–161 (1934), doi:10.1524/zkri.1934.88.1.150.

Found in:

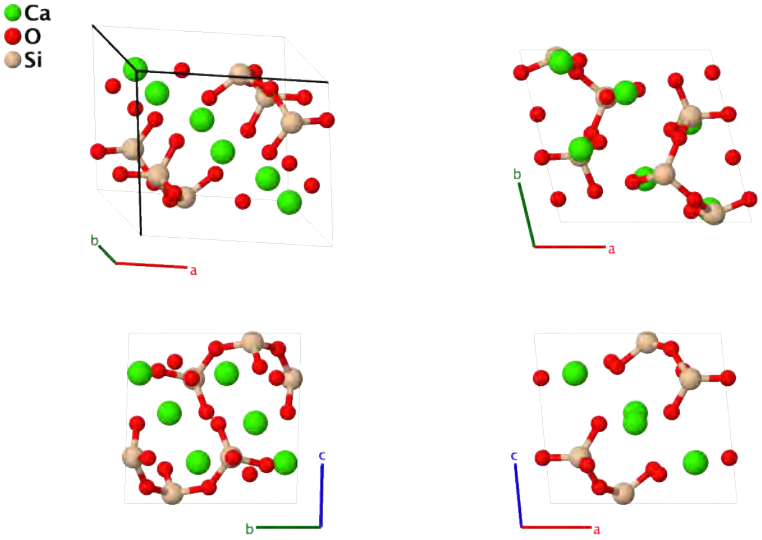
- C. Gottfried and F. Schossberger, eds., *Strukturbericht Band III 1933-1935* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1504
- POSCAR: pp. 1505

Wollastonite ($CaSiO_3$) Structure: AB3C_aP30_2_3i_9i_3i

http://aflow.org/prototype-encyclopedia/AB3C_aP30_2_3i_9i_3i



Prototype	:	CaO_3Si
AFLOW prototype label	:	AB3C_aP30_2_3i_9i_3i
***Strukturbericht* designation**	:	None
Pearson symbol	:	aP30
Space group number	:	2
Space group symbol	:	$P\bar{1}$
AFLOW prototype command	:	`aflow --proto=AB3C_aP30_2_3i_9i_3i --params=`$a, b/a, c/a, \alpha, \beta, \gamma, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}$

Other compounds with this structure

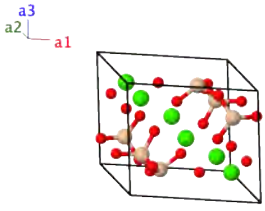
- $Ca_2NaHSi_2O_9$ (pectolite)

- (Barnick, 1936) found a monoclinic structure for wollastonite, but that structure is now referred to as parawollastonite, with the current triclinic structure receiving the original name.

Triclinic primitive vectors:

$$\mathbf{a}_1 = a\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\cos\gamma\,\hat{\mathbf{x}} + b\sin\gamma\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c_x\hat{\mathbf{x}} + c_y\,\hat{\mathbf{y}} + c_z\,\hat{\mathbf{z}}$$

$$c_x = c\cos\beta$$
$$c_y = c\left(\cos\alpha - \cos\beta\cos\gamma\right)/\sin\gamma$$
$$c_z = \sqrt{c^2 - c_x^2 - c_y^2}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1a + y_1b\cos\gamma + z_1c_x)\,\hat{\mathbf{x}} + \left(y_1b\sin\gamma + z_1c_y\right)\hat{\mathbf{y}} + z_1c_z\,\hat{\mathbf{z}}$	(2*i*)	Ca I
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$(-x_1a - y_1b\cos\gamma - z_1c_x)\,\hat{\mathbf{x}} + \left(-y_1b\sin\gamma - z_1c_y\right)\hat{\mathbf{y}} - z_1c_z\,\hat{\mathbf{z}}$	(2*i*)	Ca I
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2a + y_2b\cos\gamma + z_2c_x)\,\hat{\mathbf{x}} + \left(y_2b\sin\gamma + z_2c_y\right)\hat{\mathbf{y}} + z_2c_z\,\hat{\mathbf{z}}$	(2*i*)	Ca II
$\mathbf{B}_4$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2a - y_2b\cos\gamma - z_2c_x)\,\hat{\mathbf{x}} + \left(-y_2b\sin\gamma - z_2c_y\right)\hat{\mathbf{y}} - z_2c_z\,\hat{\mathbf{z}}$	(2*i*)	Ca II
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3a + y_3b\cos\gamma + z_3c_x)\,\hat{\mathbf{x}} + \left(y_3b\sin\gamma + z_3c_y\right)\hat{\mathbf{y}} + z_3c_z\,\hat{\mathbf{z}}$	(2*i*)	Ca III
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3a - y_3b\cos\gamma - z_3c_x)\,\hat{\mathbf{x}} + \left(-y_3b\sin\gamma - z_3c_y\right)\hat{\mathbf{y}} - z_3c_z\,\hat{\mathbf{z}}$	(2*i*)	Ca III
$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4a + y_4b\cos\gamma + z_4c_x)\,\hat{\mathbf{x}} + \left(y_4b\sin\gamma + z_4c_y\right)\hat{\mathbf{y}} + z_4c_z\,\hat{\mathbf{z}}$	(2*i*)	O I
$\mathbf{B}_8$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4a - y_4b\cos\gamma - z_4c_x)\,\hat{\mathbf{x}} + \left(-y_4b\sin\gamma - z_4c_y\right)\hat{\mathbf{y}} - z_4c_z\,\hat{\mathbf{z}}$	(2*i*)	O I
$\mathbf{B}_9$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5a + y_5b\cos\gamma + z_5c_x)\,\hat{\mathbf{x}} + \left(y_5b\sin\gamma + z_5c_y\right)\hat{\mathbf{y}} + z_5c_z\,\hat{\mathbf{z}}$	(2*i*)	O II
$\mathbf{B}_{10}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5a - y_5b\cos\gamma - z_5c_x)\,\hat{\mathbf{x}} + \left(-y_5b\sin\gamma - z_5c_y\right)\hat{\mathbf{y}} - z_5c_z\,\hat{\mathbf{z}}$	(2*i*)	O II
$\mathbf{B}_{11}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6a + y_6b\cos\gamma + z_6c_x)\,\hat{\mathbf{x}} + \left(y_6b\sin\gamma + z_6c_y\right)\hat{\mathbf{y}} + z_6c_z\,\hat{\mathbf{z}}$	(2*i*)	O III
$\mathbf{B}_{12}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6a - y_6b\cos\gamma - z_6c_x)\,\hat{\mathbf{x}} + \left(-y_6b\sin\gamma - z_6c_y\right)\hat{\mathbf{y}} - z_6c_z\,\hat{\mathbf{z}}$	(2*i*)	O III
$\mathbf{B}_{13}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7a + y_7b\cos\gamma + z_7c_x)\,\hat{\mathbf{x}} + \left(y_7b\sin\gamma + z_7c_y\right)\hat{\mathbf{y}} + z_7c_z\,\hat{\mathbf{z}}$	(2*i*)	O IV
$\mathbf{B}_{14}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7a - y_7b\cos\gamma - z_7c_x)\,\hat{\mathbf{x}} + \left(-y_7b\sin\gamma - z_7c_y\right)\hat{\mathbf{y}} - z_7c_z\,\hat{\mathbf{z}}$	(2*i*)	O IV
$\mathbf{B}_{15}$	=	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8a + y_8b\cos\gamma + z_8c_x)\,\hat{\mathbf{x}} + \left(y_8b\sin\gamma + z_8c_y\right)\hat{\mathbf{y}} + z_8c_z\,\hat{\mathbf{z}}$	(2*i*)	O V
$\mathbf{B}_{16}$	=	$-x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8a - y_8b\cos\gamma - z_8c_x)\,\hat{\mathbf{x}} + \left(-y_8b\sin\gamma - z_8c_y\right)\hat{\mathbf{y}} - z_8c_z\,\hat{\mathbf{z}}$	(2*i*)	O V
$\mathbf{B}_{17}$	=	$x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$(x_9a + y_9b\cos\gamma + z_9c_x)\,\hat{\mathbf{x}} + \left(y_9b\sin\gamma + z_9c_y\right)\hat{\mathbf{y}} + z_9c_z\,\hat{\mathbf{z}}$	(2*i*)	O VI
$\mathbf{B}_{18}$	=	$-x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$(-x_9a - y_9b\cos\gamma - z_9c_x)\,\hat{\mathbf{x}} + \left(-y_9b\sin\gamma - z_9c_y\right)\hat{\mathbf{y}} - z_9c_z\,\hat{\mathbf{z}}$	(2*i*)	O VI
$\mathbf{B}_{19}$	=	$x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$(x_{10}a + y_{10}b\cos\gamma + z_{10}c_x)\,\hat{\mathbf{x}} + \left(y_{10}b\sin\gamma + z_{10}c_y\right)\hat{\mathbf{y}} + z_{10}c_z\,\hat{\mathbf{z}}$	(2*i*)	O VII
$\mathbf{B}_{20}$	=	$-x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3$	=	$(-x_{10}a - y_{10}b\cos\gamma - z_{10}c_x)\,\hat{\mathbf{x}} + \left(-y_{10}b\sin\gamma - z_{10}c_y\right)\hat{\mathbf{y}} - z_{10}c_z\,\hat{\mathbf{z}}$	(2*i*)	O VII
$\mathbf{B}_{21}$	=	$x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$(x_{11}a + y_{11}b\cos\gamma + z_{11}c_x)\,\hat{\mathbf{x}} + \left(y_{11}b\sin\gamma + z_{11}c_y\right)\hat{\mathbf{y}} + z_{11}c_z\,\hat{\mathbf{z}}$	(2*i*)	O VIII

$$\mathbf{B}_{22} = -x_{11}\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3 = (-x_{11}a - y_{11}b\cos\gamma - z_{11}c_x)\,\hat{\mathbf{x}} + \left(-y_{11}b\sin\gamma - z_{11}c_y\right)\hat{\mathbf{y}} - z_{11}c_z\,\hat{\mathbf{z}}$$ (2*i*) O VIII

$$\mathbf{B}_{23} = x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 = (x_{12}a + y_{12}b\cos\gamma + z_{12}c_x)\,\hat{\mathbf{x}} + \left(y_{12}b\sin\gamma + z_{12}c_y\right)\hat{\mathbf{y}} + z_{12}c_z\,\hat{\mathbf{z}}$$ (2*i*) O IX

$$\mathbf{B}_{24} = -x_{12}\,\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3 = (-x_{12}a - y_{12}b\cos\gamma - z_{12}c_x)\,\hat{\mathbf{x}} + \left(-y_{12}b\sin\gamma - z_{12}c_y\right)\hat{\mathbf{y}} - z_{12}c_z\,\hat{\mathbf{z}}$$ (2*i*) O IX

$$\mathbf{B}_{25} = x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3 = (x_{13}a + y_{13}b\cos\gamma + z_{13}c_x)\,\hat{\mathbf{x}} + \left(y_{13}b\sin\gamma + z_{13}c_y\right)\hat{\mathbf{y}} + z_{13}c_z\,\hat{\mathbf{z}}$$ (2*i*) Si I

$$\mathbf{B}_{26} = -x_{13}\,\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 - z_{13}\,\mathbf{a}_3 = (-x_{13}a - y_{13}b\cos\gamma - z_{13}c_x)\,\hat{\mathbf{x}} + \left(-y_{13}b\sin\gamma - z_{13}c_y\right)\hat{\mathbf{y}} - z_{13}c_z\,\hat{\mathbf{z}}$$ (2*i*) Si I

$$\mathbf{B}_{27} = x_{14}\,\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3 = (x_{14}a + y_{14}b\cos\gamma + z_{14}c_x)\,\hat{\mathbf{x}} + \left(y_{14}b\sin\gamma + z_{14}c_y\right)\hat{\mathbf{y}} + z_{14}c_z\,\hat{\mathbf{z}}$$ (2*i*) Si II

$$\mathbf{B}_{28} = -x_{14}\,\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 - z_{14}\,\mathbf{a}_3 = (-x_{14}a - y_{14}b\cos\gamma - z_{14}c_x)\,\hat{\mathbf{x}} + \left(-y_{14}b\sin\gamma - z_{14}c_y\right)\hat{\mathbf{y}} - z_{14}c_z\,\hat{\mathbf{z}}$$ (2*i*) Si II

$$\mathbf{B}_{29} = x_{15}\,\mathbf{a}_1 + y_{15}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 = (x_{15}a + y_{15}b\cos\gamma + z_{15}c_x)\,\hat{\mathbf{x}} + \left(y_{15}b\sin\gamma + z_{15}c_y\right)\hat{\mathbf{y}} + z_{15}c_z\,\hat{\mathbf{z}}$$ (2*i*) Si III

$$\mathbf{B}_{30} = -x_{15}\,\mathbf{a}_1 - y_{15}\,\mathbf{a}_2 - z_{15}\,\mathbf{a}_3 = (-x_{15}a - y_{15}b\cos\gamma - z_{15}c_x)\,\hat{\mathbf{x}} + \left(-y_{15}b\sin\gamma - z_{15}c_y\right)\hat{\mathbf{y}} - z_{15}c_z\,\hat{\mathbf{z}}$$ (2*i*) Si III

References:

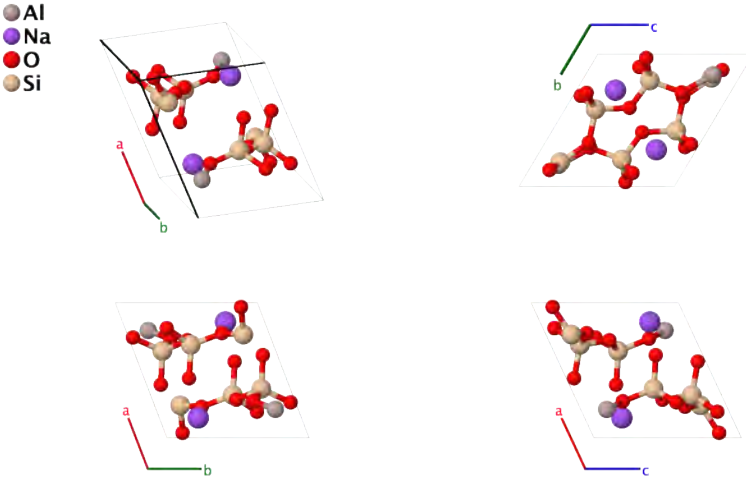
- M. J. Buerger and C. T. Prewitt, *The Crystal Structures of Wollastonite and Pectolite*, Proc. Natl. Acad. Sci. **47**, 1884–1888 (1961), doi:10.1073/pnas.47.12.1884.
- M. Barnick, *Strukturuntersuchung des natürlichen Wollastonits* (1936). Dissertation.

Geometry files:

- CIF: pp. 1505
- POSCAR: pp. 1505

Albite ($NaAlSi_3O_8$, $S6_8$) Structure: ABC8D3_aP26_2_i_i_8i_3i

http://aflow.org/prototype-encyclopedia/ABC8D3_aP26_2_i_i_8i_3i



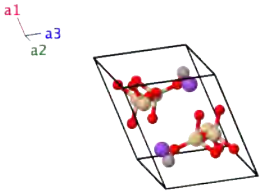
Prototype	:	$AlNaO_8Si_3$
AFLOW prototype label	:	ABC8D3_aP26_2_i_i_8i_3i
***Strukturbericht* designation**	:	$S6_8$
Pearson symbol	:	aP26
Space group number	:	2
Space group symbol	:	$P\bar{1}$
AFLOW prototype command	:	`aflow --proto=ABC8D3_aP26_2_i_i_8i_3i --params=`$a, b/a, c/a, \alpha, \beta, \gamma, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}$

- We used the 13 K data from (Smith, 1986), however they present their data in space group $C\bar{1}$, which doubles the primitive unit cell compared to the standard space group $P\bar{1}$ #2. We used FINDSYM to convert from the presented cell to the conventional cell. This involved a rotation of the cell, *e.g.*, the original c axis is the a axis in our standard primitive cell.
- Technically this is "low" albite. In high albite the silicon and aluminum atoms are mixed over all four sites, as in sandine ($S6_7$). Indeed, under some conditions albite crystals are seen in the sandine structure (Winter, 1979). See the albite entry in (Downs, 2003) for other experimental work.

Triclinic primitive vectors:

$$
\begin{aligned}
\mathbf{a}_1 &= a\hat{\mathbf{x}} \\
\mathbf{a}_2 &= b\cos\gamma\,\hat{\mathbf{x}} + b\sin\gamma\,\hat{\mathbf{y}} \\
\mathbf{a}_3 &= c_x\hat{\mathbf{x}} + c_y\,\hat{\mathbf{y}} + c_z\,\hat{\mathbf{z}}
\end{aligned}
$$

$$
\begin{aligned}
c_x &= c\,\cos\beta \\
c_y &= c\,(\cos\alpha - \cos\beta\cos\gamma)/\sin\gamma \\
c_z &= \sqrt{c^2 - c_x^2 - c_y^2}
\end{aligned}
$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1a + y_1b\cos\gamma + z_1c_x)\,\hat{\mathbf{x}} + \left(y_1b\sin\gamma + z_1c_y\right)\,\hat{\mathbf{y}} + z_1c_z\,\hat{\mathbf{z}}$	$(2i)$	Al
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$(-x_1a - y_1b\cos\gamma - z_1c_x)\,\hat{\mathbf{x}} + \left(-y_1b\sin\gamma - z_1c_y\right)\,\hat{\mathbf{y}} - z_1c_z\,\hat{\mathbf{z}}$	$(2i)$	Al
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2a + y_2b\cos\gamma + z_2c_x)\,\hat{\mathbf{x}} + \left(y_2b\sin\gamma + z_2c_y\right)\,\hat{\mathbf{y}} + z_2c_z\,\hat{\mathbf{z}}$	$(2i)$	Na
$\mathbf{B}_4$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2a - y_2b\cos\gamma - z_2c_x)\,\hat{\mathbf{x}} + \left(-y_2b\sin\gamma - z_2c_y\right)\,\hat{\mathbf{y}} - z_2c_z\,\hat{\mathbf{z}}$	$(2i)$	Na
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3a + y_3b\cos\gamma + z_3c_x)\,\hat{\mathbf{x}} + \left(y_3b\sin\gamma + z_3c_y\right)\,\hat{\mathbf{y}} + z_3c_z\,\hat{\mathbf{z}}$	$(2i)$	O I
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3a - y_3b\cos\gamma - z_3c_x)\,\hat{\mathbf{x}} + \left(-y_3b\sin\gamma - z_3c_y\right)\,\hat{\mathbf{y}} - z_3c_z\,\hat{\mathbf{z}}$	$(2i)$	O I
$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4a + y_4b\cos\gamma + z_4c_x)\,\hat{\mathbf{x}} + \left(y_4b\sin\gamma + z_4c_y\right)\,\hat{\mathbf{y}} + z_4c_z\,\hat{\mathbf{z}}$	$(2i)$	O II
$\mathbf{B}_8$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4a - y_4b\cos\gamma - z_4c_x)\,\hat{\mathbf{x}} + \left(-y_4b\sin\gamma - z_4c_y\right)\,\hat{\mathbf{y}} - z_4c_z\,\hat{\mathbf{z}}$	$(2i)$	O II
$\mathbf{B}_9$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5a + y_5b\cos\gamma + z_5c_x)\,\hat{\mathbf{x}} + \left(y_5b\sin\gamma + z_5c_y\right)\,\hat{\mathbf{y}} + z_5c_z\,\hat{\mathbf{z}}$	$(2i)$	O III
$\mathbf{B}_{10}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5a - y_5b\cos\gamma - z_5c_x)\,\hat{\mathbf{x}} + \left(-y_5b\sin\gamma - z_5c_y\right)\,\hat{\mathbf{y}} - z_5c_z\,\hat{\mathbf{z}}$	$(2i)$	O III
$\mathbf{B}_{11}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6a + y_6b\cos\gamma + z_6c_x)\,\hat{\mathbf{x}} + \left(y_6b\sin\gamma + z_6c_y\right)\,\hat{\mathbf{y}} + z_6c_z\,\hat{\mathbf{z}}$	$(2i)$	O IV
$\mathbf{B}_{12}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6a - y_6b\cos\gamma - z_6c_x)\,\hat{\mathbf{x}} + \left(-y_6b\sin\gamma - z_6c_y\right)\,\hat{\mathbf{y}} - z_6c_z\,\hat{\mathbf{z}}$	$(2i)$	O IV
$\mathbf{B}_{13}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7a + y_7b\cos\gamma + z_7c_x)\,\hat{\mathbf{x}} + \left(y_7b\sin\gamma + z_7c_y\right)\,\hat{\mathbf{y}} + z_7c_z\,\hat{\mathbf{z}}$	$(2i)$	O V
$\mathbf{B}_{14}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7a - y_7b\cos\gamma - z_7c_x)\,\hat{\mathbf{x}} + \left(-y_7b\sin\gamma - z_7c_y\right)\,\hat{\mathbf{y}} - z_7c_z\,\hat{\mathbf{z}}$	$(2i)$	O V
$\mathbf{B}_{15}$	=	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8a + y_8b\cos\gamma + z_8c_x)\,\hat{\mathbf{x}} + \left(y_8b\sin\gamma + z_8c_y\right)\,\hat{\mathbf{y}} + z_8c_z\,\hat{\mathbf{z}}$	$(2i)$	O VI
$\mathbf{B}_{16}$	=	$-x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8a - y_8b\cos\gamma - z_8c_x)\,\hat{\mathbf{x}} + \left(-y_8b\sin\gamma - z_8c_y\right)\,\hat{\mathbf{y}} - z_8c_z\,\hat{\mathbf{z}}$	$(2i)$	O VI

$\mathbf{B}_{17}$	=	$x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$(x_9a + y_9b\cos\gamma + z_9c_x)\,\hat{\mathbf{x}} + \left(y_9b\sin\gamma + z_9c_y\right)\hat{\mathbf{y}} + z_9c_z\,\hat{\mathbf{z}}$	(2*i*)	O VII
$\mathbf{B}_{18}$	=	$-x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$(-x_9a - y_9b\cos\gamma - z_9c_x)\,\hat{\mathbf{x}} + \left(-y_9b\sin\gamma - z_9c_y\right)\hat{\mathbf{y}} - z_9c_z\,\hat{\mathbf{z}}$	(2*i*)	O VII
$\mathbf{B}_{19}$	=	$x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$(x_{10}a + y_{10}b\cos\gamma + z_{10}c_x)\,\hat{\mathbf{x}} + \left(y_{10}b\sin\gamma + z_{10}c_y\right)\hat{\mathbf{y}} + z_{10}c_z\,\hat{\mathbf{z}}$	(2*i*)	O VIII
$\mathbf{B}_{20}$	=	$-x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3$	=	$(-x_{10}a - y_{10}b\cos\gamma - z_{10}c_x)\,\hat{\mathbf{x}} + \left(-y_{10}b\sin\gamma - z_{10}c_y\right)\hat{\mathbf{y}} - z_{10}c_z\,\hat{\mathbf{z}}$	(2*i*)	O VIII
$\mathbf{B}_{21}$	=	$x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$(x_{11}a + y_{11}b\cos\gamma + z_{11}c_x)\,\hat{\mathbf{x}} + \left(y_{11}b\sin\gamma + z_{11}c_y\right)\hat{\mathbf{y}} + z_{11}c_z\,\hat{\mathbf{z}}$	(2*i*)	Si I
$\mathbf{B}_{22}$	=	$-x_{11}\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3$	=	$(-x_{11}a - y_{11}b\cos\gamma - z_{11}c_x)\,\hat{\mathbf{x}} + \left(-y_{11}b\sin\gamma - z_{11}c_y\right)\hat{\mathbf{y}} - z_{11}c_z\,\hat{\mathbf{z}}$	(2*i*)	Si I
$\mathbf{B}_{23}$	=	$x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3$	=	$(x_{12}a + y_{12}b\cos\gamma + z_{12}c_x)\,\hat{\mathbf{x}} + \left(y_{12}b\sin\gamma + z_{12}c_y\right)\hat{\mathbf{y}} + z_{12}c_z\,\hat{\mathbf{z}}$	(2*i*)	Si II
$\mathbf{B}_{24}$	=	$-x_{12}\,\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3$	=	$(-x_{12}a - y_{12}b\cos\gamma - z_{12}c_x)\,\hat{\mathbf{x}} + \left(-y_{12}b\sin\gamma - z_{12}c_y\right)\hat{\mathbf{y}} - z_{12}c_z\,\hat{\mathbf{z}}$	(2*i*)	Si II
$\mathbf{B}_{25}$	=	$x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3$	=	$(x_{13}a + y_{13}b\cos\gamma + z_{13}c_x)\,\hat{\mathbf{x}} + \left(y_{13}b\sin\gamma + z_{13}c_y\right)\hat{\mathbf{y}} + z_{13}c_z\,\hat{\mathbf{z}}$	(2*i*)	Si III
$\mathbf{B}_{26}$	=	$-x_{13}\,\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 - z_{13}\,\mathbf{a}_3$	=	$(-x_{13}a - y_{13}b\cos\gamma - z_{13}c_x)\,\hat{\mathbf{x}} + \left(-y_{13}b\sin\gamma - z_{13}c_y\right)\hat{\mathbf{y}} - z_{13}c_z\,\hat{\mathbf{z}}$	(2*i*)	Si III

References:

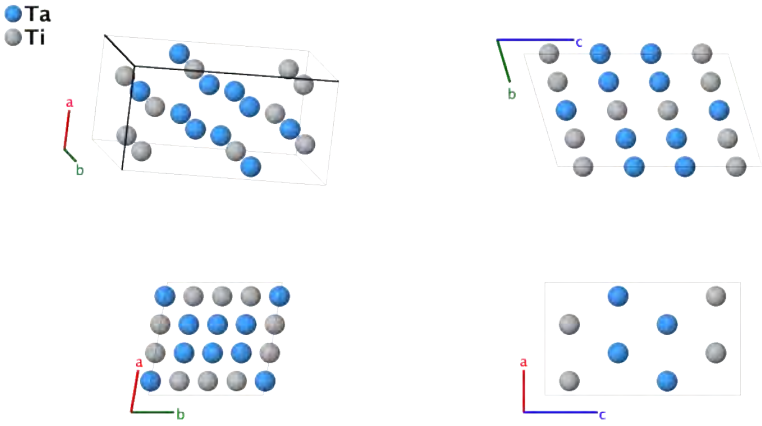
- J. V. Smith, G. Artioli, and Å. Kvick, *Low albite, $NaAlSi_3O_8$: Neutron diffraction study of crystal structure at 13 K*, Am. Mineral. **71**, 727–733 (1986).
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).
- J. K. Winter, F. P. Okamura, and S. Ghose, *A high-temperature structural study of high albite, monalbite, and the analbite → monalbite phase transition*, Am. Mineral. **64**, 409–423 (1979).

Geometry files:

- CIF: pp. 1506
- POSCAR: pp. 1506

TaTi (BCC SQS-16) Structure: AB_aP16_2_4i_4i

http://aflow.org/prototype-encyclopedia/AB_aP16_2_4i_4i



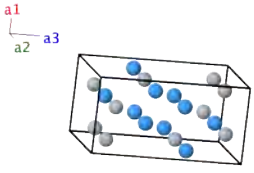
Prototype	:	TaTi
AFLOW prototype label	:	AB_aP16_2_4i_4i
***Strukturbericht* designation**	:	None
Pearson symbol	:	aP16
Space group number	:	2
Space group symbol	:	$P\bar{1}$
AFLOW prototype command	:	`aflow --proto=AB_aP16_2_4i_4i --params=`$a, b/a, c/a, \alpha, \beta, \gamma, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8$

- This is a special quasirandom structure with 16 atoms per unit cell (SQS-16) for a bcc binary substitutional alloy A_xB_{1-x} (Jiang, 2004). This prototype is the equicompositional structure ($x = 0.5$). The $x = 0.25$ and $x = 0.75$ structures are given by AB3_mC32_8_4a_12a.

Triclinic primitive vectors:

$$\mathbf{a}_1 = a\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\cos\gamma\,\hat{\mathbf{x}} + b\sin\gamma\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c_x\hat{\mathbf{x}} + c_y\,\hat{\mathbf{y}} + c_z\,\hat{\mathbf{z}}$$

$$c_x = c\,\cos\beta$$
$$c_y = c\,(\cos\alpha - \cos\beta\cos\gamma)/\sin\gamma$$
$$c_z = \sqrt{c^2 - c_x^2 - c_y^2}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$= x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	$= (x_1a + y_1b\cos\gamma + z_1c_x)\,\hat{\mathbf{x}} + \left(y_1b\sin\gamma + z_1c_y\right)\hat{\mathbf{y}} + z_1c_z\,\hat{\mathbf{z}}$	$(2i)$	Ta I
$\mathbf{B}_2$	$= -x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	$= (-x_1a - y_1b\cos\gamma - z_1c_x)\,\hat{\mathbf{x}} + \left(-y_1b\sin\gamma - z_1c_y\right)\hat{\mathbf{y}} - z_1c_z\,\hat{\mathbf{z}}$	$(2i)$	Ta I

$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2a + y_2b\cos\gamma + z_2c_x)\,\hat{\mathbf{x}} +$ $\left(y_2b\sin\gamma + z_2c_y\right)\,\hat{\mathbf{y}} + z_2c_z\,\hat{\mathbf{z}}$	$(2i)$	Ta II
$\mathbf{B}_4$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2a - y_2b\cos\gamma - z_2c_x)\,\hat{\mathbf{x}} +$ $\left(-y_2b\sin\gamma - z_2c_y\right)\,\hat{\mathbf{y}} - z_2c_z\,\hat{\mathbf{z}}$	$(2i)$	Ta II
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3a + y_3b\cos\gamma + z_3c_x)\,\hat{\mathbf{x}} +$ $\left(y_3b\sin\gamma + z_3c_y\right)\,\hat{\mathbf{y}} + z_3c_z\,\hat{\mathbf{z}}$	$(2i)$	Ta III
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3a - y_3b\cos\gamma - z_3c_x)\,\hat{\mathbf{x}} +$ $\left(-y_3b\sin\gamma - z_3c_y\right)\,\hat{\mathbf{y}} - z_3c_z\,\hat{\mathbf{z}}$	$(2i)$	Ta III
$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4a + y_4b\cos\gamma + z_4c_x)\,\hat{\mathbf{x}} +$ $\left(y_4b\sin\gamma + z_4c_y\right)\,\hat{\mathbf{y}} + z_4c_z\,\hat{\mathbf{z}}$	$(2i)$	Ta IV
$\mathbf{B}_8$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4a - y_4b\cos\gamma - z_4c_x)\,\hat{\mathbf{x}} +$ $\left(-y_4b\sin\gamma - z_4c_y\right)\,\hat{\mathbf{y}} - z_4c_z\,\hat{\mathbf{z}}$	$(2i)$	Ta IV
$\mathbf{B}_9$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5a + y_5b\cos\gamma + z_5c_x)\,\hat{\mathbf{x}} +$ $\left(y_5b\sin\gamma + z_5c_y\right)\,\hat{\mathbf{y}} + z_5c_z\,\hat{\mathbf{z}}$	$(2i)$	Ti I
$\mathbf{B}_{10}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5a - y_5b\cos\gamma - z_5c_x)\,\hat{\mathbf{x}} +$ $\left(-y_5b\sin\gamma - z_5c_y\right)\,\hat{\mathbf{y}} - z_5c_z\,\hat{\mathbf{z}}$	$(2i)$	Ti I
$\mathbf{B}_{11}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6a + y_6b\cos\gamma + z_6c_x)\,\hat{\mathbf{x}} +$ $\left(y_6b\sin\gamma + z_6c_y\right)\,\hat{\mathbf{y}} + z_6c_z\,\hat{\mathbf{z}}$	$(2i)$	Ti II
$\mathbf{B}_{12}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6a - y_6b\cos\gamma - z_6c_x)\,\hat{\mathbf{x}} +$ $\left(-y_6b\sin\gamma - z_6c_y\right)\,\hat{\mathbf{y}} - z_6c_z\,\hat{\mathbf{z}}$	$(2i)$	Ti II
$\mathbf{B}_{13}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7a + y_7b\cos\gamma + z_7c_x)\,\hat{\mathbf{x}} +$ $\left(y_7b\sin\gamma + z_7c_y\right)\,\hat{\mathbf{y}} + z_7c_z\,\hat{\mathbf{z}}$	$(2i)$	Ti III
$\mathbf{B}_{14}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7a - y_7b\cos\gamma - z_7c_x)\,\hat{\mathbf{x}} +$ $\left(-y_7b\sin\gamma - z_7c_y\right)\,\hat{\mathbf{y}} - z_7c_z\,\hat{\mathbf{z}}$	$(2i)$	Ti III
$\mathbf{B}_{15}$	=	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8a + y_8b\cos\gamma + z_8c_x)\,\hat{\mathbf{x}} +$ $\left(y_8b\sin\gamma + z_8c_y\right)\,\hat{\mathbf{y}} + z_8c_z\,\hat{\mathbf{z}}$	$(2i)$	Ti IV
$\mathbf{B}_{16}$	=	$-x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8a - y_8b\cos\gamma - z_8c_x)\,\hat{\mathbf{x}} +$ $\left(-y_8b\sin\gamma - z_8c_y\right)\,\hat{\mathbf{y}} - z_8c_z\,\hat{\mathbf{z}}$	$(2i)$	Ti IV

References:

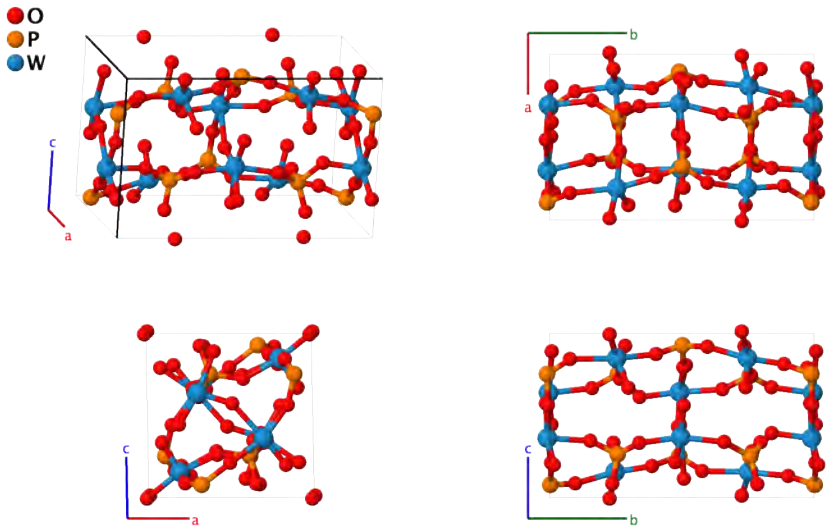
- C. Jiang, C. Wolverton, J. Sofo, L.-Q. Chen, and Z.-K. Liu, *First-principles study of binary bcc alloys using special quasirandom structures*, Phys. Rev. B **69**, 214202 (2004), doi:10.1103/PhysRevB.69.214202.

Geometry files:

- CIF: pp. 1506
- POSCAR: pp. 1507

$W_2O_3(PO_4)_2$ Structure: A11B2C2_mP60_4_22a_4a_4a

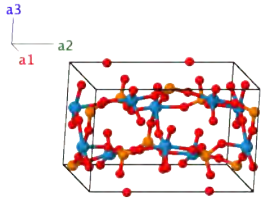
http://aflow.org/prototype-encyclopedia/A11B2C2_mP60_4_22a_4a_4a



Prototype	:	$O_{11}P_2W_2$
AFLOW prototype label	:	A11B2C2_mP60_4_22a_4a_4a
***Strukturbericht* designation**	:	None
Pearson symbol	:	mP60
Space group number	:	4
Space group symbol	:	$P2_1$
AFLOW prototype command	:	`aflow --proto=A11B2C2_mP60_4_22a_4a_4a --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}, x_{17}, y_{17}, z_{17}, x_{18}, y_{18}, z_{18}, x_{19}, y_{19}, z_{19}, x_{20}, y_{20}, z_{20}, x_{21}, y_{21}, z_{21}, x_{22}, y_{22}, z_{22}, x_{23}, y_{23}, z_{23}, x_{24}, y_{24}, z_{24}, x_{25}, y_{25}, z_{25}, x_{26}, y_{26}, z_{26}, x_{27}, y_{27}, z_{27}, x_{28}, y_{28}, z_{28}, x_{29}, y_{29}, z_{29}, x_{30}, y_{30}, z_{30}$

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	O I
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2} + y_1\right)\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$(-x_1 a - z_1 c\cos\beta)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_1\right) b\,\hat{\mathbf{y}} - z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	O I
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	O II

$$\mathbf{B}_4 = -x_2\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_2\right)\mathbf{a}_2 - z_2\,\mathbf{a}_3 = (-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_2\right) b\,\hat{\mathbf{y}} - z_2 c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O II}$$

$$\mathbf{B}_5 = x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3 = (x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O III}$$

$$\mathbf{B}_6 = -x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_3\right)\mathbf{a}_2 - z_3\,\mathbf{a}_3 = (-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_3\right) b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O III}$$

$$\mathbf{B}_7 = x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3 = (x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O IV}$$

$$\mathbf{B}_8 = -x_4\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3 = (-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_4\right) b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O IV}$$

$$\mathbf{B}_9 = x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3 = (x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O V}$$

$$\mathbf{B}_{10} = -x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3 = (-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_5\right) b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O V}$$

$$\mathbf{B}_{11} = x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 = (x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O VI}$$

$$\mathbf{B}_{12} = -x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3 = (-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_6\right) b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O VI}$$

$$\mathbf{B}_{13} = x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 = (x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O VII}$$

$$\mathbf{B}_{14} = -x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_7\right)\mathbf{a}_2 - z_7\,\mathbf{a}_3 = (-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_7\right) b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O VII}$$

$$\mathbf{B}_{15} = x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 = (x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O VIII}$$

$$\mathbf{B}_{16} = -x_8\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_8\right)\mathbf{a}_2 - z_8\,\mathbf{a}_3 = (-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_8\right) b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O VIII}$$

$$\mathbf{B}_{17} = x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 = (x_9 a + z_9 c\cos\beta)\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O IX}$$

$$\mathbf{B}_{18} = -x_9\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_9\right)\mathbf{a}_2 - z_9\,\mathbf{a}_3 = (-x_9 a - z_9 c\cos\beta)\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_9\right) b\,\hat{\mathbf{y}} - z_9 c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O IX}$$

$$\mathbf{B}_{19} = x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 = (x_{10} a + z_{10} c\cos\beta)\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O X}$$

$$\mathbf{B}_{20} = -x_{10}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{10}\right)\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 = (-x_{10} a - z_{10} c\cos\beta)\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{10}\right) b\,\hat{\mathbf{y}} - z_{10} c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O X}$$

$$\mathbf{B}_{21} = x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 = (x_{11} a + z_{11} c\cos\beta)\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + z_{11} c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XI}$$

$$\mathbf{B}_{22} = -x_{11}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{11}\right)\mathbf{a}_2 - z_{11}\,\mathbf{a}_3 = (-x_{11} a - z_{11} c\cos\beta)\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{11}\right) b\,\hat{\mathbf{y}} - z_{11} c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XI}$$

$$\mathbf{B}_{23} = x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 = (x_{12} a + z_{12} c\cos\beta)\,\hat{\mathbf{x}} + y_{12} b\,\hat{\mathbf{y}} + z_{12} c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XII}$$

$$\mathbf{B}_{24} = -x_{12}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{12}\right)\mathbf{a}_2 - z_{12}\,\mathbf{a}_3 = (-x_{12} a - z_{12} c\cos\beta)\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{12}\right) b\,\hat{\mathbf{y}} - z_{12} c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XII}$$

$$\mathbf{B}_{25} = x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3 = (x_{13} a + z_{13} c\cos\beta)\,\hat{\mathbf{x}} + y_{13} b\,\hat{\mathbf{y}} + z_{13} c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XIII}$$

$$\mathbf{B}_{26} = -x_{13}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{13}\right)\mathbf{a}_2 - z_{13}\,\mathbf{a}_3 = (-x_{13} a - z_{13} c\cos\beta)\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{13}\right) b\,\hat{\mathbf{y}} - z_{13} c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XIII}$$

$$\mathbf{B}_{27} = x_{14}\,\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3 = (x_{14} a + z_{14} c\cos\beta)\,\hat{\mathbf{x}} + y_{14} b\,\hat{\mathbf{y}} + z_{14} c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XIV}$$

$$\mathbf{B}_{28} = -x_{14}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{14}\right)\mathbf{a}_2 - z_{14}\,\mathbf{a}_3 = (-x_{14} a - z_{14} c\cos\beta)\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{14}\right) b\,\hat{\mathbf{y}} - z_{14} c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XIV}$$

$$\mathbf{B}_{29} = x_{15}\,\mathbf{a}_1 + y_{15}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 = (x_{15} a + z_{15} c\cos\beta)\,\hat{\mathbf{x}} + y_{15} b\,\hat{\mathbf{y}} + z_{15} c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XV}$$

$$\mathbf{B}_{30} = -x_{15}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{15}\right)\mathbf{a}_2 - z_{15}\,\mathbf{a}_3 = (-x_{15}a - z_{15}c\cos\beta)\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{15}\right)b\,\hat{\mathbf{y}} - z_{15}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XV}$$

$$\mathbf{B}_{31} = x_{16}\,\mathbf{a}_1 + y_{16}\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 = (x_{16}a + z_{16}c\cos\beta)\,\hat{\mathbf{x}} + y_{16}b\,\hat{\mathbf{y}} + z_{16}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XVI}$$

$$\mathbf{B}_{32} = -x_{16}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{16}\right)\mathbf{a}_2 - z_{16}\,\mathbf{a}_3 = (-x_{16}a - z_{16}c\cos\beta)\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{16}\right)b\,\hat{\mathbf{y}} - z_{16}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XVI}$$

$$\mathbf{B}_{33} = x_{17}\,\mathbf{a}_1 + y_{17}\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 = (x_{17}a + z_{17}c\cos\beta)\,\hat{\mathbf{x}} + y_{17}b\,\hat{\mathbf{y}} + z_{17}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XVII}$$

$$\mathbf{B}_{34} = -x_{17}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{17}\right)\mathbf{a}_2 - z_{17}\,\mathbf{a}_3 = (-x_{17}a - z_{17}c\cos\beta)\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{17}\right)b\,\hat{\mathbf{y}} - z_{17}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XVII}$$

$$\mathbf{B}_{35} = x_{18}\,\mathbf{a}_1 + y_{18}\,\mathbf{a}_2 + z_{18}\,\mathbf{a}_3 = (x_{18}a + z_{18}c\cos\beta)\,\hat{\mathbf{x}} + y_{18}b\,\hat{\mathbf{y}} + z_{18}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XVIII}$$

$$\mathbf{B}_{36} = -x_{18}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{18}\right)\mathbf{a}_2 - z_{18}\,\mathbf{a}_3 = (-x_{18}a - z_{18}c\cos\beta)\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{18}\right)b\,\hat{\mathbf{y}} - z_{18}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XVIII}$$

$$\mathbf{B}_{37} = x_{19}\,\mathbf{a}_1 + y_{19}\,\mathbf{a}_2 + z_{19}\,\mathbf{a}_3 = (x_{19}a + z_{19}c\cos\beta)\,\hat{\mathbf{x}} + y_{19}b\,\hat{\mathbf{y}} + z_{19}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XIX}$$

$$\mathbf{B}_{38} = -x_{19}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{19}\right)\mathbf{a}_2 - z_{19}\,\mathbf{a}_3 = (-x_{19}a - z_{19}c\cos\beta)\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{19}\right)b\,\hat{\mathbf{y}} - z_{19}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XIX}$$

$$\mathbf{B}_{39} = x_{20}\,\mathbf{a}_1 + y_{20}\,\mathbf{a}_2 + z_{20}\,\mathbf{a}_3 = (x_{20}a + z_{20}c\cos\beta)\,\hat{\mathbf{x}} + y_{20}b\,\hat{\mathbf{y}} + z_{20}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XX}$$

$$\mathbf{B}_{40} = -x_{20}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{20}\right)\mathbf{a}_2 - z_{20}\,\mathbf{a}_3 = (-x_{20}a - z_{20}c\cos\beta)\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{20}\right)b\,\hat{\mathbf{y}} - z_{20}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XX}$$

$$\mathbf{B}_{41} = x_{21}\,\mathbf{a}_1 + y_{21}\,\mathbf{a}_2 + z_{21}\,\mathbf{a}_3 = (x_{21}a + z_{21}c\cos\beta)\,\hat{\mathbf{x}} + y_{21}b\,\hat{\mathbf{y}} + z_{21}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXI}$$

$$\mathbf{B}_{42} = -x_{21}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{21}\right)\mathbf{a}_2 - z_{21}\,\mathbf{a}_3 = (-x_{21}a - z_{21}c\cos\beta)\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{21}\right)b\,\hat{\mathbf{y}} - z_{21}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXI}$$

$$\mathbf{B}_{43} = x_{22}\,\mathbf{a}_1 + y_{22}\,\mathbf{a}_2 + z_{22}\,\mathbf{a}_3 = (x_{22}a + z_{22}c\cos\beta)\,\hat{\mathbf{x}} + y_{22}b\,\hat{\mathbf{y}} + z_{22}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXII}$$

$$\mathbf{B}_{44} = -x_{22}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{22}\right)\mathbf{a}_2 - z_{22}\,\mathbf{a}_3 = (-x_{22}a - z_{22}c\cos\beta)\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{22}\right)b\,\hat{\mathbf{y}} - z_{22}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXII}$$

$$\mathbf{B}_{45} = x_{23}\,\mathbf{a}_1 + y_{23}\,\mathbf{a}_2 + z_{23}\,\mathbf{a}_3 = (x_{23}a + z_{23}c\cos\beta)\,\hat{\mathbf{x}} + y_{23}b\,\hat{\mathbf{y}} + z_{23}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{P I}$$

$$\mathbf{B}_{46} = -x_{23}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{23}\right)\mathbf{a}_2 - z_{23}\,\mathbf{a}_3 = (-x_{23}a - z_{23}c\cos\beta)\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{23}\right)b\,\hat{\mathbf{y}} - z_{23}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{P I}$$

$$\mathbf{B}_{47} = x_{24}\,\mathbf{a}_1 + y_{24}\,\mathbf{a}_2 + z_{24}\,\mathbf{a}_3 = (x_{24}a + z_{24}c\cos\beta)\,\hat{\mathbf{x}} + y_{24}b\,\hat{\mathbf{y}} + z_{24}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{P II}$$

$$\mathbf{B}_{48} = -x_{24}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{24}\right)\mathbf{a}_2 - z_{24}\,\mathbf{a}_3 = (-x_{24}a - z_{24}c\cos\beta)\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{24}\right)b\,\hat{\mathbf{y}} - z_{24}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{P II}$$

$$\mathbf{B}_{49} = x_{25}\,\mathbf{a}_1 + y_{25}\,\mathbf{a}_2 + z_{25}\,\mathbf{a}_3 = (x_{25}a + z_{25}c\cos\beta)\,\hat{\mathbf{x}} + y_{25}b\,\hat{\mathbf{y}} + z_{25}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{P III}$$

$$\mathbf{B}_{50} = -x_{25}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{25}\right)\mathbf{a}_2 - z_{25}\,\mathbf{a}_3 = (-x_{25}a - z_{25}c\cos\beta)\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{25}\right)b\,\hat{\mathbf{y}} - z_{25}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{P III}$$

$$\mathbf{B}_{51} = x_{26}\,\mathbf{a}_1 + y_{26}\,\mathbf{a}_2 + z_{26}\,\mathbf{a}_3 = (x_{26}a + z_{26}c\cos\beta)\,\hat{\mathbf{x}} + y_{26}b\,\hat{\mathbf{y}} + z_{26}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{P IV}$$

$$\mathbf{B}_{52} = -x_{26}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{26}\right)\mathbf{a}_2 - z_{26}\,\mathbf{a}_3 = (-x_{26}a - z_{26}c\cos\beta)\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{26}\right)b\,\hat{\mathbf{y}} - z_{26}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{P IV}$$

$$\mathbf{B}_{53} = x_{27}\,\mathbf{a}_1 + y_{27}\,\mathbf{a}_2 + z_{27}\,\mathbf{a}_3 = (x_{27}a + z_{27}c\cos\beta)\,\hat{\mathbf{x}} + y_{27}b\,\hat{\mathbf{y}} + z_{27}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{W I}$$

$$\mathbf{B}_{54} = -x_{27}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{27}\right)\mathbf{a}_2 - z_{27}\,\mathbf{a}_3 = (-x_{27}a - z_{27}c\cos\beta)\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{27}\right)b\,\hat{\mathbf{y}} - z_{27}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{W I}$$

$$\mathbf{B}_{55} = x_{28}\,\mathbf{a}_1 + y_{28}\,\mathbf{a}_2 + z_{28}\,\mathbf{a}_3 = (x_{28}a + z_{28}c\cos\beta)\,\hat{\mathbf{x}} + y_{28}b\,\hat{\mathbf{y}} + z_{28}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{W II}$$

$$\mathbf{B}_{56} = -x_{28}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{28}\right)\mathbf{a}_2 - z_{28}\,\mathbf{a}_3 = (-x_{28}a - z_{28}c\cos\beta)\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{28}\right)b\,\hat{\mathbf{y}} - z_{28}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{W II}$$

$$\mathbf{B}_{57} = x_{29}\,\mathbf{a}_1 + y_{29}\,\mathbf{a}_2 + z_{29}\,\mathbf{a}_3 = (x_{29}a + z_{29}c\cos\beta)\,\hat{\mathbf{x}} + y_{29}b\,\hat{\mathbf{y}} + z_{29}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{W III}$$

$$\mathbf{B}_{58} = -x_{29}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{29}\right)\mathbf{a}_2 - z_{29}\,\mathbf{a}_3 = (-x_{29}a - z_{29}c\cos\beta)\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{29}\right)b\,\hat{\mathbf{y}} - z_{29}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{W III}$$

$$\mathbf{B}_{59} = x_{30}\,\mathbf{a}_1 + y_{30}\,\mathbf{a}_2 + z_{30}\,\mathbf{a}_3 = (x_{30}a + z_{30}c\cos\beta)\,\hat{\mathbf{x}} + y_{30}b\,\hat{\mathbf{y}} + z_{30}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{W IV}$$

$$\mathbf{B}_{60} = -x_{30}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{30}\right)\mathbf{a}_2 - z_{30}\,\mathbf{a}_3 = (-x_{30}a - z_{30}c\cos\beta)\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{30}\right)b\,\hat{\mathbf{y}} - z_{30}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{W IV}$$

References:

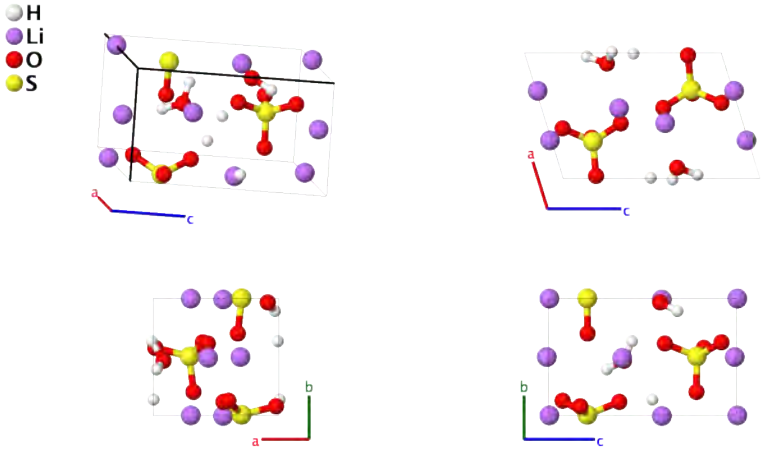
- P. Kierkegaard and S. Åsbrink, *The Crystal Structure of $W_2O_3(PO_4)_2$. Determination of a Superstructure by Means of Least-Squares Calculations*, Acta Chem. Scand. **18**, 2329–2338 (1964), doi:10.3891/acta.chem.scand.18-2329.

Geometry files:

- CIF: pp. 1507
- POSCAR: pp. 1507

$Li_2SO_4 \cdot H_2O$ ($H4_8$) Structure: A2B2C5D_mP20_4_2a_2a_5a_a

http://aflow.org/prototype-encyclopedia/A2B2C5D_mP20_4_2a_2a_5a_a

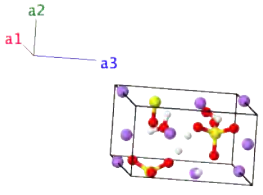


Prototype	:	$H_2Li_2O_5S$
AFLOW prototype label	:	A2B2C5D_mP20_4_2a_2a_5a_a
***Strukturbericht* designation**	:	$H4_8$
Pearson symbol	:	mP20
Space group number	:	4
Space group symbol	:	$P2_1$
AFLOW prototype command	:	`aflow --proto=A2B2C5D_mP20_4_2a_2a_5a_a --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}$

- We use the data from (Lundgren, 1984) at 20 K, including the positions of the hydrogen atoms not found in the original $H4_8$ structure in (Gottfried, 1937).
- Space group $P2_1$ #4 allows the y coordinates to have an arbitrary origin. We follow (Lundgren, 1984) and set the y coordinate of the sulfur atom, y_{10}, to zero.

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	H I

$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2} + y_1\right)\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$(-x_1 a - z_1 c\cos\beta)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_1\right) b\,\hat{\mathbf{y}} - z_1 c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	H I
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	H II
$\mathbf{B}_4$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right) b\,\hat{\mathbf{y}} - z_2 c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	H II
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	Li I
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right) b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	Li I
$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	Li II
$\mathbf{B}_8$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right) b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	Li II
$\mathbf{B}_9$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	O I
$\mathbf{B}_{10}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right) b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	O I
$\mathbf{B}_{11}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	O II
$\mathbf{B}_{12}$	=	$-x_6\,\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right) b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	O II
$\mathbf{B}_{13}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	O III
$\mathbf{B}_{14}$	=	$-x_7\,\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right) b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	O III
$\mathbf{B}_{15}$	=	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	O IV
$\mathbf{B}_{16}$	=	$-x_8\,\mathbf{a}_1 + \left(\frac{1}{2} + y_8\right)\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_8\right) b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	O IV
$\mathbf{B}_{17}$	=	$x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$(x_9 a + z_9 c\cos\beta)\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	O V
$\mathbf{B}_{18}$	=	$-x_9\,\mathbf{a}_1 + \left(\frac{1}{2} + y_9\right)\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$(-x_9 a - z_9 c\cos\beta)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_9\right) b\,\hat{\mathbf{y}} - z_9 c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	O V
$\mathbf{B}_{19}$	=	$x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$(x_{10} a + z_{10} c\cos\beta)\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	S
$\mathbf{B}_{20}$	=	$-x_{10}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{10}\right)\mathbf{a}_2 - z_{10}\,\mathbf{a}_3$	=	$(-x_{10} a - z_{10} c\cos\beta)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{10}\right) b\,\hat{\mathbf{y}} - z_{10} c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	S

References:

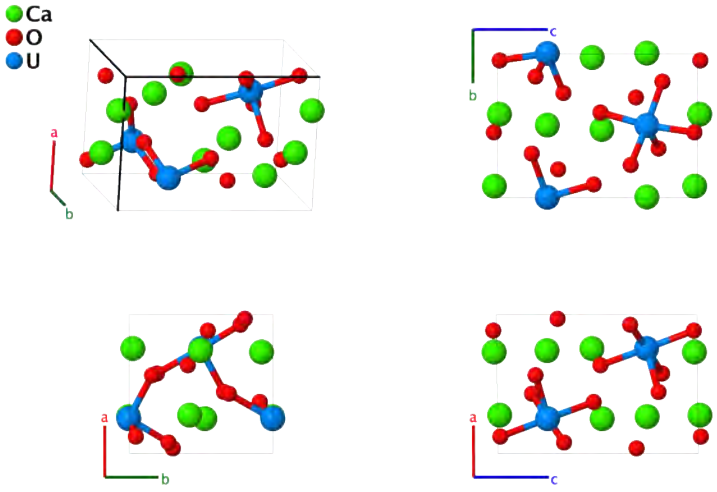
- J.-O. Lundgren, Å. Kvick, M. Karppinen, R. Liminga, and S. C. Abrahams, *Neutron diffraction structural study of pyroelectric $Li_2SO_4{\cdot}H_2O$ at 293, 80, and 20 K*, J. Chem. Phys. **80**, 423–430 (1984), doi:10.1063/1.446465.
- C. Gottfried and F. Schossberger, eds., *Strukturbericht Band III 1933-1935* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1508
- POSCAR: pp. 1508

Ca_3UO_6 Structure: A3B6C_mP20_4_3a_6a_a

http://aflow.org/prototype-encyclopedia/A3B6C_mP20_4_3a_6a_a



Prototype	:	Ca_3O_6U
AFLOW prototype label	:	A3B6C_mP20_4_3a_6a_a
***Strukturbericht* designation**	:	None
Pearson symbol	:	mP20
Space group number	:	4
Space group symbol	:	$P2_1$
AFLOW prototype command	:	`aflow --proto=A3B6C_mP20_4_3a_6a_a --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}$

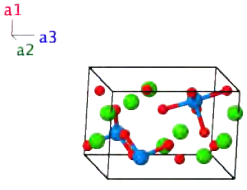
Other compounds with this structure

- Sr_3UO_6

- Space group $P2_1$ #4 allows the y coordinates to have an arbitrary origin. Here, this freedom is used to set $y_{10} = 0$ for the uranium position.

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$ =	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Ca I

$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2} + y_1\right)\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$(-x_1 a - z_1 c\cos\beta)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_1\right) b\,\hat{\mathbf{y}} - z_1 c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	Ca I
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	Ca II
$\mathbf{B}_4$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right) b\,\hat{\mathbf{y}} - z_2 c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	Ca II
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	Ca III
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right) b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	Ca III
$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	O I
$\mathbf{B}_8$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right) b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	O I
$\mathbf{B}_9$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	O II
$\mathbf{B}_{10}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right) b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	O II
$\mathbf{B}_{11}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	O III
$\mathbf{B}_{12}$	=	$-x_6\,\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right) b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	O III
$\mathbf{B}_{13}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	O IV
$\mathbf{B}_{14}$	=	$-x_7\,\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right) b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	O IV
$\mathbf{B}_{15}$	=	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	O V
$\mathbf{B}_{16}$	=	$-x_8\,\mathbf{a}_1 + \left(\frac{1}{2} + y_8\right)\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_8\right) b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	O V
$\mathbf{B}_{17}$	=	$x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$(x_9 a + z_9 c\cos\beta)\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	O VI
$\mathbf{B}_{18}$	=	$-x_9\,\mathbf{a}_1 + \left(\frac{1}{2} + y_9\right)\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$(-x_9 a - z_9 c\cos\beta)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_9\right) b\,\hat{\mathbf{y}} - z_9 c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	O VI
$\mathbf{B}_{19}$	=	$x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$(x_{10} a + z_{10} c\cos\beta)\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	U
$\mathbf{B}_{20}$	=	$-x_{10}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{10}\right)\mathbf{a}_2 - z_{10}\,\mathbf{a}_3$	=	$(-x_{10} a - z_{10} c\cos\beta)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{10}\right) b\,\hat{\mathbf{y}} - z_{10} c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	U

References:

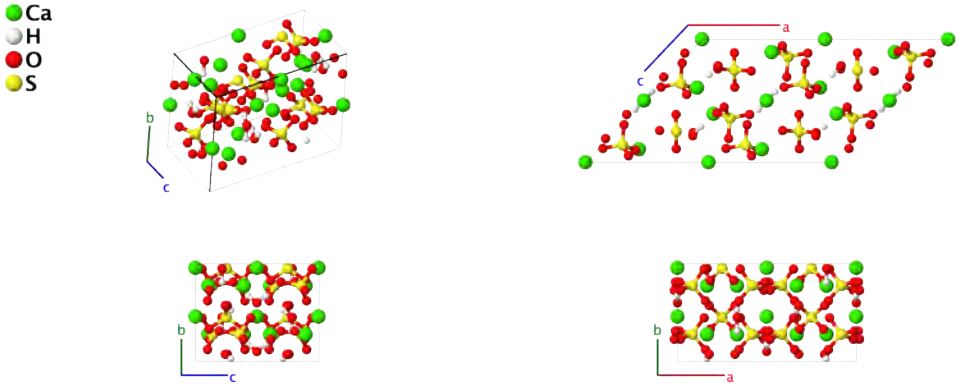
- B. O. Loopstra and H. M. Rietveld, *The structure of some alkaline-earth metal uranates*, Acta Crystallogr. Sect. B Struct. Sci. **25**, 787–791 (1969), doi:10.1107/S0567740869002974.

Geometry files:

- CIF: pp. 1508
- POSCAR: pp. 1509

Bassanite [$CaSO_4(H_2O)_{0.5}$, $H4_7$] Structure: A2B2C9D2_mC90_5_ab2c_3c_b13c_3c

http://aflow.org/prototype-encyclopedia/A2B2C9D2_mC90_5_ab2c_3c_b13c_3c

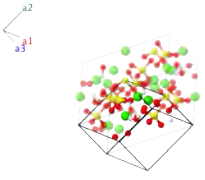


Prototype	:	$Ca_2H_2O_9S_2$
AFLOW prototype label	:	A2B2C9D2_mC90_5_ab2c_3c_b13c_3c
***Strukturbericht* designation**	:	$H4_7$
Pearson symbol	:	mC90
Space group number	:	5
Space group symbol	:	$C2$
AFLOW prototype command	:	`aflow --proto=A2B2C9D2_mC90_5_ab2c_3c_b13c_3c --params=`$a, b/a, c/a, \beta, y_1, y_2, y_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}, x_{17}, y_{17}, z_{17}, x_{18}, y_{18}, z_{18}, x_{19}, y_{19}, z_{19}, x_{20}, y_{20}, z_{20}, x_{21}, y_{21}, z_{21}, x_{22}, y_{22}, z_{22}, x_{23}, y_{23}, z_{23}, x_{24}, y_{24}, z_{24}$

- (Gottfried, 1937) gave this the *Strukturbericht* designation $H4_7$. They listed the system as monoclinic, with space group $C2$ #5, but noted that it was pseudo-hexagonal and gave the coordinates for all of the atoms except the water molecules in terms of the trigonal space group $P3_121$ #152. (Abriel, 1993) found a complete determination of the structure, in space group $I2$ #5, which we have converted to the standard $C2$ setting. The exact structure of this system seems to depend on the actual amount of water and the preparation mechanism (Singh, 2007).
- The $P3_121$ structure can be obtained from these positions by removing all of the water molecules (the hydrogen atoms plus O-I and O-XIV), and allowing for a small uncertainty in the positions of the remaining atoms.

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2$	=	$y_1 b\,\hat{\mathbf{y}}$	$(2a)$	Ca I

$\mathbf{B}_2$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	(2*b*)	Ca II
$\mathbf{B}_3$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	(2*b*)	O I
$\mathbf{B}_4$	=	$(x_4 - y_4)\,\mathbf{a}_1 + (x_4 + y_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	Ca III
$\mathbf{B}_5$	=	$(-x_4 - y_4)\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	Ca III
$\mathbf{B}_6$	=	$(x_5 - y_5)\,\mathbf{a}_1 + (x_5 + y_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	Ca IV
$\mathbf{B}_7$	=	$(-x_5 - y_5)\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	Ca IV
$\mathbf{B}_8$	=	$(x_6 - y_6)\,\mathbf{a}_1 + (x_6 + y_6)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	H I
$\mathbf{B}_9$	=	$(-x_6 - y_6)\,\mathbf{a}_1 + (-x_6 + y_6)\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	H I
$\mathbf{B}_{10}$	=	$(x_7 - y_7)\,\mathbf{a}_1 + (x_7 + y_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	H II
$\mathbf{B}_{11}$	=	$(-x_7 - y_7)\,\mathbf{a}_1 + (-x_7 + y_7)\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	H II
$\mathbf{B}_{12}$	=	$(x_8 - y_8)\,\mathbf{a}_1 + (x_8 + y_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	H III
$\mathbf{B}_{13}$	=	$(-x_8 - y_8)\,\mathbf{a}_1 + (-x_8 + y_8)\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	H III
$\mathbf{B}_{14}$	=	$(x_9 - y_9)\,\mathbf{a}_1 + (x_9 + y_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$(x_9 a + z_9 c\cos\beta)\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_{15}$	=	$(-x_9 - y_9)\,\mathbf{a}_1 + (-x_9 + y_9)\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$(-x_9 a - z_9 c\cos\beta)\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} - z_9 c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_{16}$	=	$(x_{10} - y_{10})\,\mathbf{a}_1 + (x_{10} + y_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$(x_{10} a + z_{10} c\cos\beta)\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	O III
$\mathbf{B}_{17}$	=	$(-x_{10} - y_{10})\,\mathbf{a}_1 + (-x_{10} + y_{10})\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3$	=	$(-x_{10} a - z_{10} c\cos\beta)\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} - z_{10} c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	O III
$\mathbf{B}_{18}$	=	$(x_{11} - y_{11})\,\mathbf{a}_1 + (x_{11} + y_{11})\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$(x_{11} a + z_{11} c\cos\beta)\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + z_{11} c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	O IV
$\mathbf{B}_{19}$	=	$(-x_{11} - y_{11})\,\mathbf{a}_1 + (-x_{11} + y_{11})\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3$	=	$(-x_{11} a - z_{11} c\cos\beta)\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} - z_{11} c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	O IV
$\mathbf{B}_{20}$	=	$(x_{12} - y_{12})\,\mathbf{a}_1 + (x_{12} + y_{12})\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3$	=	$(x_{12} a + z_{12} c\cos\beta)\,\hat{\mathbf{x}} + y_{12} b\,\hat{\mathbf{y}} + z_{12} c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	O V
$\mathbf{B}_{21}$	=	$(-x_{12} - y_{12})\,\mathbf{a}_1 + (-x_{12} + y_{12})\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3$	=	$(-x_{12} a - z_{12} c\cos\beta)\,\hat{\mathbf{x}} + y_{12} b\,\hat{\mathbf{y}} - z_{12} c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	O V
$\mathbf{B}_{22}$	=	$(x_{13} - y_{13})\,\mathbf{a}_1 + (x_{13} + y_{13})\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3$	=	$(x_{13} a + z_{13} c\cos\beta)\,\hat{\mathbf{x}} + y_{13} b\,\hat{\mathbf{y}} + z_{13} c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	O VI
$\mathbf{B}_{23}$	=	$(-x_{13} - y_{13})\,\mathbf{a}_1 + (-x_{13} + y_{13})\,\mathbf{a}_2 - z_{13}\,\mathbf{a}_3$	=	$(-x_{13} a - z_{13} c\cos\beta)\,\hat{\mathbf{x}} + y_{13} b\,\hat{\mathbf{y}} - z_{13} c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	O VI
$\mathbf{B}_{24}$	=	$(x_{14} - y_{14})\,\mathbf{a}_1 + (x_{14} + y_{14})\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3$	=	$(x_{14} a + z_{14} c\cos\beta)\,\hat{\mathbf{x}} + y_{14} b\,\hat{\mathbf{y}} + z_{14} c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	O VII
$\mathbf{B}_{25}$	=	$(-x_{14} - y_{14})\,\mathbf{a}_1 + (-x_{14} + y_{14})\,\mathbf{a}_2 - z_{14}\,\mathbf{a}_3$	=	$(-x_{14} a - z_{14} c\cos\beta)\,\hat{\mathbf{x}} + y_{14} b\,\hat{\mathbf{y}} - z_{14} c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	O VII
$\mathbf{B}_{26}$	=	$(x_{15} - y_{15})\,\mathbf{a}_1 + (x_{15} + y_{15})\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3$	=	$(x_{15} a + z_{15} c\cos\beta)\,\hat{\mathbf{x}} + y_{15} b\,\hat{\mathbf{y}} + z_{15} c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	O VIII

$\mathbf{B}_{27}$	$= (-x_{15} - y_{15})\,\mathbf{a}_1 + (-x_{15} + y_{15})\,\mathbf{a}_2 - z_{15}\,\mathbf{a}_3$	$= (-x_{15}a - z_{15}c\cos\beta)\,\hat{\mathbf{x}} + y_{15}b\,\hat{\mathbf{y}} - z_{15}c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	O VIII
$\mathbf{B}_{28}$	$= (x_{16} - y_{16})\,\mathbf{a}_1 + (x_{16} + y_{16})\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3$	$= (x_{16}a + z_{16}c\cos\beta)\,\hat{\mathbf{x}} + y_{16}b\,\hat{\mathbf{y}} + z_{16}c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	O IX
$\mathbf{B}_{29}$	$= (-x_{16} - y_{16})\,\mathbf{a}_1 + (-x_{16} + y_{16})\,\mathbf{a}_2 - z_{16}\,\mathbf{a}_3$	$= (-x_{16}a - z_{16}c\cos\beta)\,\hat{\mathbf{x}} + y_{16}b\,\hat{\mathbf{y}} - z_{16}c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	O IX
$\mathbf{B}_{30}$	$= (x_{17} - y_{17})\,\mathbf{a}_1 + (x_{17} + y_{17})\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3$	$= (x_{17}a + z_{17}c\cos\beta)\,\hat{\mathbf{x}} + y_{17}b\,\hat{\mathbf{y}} + z_{17}c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	O X
$\mathbf{B}_{31}$	$= (-x_{17} - y_{17})\,\mathbf{a}_1 + (-x_{17} + y_{17})\,\mathbf{a}_2 - z_{17}\,\mathbf{a}_3$	$= (-x_{17}a - z_{17}c\cos\beta)\,\hat{\mathbf{x}} + y_{17}b\,\hat{\mathbf{y}} - z_{17}c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	O X
$\mathbf{B}_{32}$	$= (x_{18} - y_{18})\,\mathbf{a}_1 + (x_{18} + y_{18})\,\mathbf{a}_2 + z_{18}\,\mathbf{a}_3$	$= (x_{18}a + z_{18}c\cos\beta)\,\hat{\mathbf{x}} + y_{18}b\,\hat{\mathbf{y}} + z_{18}c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	O XI
$\mathbf{B}_{33}$	$= (-x_{18} - y_{18})\,\mathbf{a}_1 + (-x_{18} + y_{18})\,\mathbf{a}_2 - z_{18}\,\mathbf{a}_3$	$= (-x_{18}a - z_{18}c\cos\beta)\,\hat{\mathbf{x}} + y_{18}b\,\hat{\mathbf{y}} - z_{18}c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	O XI
$\mathbf{B}_{34}$	$= (x_{19} - y_{19})\,\mathbf{a}_1 + (x_{19} + y_{19})\,\mathbf{a}_2 + z_{19}\,\mathbf{a}_3$	$= (x_{19}a + z_{19}c\cos\beta)\,\hat{\mathbf{x}} + y_{19}b\,\hat{\mathbf{y}} + z_{19}c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	O XII
$\mathbf{B}_{35}$	$= (-x_{19} - y_{19})\,\mathbf{a}_1 + (-x_{19} + y_{19})\,\mathbf{a}_2 - z_{19}\,\mathbf{a}_3$	$= (-x_{19}a - z_{19}c\cos\beta)\,\hat{\mathbf{x}} + y_{19}b\,\hat{\mathbf{y}} - z_{19}c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	O XII
$\mathbf{B}_{36}$	$= (x_{20} - y_{20})\,\mathbf{a}_1 + (x_{20} + y_{20})\,\mathbf{a}_2 + z_{20}\,\mathbf{a}_3$	$= (x_{20}a + z_{20}c\cos\beta)\,\hat{\mathbf{x}} + y_{20}b\,\hat{\mathbf{y}} + z_{20}c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	O XIII
$\mathbf{B}_{37}$	$= (-x_{20} - y_{20})\,\mathbf{a}_1 + (-x_{20} + y_{20})\,\mathbf{a}_2 - z_{20}\,\mathbf{a}_3$	$= (-x_{20}a - z_{20}c\cos\beta)\,\hat{\mathbf{x}} + y_{20}b\,\hat{\mathbf{y}} - z_{20}c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	O XIII
$\mathbf{B}_{38}$	$= (x_{21} - y_{21})\,\mathbf{a}_1 + (x_{21} + y_{21})\,\mathbf{a}_2 + z_{21}\,\mathbf{a}_3$	$= (x_{21}a + z_{21}c\cos\beta)\,\hat{\mathbf{x}} + y_{21}b\,\hat{\mathbf{y}} + z_{21}c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	O XIV
$\mathbf{B}_{39}$	$= (-x_{21} - y_{21})\,\mathbf{a}_1 + (-x_{21} + y_{21})\,\mathbf{a}_2 - z_{21}\,\mathbf{a}_3$	$= (-x_{21}a - z_{21}c\cos\beta)\,\hat{\mathbf{x}} + y_{21}b\,\hat{\mathbf{y}} - z_{21}c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	O XIV
$\mathbf{B}_{40}$	$= (x_{22} - y_{22})\,\mathbf{a}_1 + (x_{22} + y_{22})\,\mathbf{a}_2 + z_{22}\,\mathbf{a}_3$	$= (x_{22}a + z_{22}c\cos\beta)\,\hat{\mathbf{x}} + y_{22}b\,\hat{\mathbf{y}} + z_{22}c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	S I
$\mathbf{B}_{41}$	$= (-x_{22} - y_{22})\,\mathbf{a}_1 + (-x_{22} + y_{22})\,\mathbf{a}_2 - z_{22}\,\mathbf{a}_3$	$= (-x_{22}a - z_{22}c\cos\beta)\,\hat{\mathbf{x}} + y_{22}b\,\hat{\mathbf{y}} - z_{22}c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	S I
$\mathbf{B}_{42}$	$= (x_{23} - y_{23})\,\mathbf{a}_1 + (x_{23} + y_{23})\,\mathbf{a}_2 + z_{23}\,\mathbf{a}_3$	$= (x_{23}a + z_{23}c\cos\beta)\,\hat{\mathbf{x}} + y_{23}b\,\hat{\mathbf{y}} + z_{23}c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	S II
$\mathbf{B}_{43}$	$= (-x_{23} - y_{23})\,\mathbf{a}_1 + (-x_{23} + y_{23})\,\mathbf{a}_2 - z_{23}\,\mathbf{a}_3$	$= (-x_{23}a - z_{23}c\cos\beta)\,\hat{\mathbf{x}} + y_{23}b\,\hat{\mathbf{y}} - z_{23}c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	S II
$\mathbf{B}_{44}$	$= (x_{24} - y_{24})\,\mathbf{a}_1 + (x_{24} + y_{24})\,\mathbf{a}_2 + z_{24}\,\mathbf{a}_3$	$= (x_{24}a + z_{24}c\cos\beta)\,\hat{\mathbf{x}} + y_{24}b\,\hat{\mathbf{y}} + z_{24}c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	S III
$\mathbf{B}_{45}$	$= (-x_{24} - y_{24})\,\mathbf{a}_1 + (-x_{24} + y_{24})\,\mathbf{a}_2 - z_{24}\,\mathbf{a}_3$	$= (-x_{24}a - z_{24}c\cos\beta)\,\hat{\mathbf{x}} + y_{24}b\,\hat{\mathbf{y}} - z_{24}c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	S III

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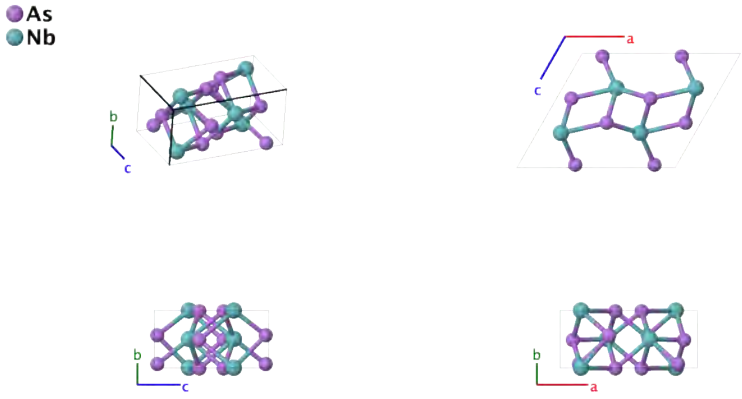
- W. Abriel and R. Nesper, *Bestimmung der Kristallstruktur von $CaSO_4(H_2O)_{0.5}$ mit Röntgenbeugungsmethoden und mit Potentialprofil-Rechnungen*, Zeitschrift für Kristallographie - Crystalline Materials **205**, 99–113 (1993), doi:10.1524/zkri.1993.205.12.99.
- C. Gottfried and F. Schossberger, eds., *Strukturbericht Band III 1933-1935* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).
- N. B. Singh and B. Middendorf, *Calcium sulphate hemihydrate hydration leading to gypsum crystallization*, Prog. Cryst. Growth Ch. **53**, 57–77 (2007), doi:10.1016/j.pcrysgrow.2007.01.002.

Found in:
- P. Ballirano, A. Maras, S. Meloni, and R. Caminiti, *The monoclinic I2 structure of bassanite, calcium sulphate hemihydrate ($CaSO_4{\cdot}0.5H_2O$)*, Eur. J. Mineral. **13**, 985–993 (2001).

Geometry files:
- CIF: pp. 1509
- POSCAR: pp. 1509

$NbAs_2$ Structure: A2B_mC12_5_2c_c

http://aflow.org/prototype-encyclopedia/A2B_mC12_5_2c_c



Prototype	:	As_2Nb
AFLOW prototype label	:	A2B_mC12_5_2c_c
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC12
Space group number	:	5
Space group symbol	:	$C2$
AFLOW prototype command	:	`aflow --proto=A2B_mC12_5_2c_c --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3$

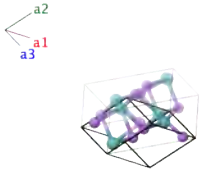
Other compounds with this structure

- $MoAs_2$, $TaAs_2$, $NbSb_2$, and $TaSb_2$

- The zero of the y-axis can be chosen arbitrarily in space group $C2$ #5. Here we chose it so that $y_3 = 1/2$ for the niobium atom.

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$= (x_1 - y_1)\,\mathbf{a}_1 + (x_1 + y_1)\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	$= (x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(4c)$	As I
$\mathbf{B}_2$	$= (-x_1 - y_1)\,\mathbf{a}_1 + (-x_1 + y_1)\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	$= (-x_1 a - z_1 c\cos\beta)\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} - z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(4c)$	As I

$\mathbf{B}_3$	=	$(x_2 - y_2)\ \mathbf{a}_1 + (x_2 + y_2)\ \mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\ \hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(4c)$	As II
$\mathbf{B}_4$	=	$(-x_2 - y_2)\ \mathbf{a}_1 + (-x_2 + y_2)\ \mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2 a - z_2 c\cos\beta)\ \hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} - z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(4c)$	As II
$\mathbf{B}_5$	=	$(x_3 - y_3)\ \mathbf{a}_1 + (x_3 + y_3)\ \mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\ \hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(4c)$	Nb
$\mathbf{B}_6$	=	$(-x_3 - y_3)\ \mathbf{a}_1 + (-x_3 + y_3)\ \mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\ \hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(4c)$	Nb

References:

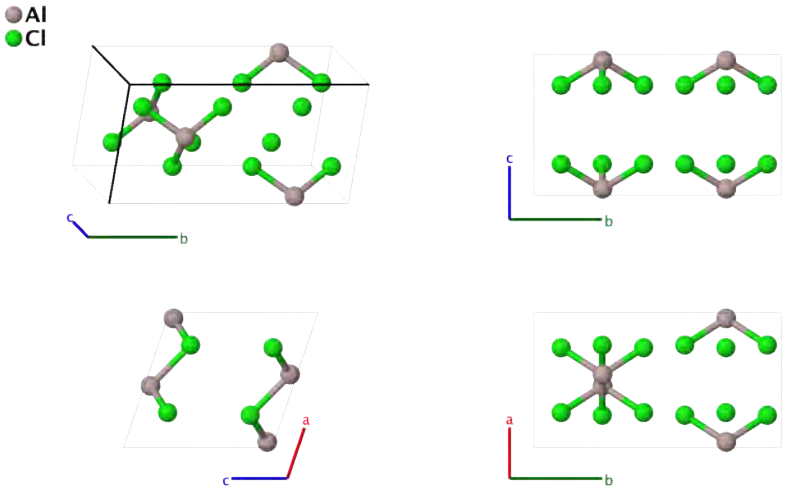
- S. Furuseth and A. Kjekshus, *The Crystal Structures of $NbAs_2$ and $NbSb_2$*, Acta Cryst. **18**, 320–324 (1965), doi:10.1107/S0365110X65000750.

Geometry files:

- CIF: pp. 1510
- POSCAR: pp. 1510

$D0_{15}$ ($AlCl_3$) (*obsolete*) Structure: AB3_mC16_5_c_3c

http://aflow.org/prototype-encyclopedia/AB3_mC16_5_c_3c

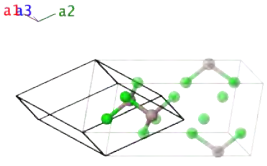


Prototype	:	$AlCl_3$
AFLOW prototype label	:	AB3_mC16_5_c_3c
***Strukturbericht* designation**	:	$D0_{15}$
Pearson symbol	:	mC16
Space group number	:	5
Space group symbol	:	$C2$
AFLOW prototype command	:	`aflow --proto=AB3_mC16_5_c_3c --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4$

- This structure was suggested by (Ketelaar, 1935) and given the *Strukturbericht* designation $D0_{15}$ by (Gottfried, 1395) with no reference to the $D0_{13}$ structure of (Laschkarew, 1930) found in (Hermann, 1937).
- Ketelaar gave a description of the unit cell a hexagonal setting, but the atomic positions resolve into the body-centered monoclinic cell shown here. This structure, however, has a lower symmetry than the currently accepted structure, which is body-centered monoclinic, space group $C2/m$ #12. We had previously designated that structure $D0_{15}$, but the current structure was given that designation by (Hermann, 1937).

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$\mathbf{B}_1$	=	$(x_1 - y_1)\,\mathbf{a}_1 + (x_1 + y_1)\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	Al
$\mathbf{B}_2$	=	$(-x_1 - y_1)\,\mathbf{a}_1 + (-x_1 + y_1)\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$(-x_1 a - z_1 c\cos\beta)\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} - z_1 c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	Al
$\mathbf{B}_3$	=	$(x_2 - y_2)\,\mathbf{a}_1 + (x_2 + y_2)\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	Cl I
$\mathbf{B}_4$	=	$(-x_2 - y_2)\,\mathbf{a}_1 + (-x_2 + y_2)\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} - z_2 c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	Cl I
$\mathbf{B}_5$	=	$(x_3 - y_3)\,\mathbf{a}_1 + (x_3 + y_3)\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	Cl II
$\mathbf{B}_6$	=	$(-x_3 - y_3)\,\mathbf{a}_1 + (-x_3 + y_3)\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	Cl II
$\mathbf{B}_7$	=	$(x_4 - y_4)\,\mathbf{a}_1 + (x_4 + y_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	Cl III
$\mathbf{B}_8$	=	$(-x_4 - y_4)\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	Cl III

References:

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- W. E. Laschkarew, *Zur Struktur $AlCl_3$*, Z. Anorg. Allg. Chem. **193**, 270–276 (1930), doi:10.1002/zaac.19301930123.
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Found in:

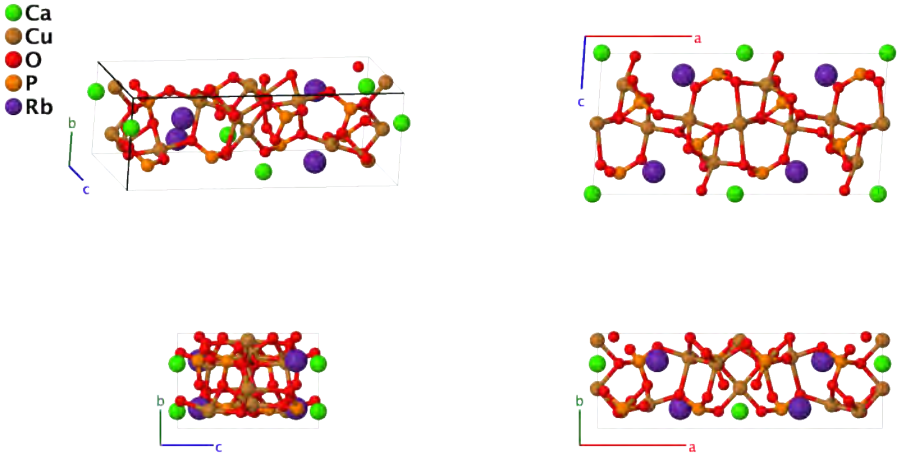
- C. Gottfried and F. Schossberger, eds., *Strukturbericht Band III 1933-1935* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1510
- POSCAR: pp. 1510

$Rb_2CaCu_6(PO_4)_4O_2$ Structure: AB6C18D4E2_mC62_5_a_2b2c_9c_2c_c

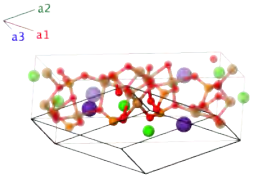
http://aflow.org/prototype-encyclopedia/AB6C18D4E2_mC62_5_a_2b2c_9c_2c_c



Prototype	:	$CaCu_6O_{18}P_4Rb_2$
AFLOW prototype label	:	AB6C18D4E2_mC62_5_a_2b2c_9c_2c_c
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC62
Space group number	:	5
Space group symbol	:	$C2$
AFLOW prototype command	:	`aflow --proto=AB6C18D4E2_mC62_5_a_2b2c_9c_2c_c --params=`$a, b/a, c/a, \beta, y_1, y_2, y_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}, x_{17}, y_{17}, z_{17}$

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2$	=	$y_1 b\,\hat{\mathbf{y}}$	$(2a)$	Ca
$\mathbf{B}_2$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	$(2b)$	Cu I
$\mathbf{B}_3$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	$(2b)$	Cu II
$\mathbf{B}_4$	=	$(x_4 - y_4)\,\mathbf{a}_1 + (x_4 + y_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(4c)$	Cu III

$\mathbf{B}_5$	$(-x_4 - y_4)\ \mathbf{a}_1 + (-x_4 + y_4)\ \mathbf{a}_2 - z_4\ \mathbf{a}_3$	$=$	$(-x_4 a - z_4 c \cos\beta)\ \hat{\mathbf{x}} + y_4 b\ \hat{\mathbf{y}} - z_4 c \sin\beta\ \hat{\mathbf{z}}$	$(4c)$	Cu III
$\mathbf{B}_6$	$(x_5 - y_5)\ \mathbf{a}_1 + (x_5 + y_5)\ \mathbf{a}_2 + z_5\ \mathbf{a}_3$	$=$	$(x_5 a + z_5 c \cos\beta)\ \hat{\mathbf{x}} + y_5 b\ \hat{\mathbf{y}} + z_5 c \sin\beta\ \hat{\mathbf{z}}$	$(4c)$	Cu IV
$\mathbf{B}_7$	$(-x_5 - y_5)\ \mathbf{a}_1 + (-x_5 + y_5)\ \mathbf{a}_2 - z_5\ \mathbf{a}_3$	$=$	$(-x_5 a - z_5 c \cos\beta)\ \hat{\mathbf{x}} + y_5 b\ \hat{\mathbf{y}} - z_5 c \sin\beta\ \hat{\mathbf{z}}$	$(4c)$	Cu IV
$\mathbf{B}_8$	$(x_6 - y_6)\ \mathbf{a}_1 + (x_6 + y_6)\ \mathbf{a}_2 + z_6\ \mathbf{a}_3$	$=$	$(x_6 a + z_6 c \cos\beta)\ \hat{\mathbf{x}} + y_6 b\ \hat{\mathbf{y}} + z_6 c \sin\beta\ \hat{\mathbf{z}}$	$(4c)$	O I
$\mathbf{B}_9$	$(-x_6 - y_6)\ \mathbf{a}_1 + (-x_6 + y_6)\ \mathbf{a}_2 - z_6\ \mathbf{a}_3$	$=$	$(-x_6 a - z_6 c \cos\beta)\ \hat{\mathbf{x}} + y_6 b\ \hat{\mathbf{y}} - z_6 c \sin\beta\ \hat{\mathbf{z}}$	$(4c)$	O I
$\mathbf{B}_{10}$	$(x_7 - y_7)\ \mathbf{a}_1 + (x_7 + y_7)\ \mathbf{a}_2 + z_7\ \mathbf{a}_3$	$=$	$(x_7 a + z_7 c \cos\beta)\ \hat{\mathbf{x}} + y_7 b\ \hat{\mathbf{y}} + z_7 c \sin\beta\ \hat{\mathbf{z}}$	$(4c)$	O II
$\mathbf{B}_{11}$	$(-x_7 - y_7)\ \mathbf{a}_1 + (-x_7 + y_7)\ \mathbf{a}_2 - z_7\ \mathbf{a}_3$	$=$	$(-x_7 a - z_7 c \cos\beta)\ \hat{\mathbf{x}} + y_7 b\ \hat{\mathbf{y}} - z_7 c \sin\beta\ \hat{\mathbf{z}}$	$(4c)$	O II
$\mathbf{B}_{12}$	$(x_8 - y_8)\ \mathbf{a}_1 + (x_8 + y_8)\ \mathbf{a}_2 + z_8\ \mathbf{a}_3$	$=$	$(x_8 a + z_8 c \cos\beta)\ \hat{\mathbf{x}} + y_8 b\ \hat{\mathbf{y}} + z_8 c \sin\beta\ \hat{\mathbf{z}}$	$(4c)$	O III
$\mathbf{B}_{13}$	$(-x_8 - y_8)\ \mathbf{a}_1 + (-x_8 + y_8)\ \mathbf{a}_2 - z_8\ \mathbf{a}_3$	$=$	$(-x_8 a - z_8 c \cos\beta)\ \hat{\mathbf{x}} + y_8 b\ \hat{\mathbf{y}} - z_8 c \sin\beta\ \hat{\mathbf{z}}$	$(4c)$	O III
$\mathbf{B}_{14}$	$(x_9 - y_9)\ \mathbf{a}_1 + (x_9 + y_9)\ \mathbf{a}_2 + z_9\ \mathbf{a}_3$	$=$	$(x_9 a + z_9 c \cos\beta)\ \hat{\mathbf{x}} + y_9 b\ \hat{\mathbf{y}} + z_9 c \sin\beta\ \hat{\mathbf{z}}$	$(4c)$	O IV
$\mathbf{B}_{15}$	$(-x_9 - y_9)\ \mathbf{a}_1 + (-x_9 + y_9)\ \mathbf{a}_2 - z_9\ \mathbf{a}_3$	$=$	$(-x_9 a - z_9 c \cos\beta)\ \hat{\mathbf{x}} + y_9 b\ \hat{\mathbf{y}} - z_9 c \sin\beta\ \hat{\mathbf{z}}$	$(4c)$	O IV
$\mathbf{B}_{16}$	$(x_{10} - y_{10})\ \mathbf{a}_1 + (x_{10} + y_{10})\ \mathbf{a}_2 + z_{10}\ \mathbf{a}_3$	$=$	$(x_{10} a + z_{10} c \cos\beta)\ \hat{\mathbf{x}} + y_{10} b\ \hat{\mathbf{y}} + z_{10} c \sin\beta\ \hat{\mathbf{z}}$	$(4c)$	O V
$\mathbf{B}_{17}$	$(-x_{10} - y_{10})\ \mathbf{a}_1 + (-x_{10} + y_{10})\ \mathbf{a}_2 - z_{10}\ \mathbf{a}_3$	$=$	$(-x_{10} a - z_{10} c \cos\beta)\ \hat{\mathbf{x}} + y_{10} b\ \hat{\mathbf{y}} - z_{10} c \sin\beta\ \hat{\mathbf{z}}$	$(4c)$	O V
$\mathbf{B}_{18}$	$(x_{11} - y_{11})\ \mathbf{a}_1 + (x_{11} + y_{11})\ \mathbf{a}_2 + z_{11}\ \mathbf{a}_3$	$=$	$(x_{11} a + z_{11} c \cos\beta)\ \hat{\mathbf{x}} + y_{11} b\ \hat{\mathbf{y}} + z_{11} c \sin\beta\ \hat{\mathbf{z}}$	$(4c)$	O VI
$\mathbf{B}_{19}$	$(-x_{11} - y_{11})\ \mathbf{a}_1 + (-x_{11} + y_{11})\ \mathbf{a}_2 - z_{11}\ \mathbf{a}_3$	$=$	$(-x_{11} a - z_{11} c \cos\beta)\ \hat{\mathbf{x}} + y_{11} b\ \hat{\mathbf{y}} - z_{11} c \sin\beta\ \hat{\mathbf{z}}$	$(4c)$	O VI
$\mathbf{B}_{20}$	$(x_{12} - y_{12})\ \mathbf{a}_1 + (x_{12} + y_{12})\ \mathbf{a}_2 + z_{12}\ \mathbf{a}_3$	$=$	$(x_{12} a + z_{12} c \cos\beta)\ \hat{\mathbf{x}} + y_{12} b\ \hat{\mathbf{y}} + z_{12} c \sin\beta\ \hat{\mathbf{z}}$	$(4c)$	O VII
$\mathbf{B}_{21}$	$(-x_{12} - y_{12})\ \mathbf{a}_1 + (-x_{12} + y_{12})\ \mathbf{a}_2 - z_{12}\ \mathbf{a}_3$	$=$	$(-x_{12} a - z_{12} c \cos\beta)\ \hat{\mathbf{x}} + y_{12} b\ \hat{\mathbf{y}} - z_{12} c \sin\beta\ \hat{\mathbf{z}}$	$(4c)$	O VII
$\mathbf{B}_{22}$	$(x_{13} - y_{13})\ \mathbf{a}_1 + (x_{13} + y_{13})\ \mathbf{a}_2 + z_{13}\ \mathbf{a}_3$	$=$	$(x_{13} a + z_{13} c \cos\beta)\ \hat{\mathbf{x}} + y_{13} b\ \hat{\mathbf{y}} + z_{13} c \sin\beta\ \hat{\mathbf{z}}$	$(4c)$	O VIII
$\mathbf{B}_{23}$	$(-x_{13} - y_{13})\ \mathbf{a}_1 + (-x_{13} + y_{13})\ \mathbf{a}_2 - z_{13}\ \mathbf{a}_3$	$=$	$(-x_{13} a - z_{13} c \cos\beta)\ \hat{\mathbf{x}} + y_{13} b\ \hat{\mathbf{y}} - z_{13} c \sin\beta\ \hat{\mathbf{z}}$	$(4c)$	O VIII
$\mathbf{B}_{24}$	$(x_{14} - y_{14})\ \mathbf{a}_1 + (x_{14} + y_{14})\ \mathbf{a}_2 + z_{14}\ \mathbf{a}_3$	$=$	$(x_{14} a + z_{14} c \cos\beta)\ \hat{\mathbf{x}} + y_{14} b\ \hat{\mathbf{y}} + z_{14} c \sin\beta\ \hat{\mathbf{z}}$	$(4c)$	O IX
$\mathbf{B}_{25}$	$(-x_{14} - y_{14})\ \mathbf{a}_1 + (-x_{14} + y_{14})\ \mathbf{a}_2 - z_{14}\ \mathbf{a}_3$	$=$	$(-x_{14} a - z_{14} c \cos\beta)\ \hat{\mathbf{x}} + y_{14} b\ \hat{\mathbf{y}} - z_{14} c \sin\beta\ \hat{\mathbf{z}}$	$(4c)$	O IX
$\mathbf{B}_{26}$	$(x_{15} - y_{15})\ \mathbf{a}_1 + (x_{15} + y_{15})\ \mathbf{a}_2 + z_{15}\ \mathbf{a}_3$	$=$	$(x_{15} a + z_{15} c \cos\beta)\ \hat{\mathbf{x}} + y_{15} b\ \hat{\mathbf{y}} + z_{15} c \sin\beta\ \hat{\mathbf{z}}$	$(4c)$	P I
$\mathbf{B}_{27}$	$(-x_{15} - y_{15})\ \mathbf{a}_1 + (-x_{15} + y_{15})\ \mathbf{a}_2 - z_{15}\ \mathbf{a}_3$	$=$	$(-x_{15} a - z_{15} c \cos\beta)\ \hat{\mathbf{x}} + y_{15} b\ \hat{\mathbf{y}} - z_{15} c \sin\beta\ \hat{\mathbf{z}}$	$(4c)$	P I
$\mathbf{B}_{28}$	$(x_{16} - y_{16})\ \mathbf{a}_1 + (x_{16} + y_{16})\ \mathbf{a}_2 + z_{16}\ \mathbf{a}_3$	$=$	$(x_{16} a + z_{16} c \cos\beta)\ \hat{\mathbf{x}} + y_{16} b\ \hat{\mathbf{y}} + z_{16} c \sin\beta\ \hat{\mathbf{z}}$	$(4c)$	P II

$\mathbf{B}_{29}$	=	$(-x_{16} - y_{16})\ \mathbf{a}_1 + (-x_{16} + y_{16})\ \mathbf{a}_2 - z_{16}\ \mathbf{a}_3$	=	$(-x_{16}a - z_{16}c\cos\beta)\ \hat{\mathbf{x}} + y_{16}b\ \hat{\mathbf{y}} - z_{16}c\sin\beta\ \hat{\mathbf{z}}$	(4*c*)	P II
$\mathbf{B}_{30}$	=	$(x_{17} - y_{17})\ \mathbf{a}_1 + (x_{17} + y_{17})\ \mathbf{a}_2 + z_{17}\ \mathbf{a}_3$	=	$(x_{17}a + z_{17}c\cos\beta)\ \hat{\mathbf{x}} + y_{17}b\ \hat{\mathbf{y}} + z_{17}c\sin\beta\ \hat{\mathbf{z}}$	(4*c*)	Rb
$\mathbf{B}_{31}$	=	$(-x_{17} - y_{17})\ \mathbf{a}_1 + (-x_{17} + y_{17})\ \mathbf{a}_2 - z_{17}\ \mathbf{a}_3$	=	$(-x_{17}a - z_{17}c\cos\beta)\ \hat{\mathbf{x}} + y_{17}b\ \hat{\mathbf{y}} - z_{17}c\sin\beta\ \hat{\mathbf{z}}$	(4*c*)	Rb

References:

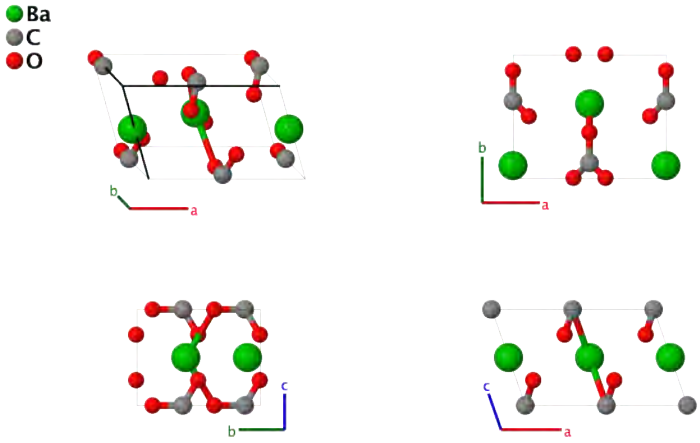
- S. M. Aksenov, E. Y. Borovikova, V. S. Mironov, N. A. Yamnova, A. S. Volkov, D. A. Ksenofontov, O. A. Gurbanova, O. V. Dimitrova, D. V. Deyneko, E. A. Zvereva, O. V. Maximova, S. V. Krivovichev, P. C. Burns, and A. N. Vasiliev, *$Rb_2CaCu_6(PO_4)_4O_2$, a novel oxophosphate with a shchurovskyite-type topology: synthesis, structure, magnetic properties and crystal chemistry of rubidium copper phosphates*, Acta Crystallogr. Sect. B Struct. Sci. **75**, 903–913 (2019), doi:10.1107/S2052520619008527.

Geometry files:

- CIF: pp. 1511
- POSCAR: pp. 1511

$C2$ (Ba,Ca)CO_3 Structure: ABC3_mC10_5_b_a_ac

http://aflow.org/prototype-encyclopedia/ABC3_mC10_5_b_a_ac

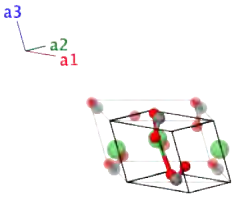


Prototype	:	(Ba,Ca)CO_3
AFLOW prototype label	:	ABC3_mC10_5_b_a_ac
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC10
Space group number	:	5
Space group symbol	:	$C2$
AFLOW prototype command	:	`aflow --proto=ABC3_mC10_5_b_a_ac --params=`$a, b/a, c/a, \beta, y_1, y_2, y_3, x_4, y_4, z_4$

- (Ba,Ca)$(CO_3)_2$ comes in a variety of crystal structures (Spahr, 2019):
  - monoclinic barytocalcite, space group $P2_1/m$ #11 (the current structure).
  - trigonal paralstonite, space group $P321$ #150,
  - Triclinic alstonite, space group $P1$ #1 or $P\bar{1}$ #2 (Sartori, 1975), and
  - a new monoclinic structure, space group $C2$ #5, synthesized by (Spahr, 2019), and lacking the centrosymmetric character of barytocalcite.
- The site we have labeled Ba is actually a mixture of barium and calcium atoms.

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$\mathbf{B}_1$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2$	=	$y_1 b\,\hat{\mathbf{y}}$	$(2a)$	C
$\mathbf{B}_2$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2$	=	$y_2 b\,\hat{\mathbf{y}}$	$(2a)$	O I
$\mathbf{B}_3$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	$(2b)$	Ba
$\mathbf{B}_4$	=	$(x_4 - y_4)\,\mathbf{a}_1 + (x_4 + y_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(4c)$	O II
$\mathbf{B}_5$	=	$(-x_4 - y_4)\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(4c)$	O II

References:

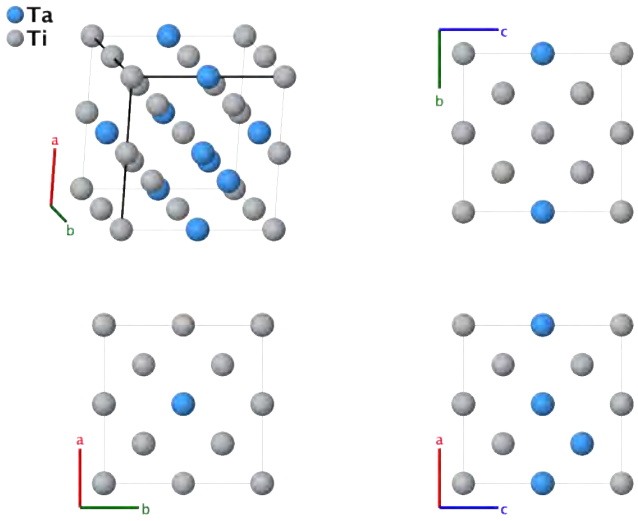
- D. Spahr, L. Bayarjargal, V. Vinograd, R. Luchitskaia, V. Milman, and B. Winkler, *A new $BaCa(CO_3)_2$ polymorph*, Acta Crystallogr. Sect. B Struct. Sci. **75**, 291–300 (2019), doi:10.1107/S2052520619003238.
- F. Sartori, *New data on alstonite*, Lithos **8**, 199–207 (1975), doi:10.1016/0024-4937(75)90036-5.

Geometry files:

- CIF: pp. 1511
- POSCAR: pp. 1512

Ta_5Ti_{11} (BCC SQS-16) Structure: A5B11_mP16_6_2abc_2a3b3c

http://aflow.org/prototype-encyclopedia/A5B11_mP16_6_2abc_2a3b3c

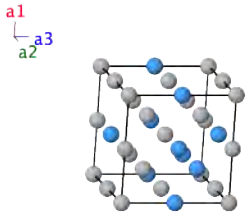


Prototype	:	Ta_5Ti_{11}
AFLOW prototype label	:	A5B11_mP16_6_2abc_2a3b3c
***Strukturbericht* designation**	:	None
Pearson symbol	:	mP16
Space group number	:	6
Space group symbol	:	*Pm*
AFLOW prototype command	:	`aflow --proto=A5B11_mP16_6_2abc_2a3b3c --params=`$a, b/a, c/a, \beta, x_1, z_1, x_2, z_2, x_3, z_3, x_4, z_4, x_5, z_5, x_6, z_6, x_7, z_7, x_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}$

- This is a special quasirandom structure with 16 atoms per unit cell (SQS-16) for the β-phase (high-temperature austenite) bcc substitutional Ti-Ta alloy (Chakraborty, 2016). This protoype contains 31.25% Ta. Prototypes are listed for other Ta-Ti concentrations: 12.5% Ta (AB7_hR16_166_c_c2h), 18.75% Ta (A3B13_oC32_38_ac_a2bcdef), 25% Ta (AB3_mC32_8_4a_4a4b), and 37.5% Ta (A3B5_oC32_38_abce_abcdf).

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + z_1\,\mathbf{a}_3$	=	$(x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(1a)$	Ta I
$\mathbf{B}_2$	=	$x_2\,\mathbf{a}_1 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(1a)$	Ta II
$\mathbf{B}_3$	=	$x_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(1a)$	Ti I
$\mathbf{B}_4$	=	$x_4\,\mathbf{a}_1 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(1a)$	Ti II
$\mathbf{B}_5$	=	$x_5\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{2} b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(1b)$	Ta III
$\mathbf{B}_6$	=	$x_6\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{2} b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(1b)$	Ti III
$\mathbf{B}_7$	=	$x_7\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{2} b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(1b)$	Ti IV
$\mathbf{B}_8$	=	$x_8\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{2} b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(1b)$	Ti V
$\mathbf{B}_9$	=	$x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$(x_9 a + z_9 c\cos\beta)\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\sin\beta\,\hat{\mathbf{z}}$	$(2c)$	Ta IV
$\mathbf{B}_{10}$	=	$x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$(x_9 a + z_9 c\cos\beta)\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + z_9 c\sin\beta\,\hat{\mathbf{z}}$	$(2c)$	Ta IV
$\mathbf{B}_{11}$	=	$x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$(x_{10} a + z_{10} c\cos\beta)\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\sin\beta\,\hat{\mathbf{z}}$	$(2c)$	Ti VI
$\mathbf{B}_{12}$	=	$x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$(x_{10} a + z_{10} c\cos\beta)\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} + z_{10} c\sin\beta\,\hat{\mathbf{z}}$	$(2c)$	Ti VI
$\mathbf{B}_{13}$	=	$x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$(x_{11} a + z_{11} c\cos\beta)\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + z_{11} c\sin\beta\,\hat{\mathbf{z}}$	$(2c)$	Ti VII
$\mathbf{B}_{14}$	=	$x_{11}\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$(x_{11} a + z_{11} c\cos\beta)\,\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} + z_{11} c\sin\beta\,\hat{\mathbf{z}}$	$(2c)$	Ti VII
$\mathbf{B}_{15}$	=	$x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3$	=	$(x_{12} a + z_{12} c\cos\beta)\,\hat{\mathbf{x}} + y_{12} b\,\hat{\mathbf{y}} + z_{12} c\sin\beta\,\hat{\mathbf{z}}$	$(2c)$	Ti VIII
$\mathbf{B}_{16}$	=	$x_{12}\,\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3$	=	$(x_{12} a + z_{12} c\cos\beta)\,\hat{\mathbf{x}} - y_{12} b\,\hat{\mathbf{y}} + z_{12} c\sin\beta\,\hat{\mathbf{z}}$	$(2c)$	Ti VIII

References:

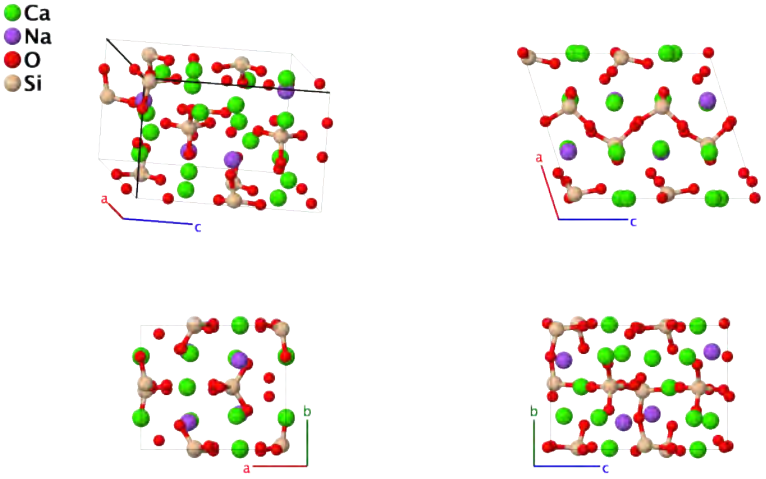
- T. Chakraborty, J. Rogal, and R. Drautz, *Unraveling the composition dependence of the martensitic transformation temperature: A first-principles study of Ti-Ta alloys*, Phys. Rev. B **94**, 224104 (2016), doi:10.1103/PhysRevB.94.224104.

Geometry files:

- CIF: pp. 1512
- POSCAR: pp. 1512

$Na_2Ca_6Si_4O_{15}$ Structure: A6B2C15D4_mP54_7_6a_2a_15a_4a

http://aflow.org/prototype-encyclopedia/A6B2C15D4_mP54_7_6a_2a_15a_4a

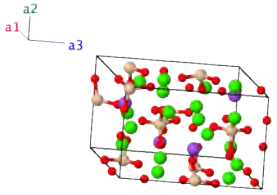


Prototype	:	$Ca_6Na_2O_{15}Si_4$
AFLOW prototype label	:	A6B2C15D4_mP54_7_6a_2a_15a_4a
***Strukturbericht* designation**	:	None
Pearson symbol	:	mP54
Space group number	:	7
Space group symbol	:	*Pc*
AFLOW prototype command	:	`aflow --proto=A6B2C15D4_mP54_7_6a_2a_15a_4a --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}, x_{17}, y_{17}, z_{17}, x_{18}, y_{18}, z_{18}, x_{19}, y_{19}, z_{19}, x_{20}, y_{20}, z_{20}, x_{21}, y_{21}, z_{21}, x_{22}, y_{22}, z_{22}, x_{23}, y_{23}, z_{23}, x_{24}, y_{24}, z_{24}, x_{25}, y_{25}, z_{25}, x_{26}, y_{26}, z_{26}, x_{27}, y_{27}, z_{27}$

- Although this structure was formed at 1300 °C, the data was taken after cooling to room temperature, 295 K.
- The true composition of the Ca-I site is $Ca_{0.84}Na_{0.16}$, and the Na-I site has composition $Na_{0.84}Ca_{0.16}$.
- The Si-III and Si-IV atoms share the O-*X* atom between them, accounting for the "missing" oxygen atom.

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1a + z_1c\cos\beta)\,\hat{\mathbf{x}} + y_1b\,\hat{\mathbf{y}} + z_1c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Ca I
$\mathbf{B}_2$	=	$x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_1a + z_1c\cos\beta\right)\hat{\mathbf{x}} - y_1b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Ca I
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2a + z_2c\cos\beta)\,\hat{\mathbf{x}} + y_2b\,\hat{\mathbf{y}} + z_2c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Ca II
$\mathbf{B}_4$	=	$x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_2a + z_2c\cos\beta\right)\hat{\mathbf{x}} - y_2b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Ca II
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3a + z_3c\cos\beta)\,\hat{\mathbf{x}} + y_3b\,\hat{\mathbf{y}} + z_3c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Ca III
$\mathbf{B}_6$	=	$x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_3a + z_3c\cos\beta\right)\hat{\mathbf{x}} - y_3b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Ca III
$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4a + z_4c\cos\beta)\,\hat{\mathbf{x}} + y_4b\,\hat{\mathbf{y}} + z_4c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Ca IV
$\mathbf{B}_8$	=	$x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_4a + z_4c\cos\beta\right)\hat{\mathbf{x}} - y_4b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Ca IV
$\mathbf{B}_9$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5a + z_5c\cos\beta)\,\hat{\mathbf{x}} + y_5b\,\hat{\mathbf{y}} + z_5c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Ca V
$\mathbf{B}_{10}$	=	$x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_5a + z_5c\cos\beta\right)\hat{\mathbf{x}} - y_5b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Ca V
$\mathbf{B}_{11}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6a + z_6c\cos\beta)\,\hat{\mathbf{x}} + y_6b\,\hat{\mathbf{y}} + z_6c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Ca VI
$\mathbf{B}_{12}$	=	$x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_6a + z_6c\cos\beta\right)\hat{\mathbf{x}} - y_6b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Ca VI
$\mathbf{B}_{13}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7a + z_7c\cos\beta)\,\hat{\mathbf{x}} + y_7b\,\hat{\mathbf{y}} + z_7c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Na I
$\mathbf{B}_{14}$	=	$x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_7a + z_7c\cos\beta\right)\hat{\mathbf{x}} - y_7b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right)c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Na I
$\mathbf{B}_{15}$	=	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8a + z_8c\cos\beta)\,\hat{\mathbf{x}} + y_8b\,\hat{\mathbf{y}} + z_8c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Na II
$\mathbf{B}_{16}$	=	$x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_8a + z_8c\cos\beta\right)\hat{\mathbf{x}} - y_8b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right)c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Na II
$\mathbf{B}_{17}$	=	$x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$(x_9a + z_9c\cos\beta)\,\hat{\mathbf{x}} + y_9b\,\hat{\mathbf{y}} + z_9c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	O I
$\mathbf{B}_{18}$	=	$x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_9a + z_9c\cos\beta\right)\hat{\mathbf{x}} - y_9b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right)c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	O I
$\mathbf{B}_{19}$	=	$x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$(x_{10}a + z_{10}c\cos\beta)\,\hat{\mathbf{x}} + y_{10}b\,\hat{\mathbf{y}} + z_{10}c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	O II
$\mathbf{B}_{20}$	=	$x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{10}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{10}a + z_{10}c\cos\beta\right)\hat{\mathbf{x}} - y_{10}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{10}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	O II
$\mathbf{B}_{21}$	=	$x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$(x_{11}a + z_{11}c\cos\beta)\,\hat{\mathbf{x}} + y_{11}b\,\hat{\mathbf{y}} + z_{11}c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	O III
$\mathbf{B}_{22}$	=	$x_{11}\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{11}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{11}a + z_{11}c\cos\beta\right)\hat{\mathbf{x}} - y_{11}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{11}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	O III
$\mathbf{B}_{23}$	=	$x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3$	=	$(x_{12}a + z_{12}c\cos\beta)\,\hat{\mathbf{x}} + y_{12}b\,\hat{\mathbf{y}} + z_{12}c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	O IV
$\mathbf{B}_{24}$	=	$x_{12}\,\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{12}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{12}a + z_{12}c\cos\beta\right)\hat{\mathbf{x}} - y_{12}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{12}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	O IV

$$\mathbf{B}_{25} = x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3 = (x_{13}a + z_{13}c\cos\beta)\,\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} + z_{13}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O V}$$

$$\mathbf{B}_{26} = x_{13}\,\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{13}a + z_{13}c\cos\beta\right)\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{13}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O V}$$

$$\mathbf{B}_{27} = x_{14}\,\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3 = (x_{14}a + z_{14}c\cos\beta)\,\hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} + z_{14}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O VI}$$

$$\mathbf{B}_{28} = x_{14}\,\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{14}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{14}a + z_{14}c\cos\beta\right)\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{14}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O VI}$$

$$\mathbf{B}_{29} = x_{15}\,\mathbf{a}_1 + y_{15}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 = (x_{15}a + z_{15}c\cos\beta)\,\hat{\mathbf{x}} + y_{15}b\,\hat{\mathbf{y}} + z_{15}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O VII}$$

$$\mathbf{B}_{30} = x_{15}\,\mathbf{a}_1 - y_{15}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{15}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{15}a + z_{15}c\cos\beta\right)\hat{\mathbf{x}} - y_{15}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{15}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O VII}$$

$$\mathbf{B}_{31} = x_{16}\,\mathbf{a}_1 + y_{16}\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 = (x_{16}a + z_{16}c\cos\beta)\,\hat{\mathbf{x}} + y_{16}b\,\hat{\mathbf{y}} + z_{16}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O VIII}$$

$$\mathbf{B}_{32} = x_{16}\,\mathbf{a}_1 - y_{16}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{16}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{16}a + z_{16}c\cos\beta\right)\hat{\mathbf{x}} - y_{16}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{16}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O VIII}$$

$$\mathbf{B}_{33} = x_{17}\,\mathbf{a}_1 + y_{17}\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 = (x_{17}a + z_{17}c\cos\beta)\,\hat{\mathbf{x}} + y_{17}b\,\hat{\mathbf{y}} + z_{17}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O IX}$$

$$\mathbf{B}_{34} = x_{17}\,\mathbf{a}_1 - y_{17}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{17}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{17}a + z_{17}c\cos\beta\right)\hat{\mathbf{x}} - y_{17}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{17}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O IX}$$

$$\mathbf{B}_{35} = x_{18}\,\mathbf{a}_1 + y_{18}\,\mathbf{a}_2 + z_{18}\,\mathbf{a}_3 = (x_{18}a + z_{18}c\cos\beta)\,\hat{\mathbf{x}} + y_{18}b\,\hat{\mathbf{y}} + z_{18}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O X}$$

$$\mathbf{B}_{36} = x_{18}\,\mathbf{a}_1 - y_{18}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{18}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{18}a + z_{18}c\cos\beta\right)\hat{\mathbf{x}} - y_{18}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{18}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O X}$$

$$\mathbf{B}_{37} = x_{19}\,\mathbf{a}_1 + y_{19}\,\mathbf{a}_2 + z_{19}\,\mathbf{a}_3 = (x_{19}a + z_{19}c\cos\beta)\,\hat{\mathbf{x}} + y_{19}b\,\hat{\mathbf{y}} + z_{19}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XI}$$

$$\mathbf{B}_{38} = x_{19}\,\mathbf{a}_1 - y_{19}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{19}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{19}a + z_{19}c\cos\beta\right)\hat{\mathbf{x}} - y_{19}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{19}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XI}$$

$$\mathbf{B}_{39} = x_{20}\,\mathbf{a}_1 + y_{20}\,\mathbf{a}_2 + z_{20}\,\mathbf{a}_3 = (x_{20}a + z_{20}c\cos\beta)\,\hat{\mathbf{x}} + y_{20}b\,\hat{\mathbf{y}} + z_{20}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XII}$$

$$\mathbf{B}_{40} = x_{20}\,\mathbf{a}_1 - y_{20}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{20}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{20}a + z_{20}c\cos\beta\right)\hat{\mathbf{x}} - y_{20}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{20}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XII}$$

$$\mathbf{B}_{41} = x_{21}\,\mathbf{a}_1 + y_{21}\,\mathbf{a}_2 + z_{21}\,\mathbf{a}_3 = (x_{21}a + z_{21}c\cos\beta)\,\hat{\mathbf{x}} + y_{21}b\,\hat{\mathbf{y}} + z_{21}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XIII}$$

$$\mathbf{B}_{42} = x_{21}\,\mathbf{a}_1 - y_{21}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{21}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{21}a + z_{21}c\cos\beta\right)\hat{\mathbf{x}} - y_{21}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{21}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XIII}$$

$$\mathbf{B}_{43} = x_{22}\,\mathbf{a}_1 + y_{22}\,\mathbf{a}_2 + z_{22}\,\mathbf{a}_3 = (x_{22}a + z_{22}c\cos\beta)\,\hat{\mathbf{x}} + y_{22}b\,\hat{\mathbf{y}} + z_{22}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XIV}$$

$$\mathbf{B}_{44} = x_{22}\,\mathbf{a}_1 - y_{22}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{22}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{22}a + z_{22}c\cos\beta\right)\hat{\mathbf{x}} - y_{22}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{22}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XIV}$$

$$\mathbf{B}_{45} = x_{23}\,\mathbf{a}_1 + y_{23}\,\mathbf{a}_2 + z_{23}\,\mathbf{a}_3 = (x_{23}a + z_{23}c\cos\beta)\,\hat{\mathbf{x}} + y_{23}b\,\hat{\mathbf{y}} + z_{23}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XV}$$

$$\mathbf{B}_{46} = x_{23}\,\mathbf{a}_1 - y_{23}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{23}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{23}a + z_{23}c\cos\beta\right)\hat{\mathbf{x}} - y_{23}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{23}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XV}$$

$$\mathbf{B}_{47} = x_{24}\,\mathbf{a}_1 + y_{24}\,\mathbf{a}_2 + z_{24}\,\mathbf{a}_3 = (x_{24}a + z_{24}c\cos\beta)\,\hat{\mathbf{x}} + y_{24}b\,\hat{\mathbf{y}} + z_{24}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{Si I}$$

$$\mathbf{B}_{48} = x_{24}\,\mathbf{a}_1 - y_{24}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{24}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{24}a + z_{24}c\cos\beta\right)\hat{\mathbf{x}} - y_{24}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{24}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{Si I}$$

$$\mathbf{B}_{49} = x_{25}\,\mathbf{a}_1 + y_{25}\,\mathbf{a}_2 + z_{25}\,\mathbf{a}_3 = (x_{25}a + z_{25}c\cos\beta)\,\hat{\mathbf{x}} + y_{25}b\,\hat{\mathbf{y}} + z_{25}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{Si II}$$

$$\mathbf{B}_{50} = x_{25}\,\mathbf{a}_1 - y_{25}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{25}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{25}a + z_{25}c\cos\beta\right)\hat{\mathbf{x}} - y_{25}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{25}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{Si II}$$

$$\mathbf{B}_{51} = x_{26}\,\mathbf{a}_1 + y_{26}\,\mathbf{a}_2 + z_{26}\,\mathbf{a}_3 = (x_{26}a + z_{26}c\cos\beta)\,\hat{\mathbf{x}} + y_{26}b\,\hat{\mathbf{y}} + z_{26}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{Si III}$$

$$\mathbf{B}_{52} = x_{26}\,\mathbf{a}_1 - y_{26}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{26}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{26}a + z_{26}c\cos\beta\right)\hat{\mathbf{x}} - y_{26}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{26}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{Si III}$$

$$\mathbf{B}_{53} = x_{27}\,\mathbf{a}_1 + y_{27}\,\mathbf{a}_2 + z_{27}\,\mathbf{a}_3 = (x_{27}a + z_{27}c\cos\beta)\,\hat{\mathbf{x}} + y_{27}b\,\hat{\mathbf{y}} + z_{27}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{Si IV}$$

$$\mathbf{B}_{54} = x_{27}\,\mathbf{a}_1 - y_{27}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{27}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{27}a + z_{27}c\cos\beta\right)\hat{\mathbf{x}} - y_{27}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{27}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{Si IV}$$

References:

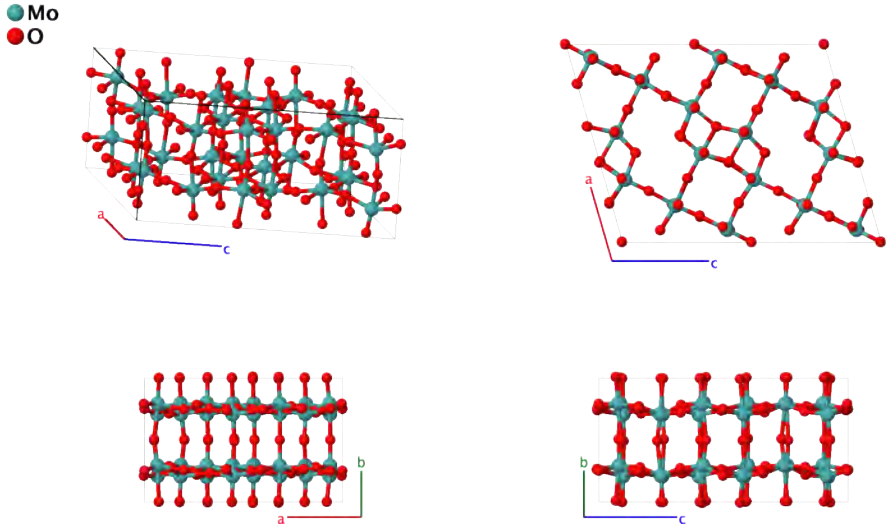
- V. Kahlenberg and M. Maier, *On the existence of a high-temperature polymorph of $Na_2Ca_6Si_4O_{15}$–implications for the phase equilibria in the system Na_2O-CaO-SiO_2*, Mineral. Petrol. **110**, 905–915 (2016), doi:10.1007/s00710-016-0447-1.

Geometry files:

- CIF: pp. 1512
- POSCAR: pp. 1513

Low-Temperature Mo_8O_{23} Structure: A8B23_mP124_7_16a_46a

http://aflow.org/prototype-encyclopedia/A8B23_mP124_7_16a_46a

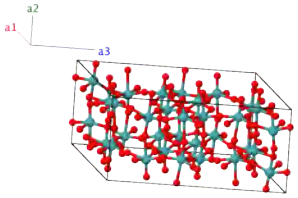


Prototype	:	Mo_8O_{23}
AFLOW prototype label	:	A8B23_mP124_7_16a_46a
***Strukturbericht* designation**	:	None
Pearson symbol	:	mP124
Space group number	:	7
Space group symbol	:	*Pc*
AFLOW prototype command	:	`aflow --proto=A8B23_mP124_7_16a_46a --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}, x_{17}, y_{17}, z_{17}, x_{18}, y_{18}, z_{18}, x_{19}, y_{19}, z_{19}, x_{20}, y_{20}, z_{20}, x_{21}, y_{21}, z_{21}, x_{22}, y_{22}, z_{22}, x_{23}, y_{23}, z_{23}, x_{24}, y_{24}, z_{24}, x_{25}, y_{25}, z_{25}, x_{26}, y_{26}, z_{26}, x_{27}, y_{27}, z_{27}, x_{28}, y_{28}, z_{28}, x_{29}, y_{29}, z_{29}, x_{30}, y_{30}, z_{30}, x_{31}, y_{31}, z_{31}, x_{32}, y_{32}, z_{32}, x_{33}, y_{33}, z_{33}, x_{34}, y_{34}, z_{34}, x_{35}, y_{35}, z_{35}, x_{36}, y_{36}, z_{36}, x_{37}, y_{37}, z_{37}, x_{38}, y_{38}, z_{38}, x_{39}, y_{39}, z_{39}, x_{40}, y_{40}, z_{40}, x_{41}, y_{41}, z_{41}, x_{42}, y_{42}, z_{42}, x_{43}, y_{43}, z_{43}, x_{44}, y_{44}, z_{44}, x_{45}, y_{45}, z_{45}, x_{46}, y_{46}, z_{46}, x_{47}, y_{47}, z_{47}, x_{48}, y_{48}, z_{48}, x_{49}, y_{49}, z_{49}, x_{50}, y_{50}, z_{50}, x_{51}, y_{51}, z_{51}, x_{52}, y_{52}, z_{52}, x_{53}, y_{53}, z_{53}, x_{54}, y_{54}, z_{54}, x_{55}, y_{55}, z_{55}, x_{56}, y_{56}, z_{56}, x_{57}, y_{57}, z_{57}, x_{58}, y_{58}, z_{58}, x_{59}, y_{59}, z_{59}, x_{60}, y_{60}, z_{60}, x_{61}, y_{61}, z_{61}, x_{62}, y_{62}, z_{62}$

- This data was taken at 100 K. Above 285 K the structure exhibits an incommensurate charge density wave, which is approximated by the high temperature Mo_8O_{23} structure.

Simple Monoclinic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= a\,\hat{\mathbf{x}}\\ \mathbf{a}_2 &= b\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	$(x_1a + z_1c\cos\beta)\,\hat{\mathbf{x}} + y_1b\,\hat{\mathbf{y}} + z_1c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Mo I
$\mathbf{B}_2$	$x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_1a + z_1c\cos\beta\right)\hat{\mathbf{x}} - y_1b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Mo I
$\mathbf{B}_3$	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	$(x_2a + z_2c\cos\beta)\,\hat{\mathbf{x}} + y_2b\,\hat{\mathbf{y}} + z_2c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Mo II
$\mathbf{B}_4$	$x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_2a + z_2c\cos\beta\right)\hat{\mathbf{x}} - y_2b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Mo II
$\mathbf{B}_5$	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$(x_3a + z_3c\cos\beta)\,\hat{\mathbf{x}} + y_3b\,\hat{\mathbf{y}} + z_3c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Mo III
$\mathbf{B}_6$	$x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_3a + z_3c\cos\beta\right)\hat{\mathbf{x}} - y_3b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Mo III
$\mathbf{B}_7$	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$(x_4a + z_4c\cos\beta)\,\hat{\mathbf{x}} + y_4b\,\hat{\mathbf{y}} + z_4c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Mo IV
$\mathbf{B}_8$	$x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_4a + z_4c\cos\beta\right)\hat{\mathbf{x}} - y_4b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Mo IV
$\mathbf{B}_9$	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	$(x_5a + z_5c\cos\beta)\,\hat{\mathbf{x}} + y_5b\,\hat{\mathbf{y}} + z_5c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Mo V
$\mathbf{B}_{10}$	$x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_5a + z_5c\cos\beta\right)\hat{\mathbf{x}} - y_5b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Mo V
$\mathbf{B}_{11}$	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	$(x_6a + z_6c\cos\beta)\,\hat{\mathbf{x}} + y_6b\,\hat{\mathbf{y}} + z_6c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Mo VI
$\mathbf{B}_{12}$	$x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_6a + z_6c\cos\beta\right)\hat{\mathbf{x}} - y_6b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Mo VI
$\mathbf{B}_{13}$	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	$(x_7a + z_7c\cos\beta)\,\hat{\mathbf{x}} + y_7b\,\hat{\mathbf{y}} + z_7c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Mo VII
$\mathbf{B}_{14}$	$x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_7a + z_7c\cos\beta\right)\hat{\mathbf{x}} - y_7b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right)c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Mo VII
$\mathbf{B}_{15}$	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	$(x_8a + z_8c\cos\beta)\,\hat{\mathbf{x}} + y_8b\,\hat{\mathbf{y}} + z_8c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Mo VIII
$\mathbf{B}_{16}$	$x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_8a + z_8c\cos\beta\right)\hat{\mathbf{x}} - y_8b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right)c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Mo VIII
$\mathbf{B}_{17}$	$x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	$(x_9a + z_9c\cos\beta)\,\hat{\mathbf{x}} + y_9b\,\hat{\mathbf{y}} + z_9c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Mo IX
$\mathbf{B}_{18}$	$x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_9a + z_9c\cos\beta\right)\hat{\mathbf{x}} - y_9b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right)c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Mo IX
$\mathbf{B}_{19}$	$x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	$(x_{10}a + z_{10}c\cos\beta)\,\hat{\mathbf{x}} + y_{10}b\,\hat{\mathbf{y}} + z_{10}c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Mo X
$\mathbf{B}_{20}$	$x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{10}\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_{10}a + z_{10}c\cos\beta\right)\hat{\mathbf{x}} - y_{10}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{10}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Mo X

$$\mathbf{B}_{21} = x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 = (x_{11}a + z_{11}c\cos\beta)\,\hat{\mathbf{x}} + y_{11}b\,\hat{\mathbf{y}} + z_{11}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{Mo XI}$$

$$\mathbf{B}_{22} = x_{11}\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{11}a + z_{11}c\cos\beta\right)\hat{\mathbf{x}} - y_{11}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{11}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{Mo XI}$$

$$\mathbf{B}_{23} = x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 = (x_{12}a + z_{12}c\cos\beta)\,\hat{\mathbf{x}} + y_{12}b\,\hat{\mathbf{y}} + z_{12}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{Mo XII}$$

$$\mathbf{B}_{24} = x_{12}\,\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{12}a + z_{12}c\cos\beta\right)\hat{\mathbf{x}} - y_{12}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{12}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{Mo XII}$$

$$\mathbf{B}_{25} = x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3 = (x_{13}a + z_{13}c\cos\beta)\,\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} + z_{13}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{Mo XIII}$$

$$\mathbf{B}_{26} = x_{13}\,\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{13}a + z_{13}c\cos\beta\right)\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{13}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{Mo XIII}$$

$$\mathbf{B}_{27} = x_{14}\,\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3 = (x_{14}a + z_{14}c\cos\beta)\,\hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} + z_{14}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{Mo XIV}$$

$$\mathbf{B}_{28} = x_{14}\,\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{14}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{14}a + z_{14}c\cos\beta\right)\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{14}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{Mo XIV}$$

$$\mathbf{B}_{29} = x_{15}\,\mathbf{a}_1 + y_{15}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 = (x_{15}a + z_{15}c\cos\beta)\,\hat{\mathbf{x}} + y_{15}b\,\hat{\mathbf{y}} + z_{15}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{Mo XV}$$

$$\mathbf{B}_{30} = x_{15}\,\mathbf{a}_1 - y_{15}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{15}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{15}a + z_{15}c\cos\beta\right)\hat{\mathbf{x}} - y_{15}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{15}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{Mo XV}$$

$$\mathbf{B}_{31} = x_{16}\,\mathbf{a}_1 + y_{16}\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 = (x_{16}a + z_{16}c\cos\beta)\,\hat{\mathbf{x}} + y_{16}b\,\hat{\mathbf{y}} + z_{16}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{Mo XVI}$$

$$\mathbf{B}_{32} = x_{16}\,\mathbf{a}_1 - y_{16}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{16}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{16}a + z_{16}c\cos\beta\right)\hat{\mathbf{x}} - y_{16}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{16}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{Mo XVI}$$

$$\mathbf{B}_{33} = x_{17}\,\mathbf{a}_1 + y_{17}\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 = (x_{17}a + z_{17}c\cos\beta)\,\hat{\mathbf{x}} + y_{17}b\,\hat{\mathbf{y}} + z_{17}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O I}$$

$$\mathbf{B}_{34} = x_{17}\,\mathbf{a}_1 - y_{17}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{17}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{17}a + z_{17}c\cos\beta\right)\hat{\mathbf{x}} - y_{17}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{17}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O I}$$

$$\mathbf{B}_{35} = x_{18}\,\mathbf{a}_1 + y_{18}\,\mathbf{a}_2 + z_{18}\,\mathbf{a}_3 = (x_{18}a + z_{18}c\cos\beta)\,\hat{\mathbf{x}} + y_{18}b\,\hat{\mathbf{y}} + z_{18}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O II}$$

$$\mathbf{B}_{36} = x_{18}\,\mathbf{a}_1 - y_{18}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{18}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{18}a + z_{18}c\cos\beta\right)\hat{\mathbf{x}} - y_{18}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{18}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O II}$$

$$\mathbf{B}_{37} = x_{19}\,\mathbf{a}_1 + y_{19}\,\mathbf{a}_2 + z_{19}\,\mathbf{a}_3 = (x_{19}a + z_{19}c\cos\beta)\,\hat{\mathbf{x}} + y_{19}b\,\hat{\mathbf{y}} + z_{19}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O III}$$

$$\mathbf{B}_{38} = x_{19}\,\mathbf{a}_1 - y_{19}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{19}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{19}a + z_{19}c\cos\beta\right)\hat{\mathbf{x}} - y_{19}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{19}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O III}$$

$$\mathbf{B}_{39} = x_{20}\,\mathbf{a}_1 + y_{20}\,\mathbf{a}_2 + z_{20}\,\mathbf{a}_3 = (x_{20}a + z_{20}c\cos\beta)\,\hat{\mathbf{x}} + y_{20}b\,\hat{\mathbf{y}} + z_{20}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O IV}$$

$$\mathbf{B}_{40} = x_{20}\,\mathbf{a}_1 - y_{20}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{20}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{20}a + z_{20}c\cos\beta\right)\hat{\mathbf{x}} - y_{20}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{20}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O IV}$$

$$\mathbf{B}_{41} = x_{21}\,\mathbf{a}_1 + y_{21}\,\mathbf{a}_2 + z_{21}\,\mathbf{a}_3 = (x_{21}a + z_{21}c\cos\beta)\,\hat{\mathbf{x}} + y_{21}b\,\hat{\mathbf{y}} + z_{21}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O V}$$

$$\mathbf{B}_{42} = x_{21}\,\mathbf{a}_1 - y_{21}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{21}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{21}a + z_{21}c\cos\beta\right)\hat{\mathbf{x}} - y_{21}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{21}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O V}$$

$$\mathbf{B}_{43} = x_{22}\,\mathbf{a}_1 + y_{22}\,\mathbf{a}_2 + z_{22}\,\mathbf{a}_3 = (x_{22}a + z_{22}c\cos\beta)\,\hat{\mathbf{x}} + y_{22}b\,\hat{\mathbf{y}} + z_{22}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O VI}$$

$$\mathbf{B}_{44} = x_{22}\,\mathbf{a}_1 - y_{22}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{22}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{22}a + z_{22}c\cos\beta\right)\hat{\mathbf{x}} - y_{22}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{22}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O VI}$$

$$\mathbf{B}_{45} = x_{23}\,\mathbf{a}_1 + y_{23}\,\mathbf{a}_2 + z_{23}\,\mathbf{a}_3 = (x_{23}a + z_{23}c\cos\beta)\,\hat{\mathbf{x}} + y_{23}b\,\hat{\mathbf{y}} + z_{23}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O VII}$$

$$\mathbf{B}_{46} = x_{23}\,\mathbf{a}_1 - y_{23}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{23}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{23}a + z_{23}c\cos\beta\right)\hat{\mathbf{x}} - y_{23}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{23}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O VII}$$

$$\mathbf{B}_{47} = x_{24}\,\mathbf{a}_1 + y_{24}\,\mathbf{a}_2 + z_{24}\,\mathbf{a}_3 = (x_{24}a + z_{24}c\cos\beta)\,\hat{\mathbf{x}} + y_{24}b\,\hat{\mathbf{y}} + z_{24}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O VIII}$$

$$\mathbf{B}_{48} = x_{24}\,\mathbf{a}_1 - y_{24}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{24}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{24}a + z_{24}c\cos\beta\right)\hat{\mathbf{x}} - y_{24}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{24}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O VIII}$$

$$\mathbf{B}_{49} = x_{25}\,\mathbf{a}_1 + y_{25}\,\mathbf{a}_2 + z_{25}\,\mathbf{a}_3 = (x_{25}a + z_{25}c\cos\beta)\,\hat{\mathbf{x}} + y_{25}b\,\hat{\mathbf{y}} + z_{25}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O IX}$$

$$\mathbf{B}_{50} = x_{25}\,\mathbf{a}_1 - y_{25}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{25}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{25}a + z_{25}c\cos\beta\right)\hat{\mathbf{x}} - y_{25}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{25}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O IX}$$

$$\mathbf{B}_{51} = x_{26}\,\mathbf{a}_1 + y_{26}\,\mathbf{a}_2 + z_{26}\,\mathbf{a}_3 = (x_{26}a + z_{26}c\cos\beta)\,\hat{\mathbf{x}} + y_{26}b\,\hat{\mathbf{y}} + z_{26}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O X}$$

$$\mathbf{B}_{52} = x_{26}\,\mathbf{a}_1 - y_{26}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{26}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{26}a + z_{26}c\cos\beta\right)\hat{\mathbf{x}} - y_{26}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{26}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O X}$$

$$\mathbf{B}_{53} = x_{27}\,\mathbf{a}_1 + y_{27}\,\mathbf{a}_2 + z_{27}\,\mathbf{a}_3 = (x_{27}a + z_{27}c\cos\beta)\,\hat{\mathbf{x}} + y_{27}b\,\hat{\mathbf{y}} + z_{27}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XI}$$

$$\mathbf{B}_{54} = x_{27}\,\mathbf{a}_1 - y_{27}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{27}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{27}a + z_{27}c\cos\beta\right)\hat{\mathbf{x}} - y_{27}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{27}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XI}$$

$$\mathbf{B}_{55} = x_{28}\,\mathbf{a}_1 + y_{28}\,\mathbf{a}_2 + z_{28}\,\mathbf{a}_3 = (x_{28}a + z_{28}c\cos\beta)\,\hat{\mathbf{x}} + y_{28}b\,\hat{\mathbf{y}} + z_{28}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XII}$$

$$\mathbf{B}_{56} = x_{28}\,\mathbf{a}_1 - y_{28}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{28}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{28}a + z_{28}c\cos\beta\right)\hat{\mathbf{x}} - y_{28}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{28}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XII}$$

$$\mathbf{B}_{57} = x_{29}\,\mathbf{a}_1 + y_{29}\,\mathbf{a}_2 + z_{29}\,\mathbf{a}_3 = (x_{29}a + z_{29}c\cos\beta)\,\hat{\mathbf{x}} + y_{29}b\,\hat{\mathbf{y}} + z_{29}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XIII}$$

$$\mathbf{B}_{58} = x_{29}\,\mathbf{a}_1 - y_{29}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{29}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{29}a + z_{29}c\cos\beta\right)\hat{\mathbf{x}} - y_{29}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{29}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XIII}$$

$$\mathbf{B}_{59} = x_{30}\,\mathbf{a}_1 + y_{30}\,\mathbf{a}_2 + z_{30}\,\mathbf{a}_3 = (x_{30}a + z_{30}c\cos\beta)\,\hat{\mathbf{x}} + y_{30}b\,\hat{\mathbf{y}} + z_{30}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XIV}$$

$$\mathbf{B}_{60} = x_{30}\,\mathbf{a}_1 - y_{30}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{30}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{30}a + z_{30}c\cos\beta\right)\hat{\mathbf{x}} - y_{30}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{30}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XIV}$$

$$\mathbf{B}_{61} = x_{31}\,\mathbf{a}_1 + y_{31}\,\mathbf{a}_2 + z_{31}\,\mathbf{a}_3 = (x_{31}a + z_{31}c\cos\beta)\,\hat{\mathbf{x}} + y_{31}b\,\hat{\mathbf{y}} + z_{31}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XV}$$

$$\mathbf{B}_{62} = x_{31}\,\mathbf{a}_1 - y_{31}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{31}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{31}a + z_{31}c\cos\beta\right)\hat{\mathbf{x}} - y_{31}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{31}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XV}$$

$$\mathbf{B}_{63} = x_{32}\,\mathbf{a}_1 + y_{32}\,\mathbf{a}_2 + z_{32}\,\mathbf{a}_3 = (x_{32}a + z_{32}c\cos\beta)\,\hat{\mathbf{x}} + y_{32}b\,\hat{\mathbf{y}} + z_{32}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XVI}$$

$$\mathbf{B}_{64} = x_{32}\,\mathbf{a}_1 - y_{32}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{32}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{32}a + z_{32}c\cos\beta\right)\hat{\mathbf{x}} - y_{32}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{32}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XVI}$$

$$\mathbf{B}_{65} = x_{33}\,\mathbf{a}_1 + y_{33}\,\mathbf{a}_2 + z_{33}\,\mathbf{a}_3 = (x_{33}a + z_{33}c\cos\beta)\,\hat{\mathbf{x}} + y_{33}b\,\hat{\mathbf{y}} + z_{33}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XVII}$$

$$\mathbf{B}_{66} = x_{33}\,\mathbf{a}_1 - y_{33}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{33}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{33}a + z_{33}c\cos\beta\right)\hat{\mathbf{x}} - y_{33}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{33}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XVII}$$

$$\mathbf{B}_{67} = x_{34}\,\mathbf{a}_1 + y_{34}\,\mathbf{a}_2 + z_{34}\,\mathbf{a}_3 = (x_{34}a + z_{34}c\cos\beta)\,\hat{\mathbf{x}} + y_{34}b\,\hat{\mathbf{y}} + z_{34}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XVIII}$$

$$\mathbf{B}_{68} = x_{34}\,\mathbf{a}_1 - y_{34}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{34}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{34}a + z_{34}c\cos\beta\right)\hat{\mathbf{x}} - y_{34}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{34}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XVIII}$$

$$\mathbf{B}_{69} = x_{35}\,\mathbf{a}_1 + y_{35}\,\mathbf{a}_2 + z_{35}\,\mathbf{a}_3 = (x_{35}a + z_{35}c\cos\beta)\,\hat{\mathbf{x}} + y_{35}b\,\hat{\mathbf{y}} + z_{35}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XIX}$$

$$\mathbf{B}_{70} = x_{35}\,\mathbf{a}_1 - y_{35}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{35}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{35}a + z_{35}c\cos\beta\right)\hat{\mathbf{x}} - y_{35}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{35}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XIX}$$

$$\mathbf{B}_{71} = x_{36}\,\mathbf{a}_1 + y_{36}\,\mathbf{a}_2 + z_{36}\,\mathbf{a}_3 = (x_{36}a + z_{36}c\cos\beta)\,\hat{\mathbf{x}} + y_{36}b\,\hat{\mathbf{y}} + z_{36}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XX}$$

$$\mathbf{B}_{72} = x_{36}\,\mathbf{a}_1 - y_{36}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{36}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{36}a + z_{36}c\cos\beta\right)\hat{\mathbf{x}} - y_{36}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{36}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XX}$$

$$\mathbf{B}_{73} = x_{37}\,\mathbf{a}_1 + y_{37}\,\mathbf{a}_2 + z_{37}\,\mathbf{a}_3 = (x_{37}a + z_{37}c\cos\beta)\,\hat{\mathbf{x}} + y_{37}b\,\hat{\mathbf{y}} + z_{37}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXI}$$

$$\mathbf{B}_{74} = x_{37}\,\mathbf{a}_1 - y_{37}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{37}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{37}a + z_{37}c\cos\beta\right)\hat{\mathbf{x}} - y_{37}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{37}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXI}$$

$$\mathbf{B}_{75} = x_{38}\,\mathbf{a}_1 + y_{38}\,\mathbf{a}_2 + z_{38}\,\mathbf{a}_3 = (x_{38}a + z_{38}c\cos\beta)\,\hat{\mathbf{x}} + y_{38}b\,\hat{\mathbf{y}} + z_{38}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXII}$$

$$\mathbf{B}_{76} = x_{38}\,\mathbf{a}_1 - y_{38}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{38}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{38}a + z_{38}c\cos\beta\right)\hat{\mathbf{x}} - y_{38}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{38}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXII}$$

$$\mathbf{B}_{77} = x_{39}\,\mathbf{a}_1 + y_{39}\,\mathbf{a}_2 + z_{39}\,\mathbf{a}_3 = (x_{39}a + z_{39}c\cos\beta)\,\hat{\mathbf{x}} + y_{39}b\,\hat{\mathbf{y}} + z_{39}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXIII}$$

$$\mathbf{B}_{78} = x_{39}\,\mathbf{a}_1 - y_{39}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{39}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{39}a + z_{39}c\cos\beta\right)\hat{\mathbf{x}} - y_{39}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{39}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXIII}$$

$$\mathbf{B}_{79} = x_{40}\,\mathbf{a}_1 + y_{40}\,\mathbf{a}_2 + z_{40}\,\mathbf{a}_3 = (x_{40}a + z_{40}c\cos\beta)\,\hat{\mathbf{x}} + y_{40}b\,\hat{\mathbf{y}} + z_{40}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXIV}$$

$$\mathbf{B}_{80} = x_{40}\,\mathbf{a}_1 - y_{40}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{40}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{40}a + z_{40}c\cos\beta\right)\hat{\mathbf{x}} - y_{40}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{40}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXIV}$$

$$\mathbf{B}_{81} = x_{41}\,\mathbf{a}_1 + y_{41}\,\mathbf{a}_2 + z_{41}\,\mathbf{a}_3 = (x_{41}a + z_{41}c\cos\beta)\,\hat{\mathbf{x}} + y_{41}b\,\hat{\mathbf{y}} + z_{41}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXV}$$

$$\mathbf{B}_{82} = x_{41}\,\mathbf{a}_1 - y_{41}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{41}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{41}a + z_{41}c\cos\beta\right)\hat{\mathbf{x}} - y_{41}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{41}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXV}$$

$$\mathbf{B}_{83} = x_{42}\,\mathbf{a}_1 + y_{42}\,\mathbf{a}_2 + z_{42}\,\mathbf{a}_3 = (x_{42}a + z_{42}c\cos\beta)\,\hat{\mathbf{x}} + y_{42}b\,\hat{\mathbf{y}} + z_{42}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXVI}$$

$$\mathbf{B}_{84} = x_{42}\,\mathbf{a}_1 - y_{42}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{42}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{42}a + z_{42}c\cos\beta\right)\hat{\mathbf{x}} - y_{42}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{42}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXVI}$$

$$\mathbf{B}_{85} = x_{43}\,\mathbf{a}_1 + y_{43}\,\mathbf{a}_2 + z_{43}\,\mathbf{a}_3 = (x_{43}a + z_{43}c\cos\beta)\,\hat{\mathbf{x}} + y_{43}b\,\hat{\mathbf{y}} + z_{43}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXVII}$$

$$\mathbf{B}_{86} = x_{43}\,\mathbf{a}_1 - y_{43}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{43}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{43}a + z_{43}c\cos\beta\right)\hat{\mathbf{x}} - y_{43}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{43}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXVII}$$

$$\mathbf{B}_{87} = x_{44}\,\mathbf{a}_1 + y_{44}\,\mathbf{a}_2 + z_{44}\,\mathbf{a}_3 = (x_{44}a + z_{44}c\cos\beta)\,\hat{\mathbf{x}} + y_{44}b\,\hat{\mathbf{y}} + z_{44}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXVIII}$$

$$\mathbf{B}_{88} = x_{44}\,\mathbf{a}_1 - y_{44}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{44}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{44}a + z_{44}c\cos\beta\right)\hat{\mathbf{x}} - y_{44}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{44}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXVIII}$$

$$\mathbf{B}_{89} = x_{45}\,\mathbf{a}_1 + y_{45}\,\mathbf{a}_2 + z_{45}\,\mathbf{a}_3 = (x_{45}a + z_{45}c\cos\beta)\,\hat{\mathbf{x}} + y_{45}b\,\hat{\mathbf{y}} + z_{45}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXIX}$$

$$\mathbf{B}_{90} = x_{45}\,\mathbf{a}_1 - y_{45}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{45}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{45}a + z_{45}c\cos\beta\right)\hat{\mathbf{x}} - y_{45}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{45}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXIX}$$

$$\mathbf{B}_{91} = x_{46}\,\mathbf{a}_1 + y_{46}\,\mathbf{a}_2 + z_{46}\,\mathbf{a}_3 = (x_{46}a + z_{46}c\cos\beta)\,\hat{\mathbf{x}} + y_{46}b\,\hat{\mathbf{y}} + z_{46}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXX}$$

$$\mathbf{B}_{92} = x_{46}\,\mathbf{a}_1 - y_{46}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{46}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{46}a + z_{46}c\cos\beta\right)\hat{\mathbf{x}} - y_{46}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{46}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXX}$$

$$\mathbf{B}_{93} = x_{47}\,\mathbf{a}_1 + y_{47}\,\mathbf{a}_2 + z_{47}\,\mathbf{a}_3 = (x_{47}a + z_{47}c\cos\beta)\,\hat{\mathbf{x}} + y_{47}b\,\hat{\mathbf{y}} + z_{47}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXXI}$$

$$\mathbf{B}_{94} = x_{47}\,\mathbf{a}_1 - y_{47}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{47}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{47}a + z_{47}c\cos\beta\right)\hat{\mathbf{x}} - y_{47}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{47}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXXI}$$

$$\mathbf{B}_{95} = x_{48}\,\mathbf{a}_1 + y_{48}\,\mathbf{a}_2 + z_{48}\,\mathbf{a}_3 = (x_{48}a + z_{48}c\cos\beta)\,\hat{\mathbf{x}} + y_{48}b\,\hat{\mathbf{y}} + z_{48}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXXII}$$

$$\mathbf{B}_{96} = x_{48}\,\mathbf{a}_1 - y_{48}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{48}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{48}a + z_{48}c\cos\beta\right)\hat{\mathbf{x}} - y_{48}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{48}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXXII}$$

$$\mathbf{B}_{97} = x_{49}\,\mathbf{a}_1 + y_{49}\,\mathbf{a}_2 + z_{49}\,\mathbf{a}_3 = (x_{49}a + z_{49}c\cos\beta)\,\hat{\mathbf{x}} + y_{49}b\,\hat{\mathbf{y}} + z_{49}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXXIII}$$

$$\mathbf{B}_{98} = x_{49}\,\mathbf{a}_1 - y_{49}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{49}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{49}a + z_{49}c\cos\beta\right)\hat{\mathbf{x}} - y_{49}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{49}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXXIII}$$

$$\mathbf{B}_{99} = x_{50}\,\mathbf{a}_1 + y_{50}\,\mathbf{a}_2 + z_{50}\,\mathbf{a}_3 = (x_{50}a + z_{50}c\cos\beta)\,\hat{\mathbf{x}} + y_{50}b\,\hat{\mathbf{y}} + z_{50}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXXIV}$$

$$\mathbf{B}_{100} = x_{50}\,\mathbf{a}_1 - y_{50}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{50}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{50}a + z_{50}c\cos\beta\right)\hat{\mathbf{x}} - y_{50}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{50}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXXIV}$$

$$\mathbf{B}_{101} = x_{51}\,\mathbf{a}_1 + y_{51}\,\mathbf{a}_2 + z_{51}\,\mathbf{a}_3 = (x_{51}a + z_{51}c\cos\beta)\,\hat{\mathbf{x}} + y_{51}b\,\hat{\mathbf{y}} + z_{51}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXXV}$$

$$\mathbf{B}_{102} = x_{51}\,\mathbf{a}_1 - y_{51}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{51}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{51}a + z_{51}c\cos\beta\right)\hat{\mathbf{x}} - y_{51}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{51}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXXV}$$

$$\mathbf{B}_{103} = x_{52}\,\mathbf{a}_1 + y_{52}\,\mathbf{a}_2 + z_{52}\,\mathbf{a}_3 = (x_{52}a + z_{52}c\cos\beta)\,\hat{\mathbf{x}} + y_{52}b\,\hat{\mathbf{y}} + z_{52}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXXVI}$$

$$\mathbf{B}_{104} = x_{52}\,\mathbf{a}_1 - y_{52}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{52}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{52}a + z_{52}c\cos\beta\right)\hat{\mathbf{x}} - y_{52}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{52}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXXVI}$$

$$\mathbf{B}_{105} = x_{53}\,\mathbf{a}_1 + y_{53}\,\mathbf{a}_2 + z_{53}\,\mathbf{a}_3 = (x_{53}a + z_{53}c\cos\beta)\,\hat{\mathbf{x}} + y_{53}b\,\hat{\mathbf{y}} + z_{53}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXXVII}$$

$$\mathbf{B}_{106} = x_{53}\,\mathbf{a}_1 - y_{53}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{53}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{53}a + z_{53}c\cos\beta\right)\hat{\mathbf{x}} - y_{53}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{53}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXXVII}$$

$$\mathbf{B}_{107} = x_{54}\,\mathbf{a}_1 + y_{54}\,\mathbf{a}_2 + z_{54}\,\mathbf{a}_3 = (x_{54}a + z_{54}c\cos\beta)\,\hat{\mathbf{x}} + y_{54}b\,\hat{\mathbf{y}} + z_{54}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXXVIII}$$

$$\mathbf{B}_{108} = x_{54}\,\mathbf{a}_1 - y_{54}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{54}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{54}a + z_{54}c\cos\beta\right)\hat{\mathbf{x}} - y_{54}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{54}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXXVIII}$$

$$\mathbf{B}_{109} = x_{55}\,\mathbf{a}_1 + y_{55}\,\mathbf{a}_2 + z_{55}\,\mathbf{a}_3 = (x_{55}a + z_{55}c\cos\beta)\,\hat{\mathbf{x}} + y_{55}b\,\hat{\mathbf{y}} + z_{55}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXXIX}$$

$$\mathbf{B}_{110} = x_{55}\,\mathbf{a}_1 - y_{55}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{55}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{55}a + z_{55}c\cos\beta\right)\hat{\mathbf{x}} - y_{55}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{55}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XXXIX}$$

$$\mathbf{B}_{111} = x_{56}\,\mathbf{a}_1 + y_{56}\,\mathbf{a}_2 + z_{56}\,\mathbf{a}_3 = (x_{56}a + z_{56}c\cos\beta)\,\hat{\mathbf{x}} + y_{56}b\,\hat{\mathbf{y}} + z_{56}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XL}$$

$$\mathbf{B}_{112} = x_{56}\,\mathbf{a}_1 - y_{56}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{56}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{56}a + z_{56}c\cos\beta\right)\hat{\mathbf{x}} - y_{56}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{56}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XL}$$

$$\mathbf{B}_{113} = x_{57}\,\mathbf{a}_1 + y_{57}\,\mathbf{a}_2 + z_{57}\,\mathbf{a}_3 = (x_{57}a + z_{57}c\cos\beta)\,\hat{\mathbf{x}} + y_{57}b\,\hat{\mathbf{y}} + z_{57}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XLI}$$

$$\mathbf{B}_{114} = x_{57}\,\mathbf{a}_1 - y_{57}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{57}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{57}a + z_{57}c\cos\beta\right)\hat{\mathbf{x}} - y_{57}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{57}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XLI}$$

$$\mathbf{B}_{115} = x_{58}\,\mathbf{a}_1 + y_{58}\,\mathbf{a}_2 + z_{58}\,\mathbf{a}_3 = (x_{58}a + z_{58}c\cos\beta)\,\hat{\mathbf{x}} + y_{58}b\,\hat{\mathbf{y}} + z_{58}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XLII}$$

$$\mathbf{B}_{116} = x_{58}\,\mathbf{a}_1 - y_{58}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{58}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{58}a + z_{58}c\cos\beta\right)\hat{\mathbf{x}} - y_{58}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{58}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XLII}$$

$$\mathbf{B}_{117} = x_{59}\,\mathbf{a}_1 + y_{59}\,\mathbf{a}_2 + z_{59}\,\mathbf{a}_3 = (x_{59}a + z_{59}c\cos\beta)\,\hat{\mathbf{x}} + y_{59}b\,\hat{\mathbf{y}} + z_{59}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XLIII}$$

$$\mathbf{B}_{118} = x_{59}\,\mathbf{a}_1 - y_{59}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{59}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{59}a + z_{59}c\cos\beta\right)\hat{\mathbf{x}} - y_{59}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{59}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XLIII}$$

$$\mathbf{B}_{119} = x_{60}\,\mathbf{a}_1 + y_{60}\,\mathbf{a}_2 + z_{60}\,\mathbf{a}_3 = (x_{60}a + z_{60}c\cos\beta)\,\hat{\mathbf{x}} + y_{60}b\,\hat{\mathbf{y}} + z_{60}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XLIV}$$

$$\mathbf{B}_{120} = x_{60}\,\mathbf{a}_1 - y_{60}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{60}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{60}a + z_{60}c\cos\beta\right)\hat{\mathbf{x}} - y_{60}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{60}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XLIV}$$

$$\mathbf{B}_{121} = x_{61}\,\mathbf{a}_1 + y_{61}\,\mathbf{a}_2 + z_{61}\,\mathbf{a}_3 = (x_{61}a + z_{61}c\cos\beta)\,\hat{\mathbf{x}} + y_{61}b\,\hat{\mathbf{y}} + z_{61}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XLV}$$

$$\mathbf{B}_{122} = x_{61}\,\mathbf{a}_1 - y_{61}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{61}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{61}a + z_{61}c\cos\beta\right)\hat{\mathbf{x}} - y_{61}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{61}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XLV}$$

$$\mathbf{B}_{123} = x_{62}\,\mathbf{a}_1 + y_{62}\,\mathbf{a}_2 + z_{62}\,\mathbf{a}_3 = (x_{62}a + z_{62}c\cos\beta)\,\hat{\mathbf{x}} + y_{62}b\,\hat{\mathbf{y}} + z_{62}c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XLVI}$$

$$\mathbf{B}_{124} = x_{62}\,\mathbf{a}_1 - y_{62}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{62}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{62}a + z_{62}c\cos\beta\right)\hat{\mathbf{x}} - y_{62}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{62}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (2a) \quad \text{O XLVI}$$

References:

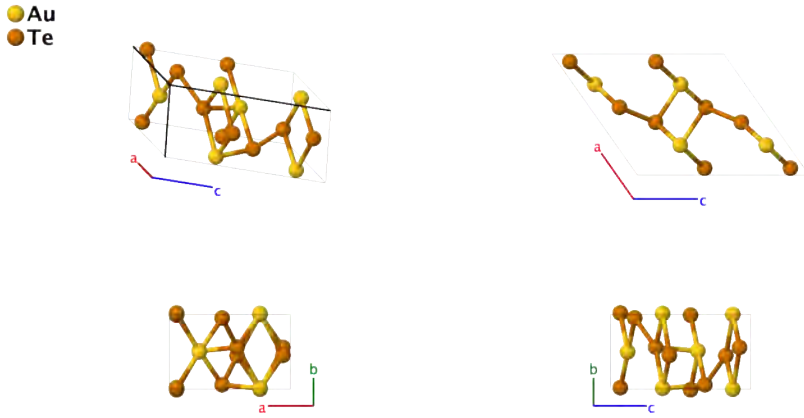
- H. Fujishita, M. Sato, S. Sato, and S. Hoshino, *Structure Determination of low-dimensional conductor Mo_8O_{23}*, J. Solid State Chem. **66**, 40–46 (1987), doi:10.1016/0022-4596(87)90218-0.

Geometry files:

- CIF: pp. 1513
- POSCAR: pp. 1514

Calaverite ($AuTe_2$) Structure: AB2_mP12_7_2a_4a

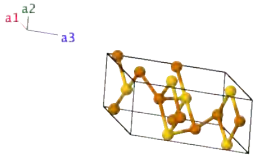
http://aflow.org/prototype-encyclopedia/AB2_mP12_7_2a_4a



Prototype	:	$AuTe_2$
AFLOW prototype label	:	AB2_mP12_7_2a_4a
***Strukturbericht* designation**	:	None
Pearson symbol	:	mP12
Space group number	:	7
Space group symbol	:	*Pc*
AFLOW prototype command	:	`aflow --proto=AB2_mP12_7_2a_4a --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6$

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	$(x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Au I
$\mathbf{B}_2$	$x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_1 a + z_1 c\cos\beta\right)\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Au I
$\mathbf{B}_3$	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Au II
$\mathbf{B}_4$	$x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_2 a + z_2 c\cos\beta\right)\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Au II
$\mathbf{B}_5$	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Te I

$\mathbf{B}_6$	=	$x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Te I
$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Te II
$\mathbf{B}_8$	=	$x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Te II
$\mathbf{B}_9$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Te III
$\mathbf{B}_{10}$	=	$x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Te III
$\mathbf{B}_{11}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Te IV
$\mathbf{B}_{12}$	=	$x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Te IV

References:

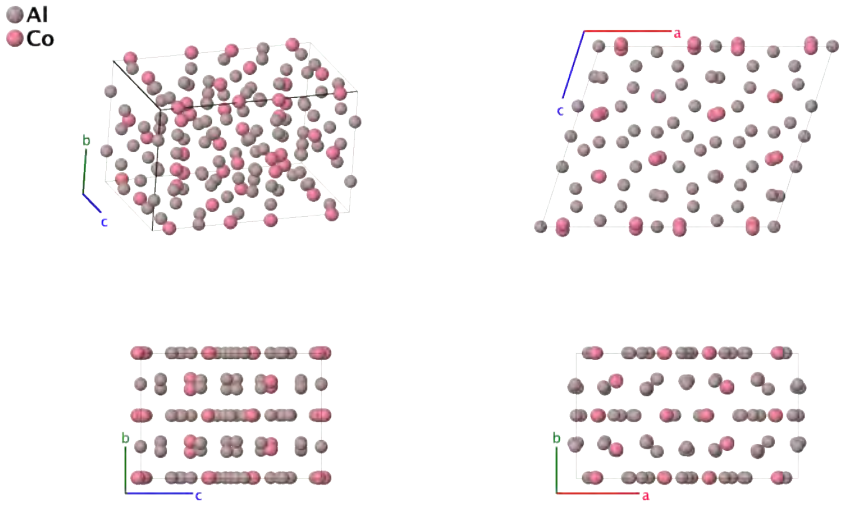
- F. Pertlik, *Kristallchemie natülicher Telluride III: Die Kristallstruktur des Minerals Calaverit, $AuTe_2$*, Zeitschrift für Kristallographie - Crystalline Materials **169**, 227–236 (1984), doi:10.1524/zkri.1984.169.14.227.

Geometry files:

- CIF: pp. 1515
- POSCAR: pp. 1515

Monoclinic Co_4Al_{13} Structure: A13B4_mC102_8_17a11b_8a2b

http://aflow.org/prototype-encyclopedia/A13B4_mC102_8_17a11b_8a2b

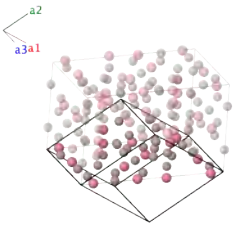


Prototype	:	$Al_{13}Co_4$
AFLOW prototype label	:	A13B4_mC102_8_17a11b_8a2b
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC102
Space group number	:	8
Space group symbol	:	*Cm*
AFLOW prototype command	:	`aflow --proto=A13B4_mC102_8_17a11b_8a2b --params=`$a, b/a, c/a, \beta, x_1, z_1, x_2, z_2, x_3, z_3, x_4, z_4, x_5, z_5, x_6, z_6, x_7, z_7, x_8, z_8, x_9, z_9, x_{10}, z_{10}, x_{11}, z_{11}, x_{12}, z_{12}, x_{13}, z_{13}, x_{14}, z_{14}, x_{15}, z_{15}, x_{16}, z_{16}, x_{17}, z_{17}, x_{18}, z_{18}, x_{19}, z_{19}, x_{20}, z_{20}, x_{21}, z_{21}, x_{22}, z_{22}, x_{23}, z_{23}, x_{24}, z_{24}, x_{25}, z_{25}, x_{26}, y_{26}, z_{26}, x_{27}, y_{27}, z_{27}, x_{28}, y_{28}, z_{28}, x_{29}, y_{29}, z_{29}, x_{30}, y_{30}, z_{30}, x_{31}, y_{31}, z_{31}, x_{32}, y_{32}, z_{32}, x_{33}, y_{33}, z_{33}, x_{34}, y_{34}, z_{34}, x_{35}, y_{35}, z_{35}, x_{36}, y_{36}, z_{36}, x_{37}, y_{37}, z_{37}, x_{38}, y_{38}, z_{38}$

- Following (Hudd, 1962), the Al-IV and Al-XIII sites are occupied 30% of the time, while the occupation of Al-VI, Al-IX, Al-XIV, and Al-XVII is 70%. This gives a nominial occupation of $Al_{91}Co_{30}$, though the authors state the actual composition is $Al_{68.3}Co_{24.4}$.
- Space group *Cm* #8 allows an arbitrary choice for the origin of the z-axis. We follow (Hudd, 1962) and set the $z_{26} = 0$.
- AFLOW-SYM yields space group *Cm* #8 consistently for small and large tolerances. However, if we allow a rather large uncertainty of 0.3 Å in the atomic positions with FINDSYM, the symmetry is set as $C2/m$ #12. That crystal has the $Al_{13}Fe_4$ prototype.

Base-centered Monoclinic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2}\,b\,\hat{\mathbf{y}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,b\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + x_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1a + z_1c\cos\beta)\,\hat{\mathbf{x}} + z_1c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Al I
$\mathbf{B}_2$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2a + z_2c\cos\beta)\,\hat{\mathbf{x}} + z_2c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Al II
$\mathbf{B}_3$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3a + z_3c\cos\beta)\,\hat{\mathbf{x}} + z_3c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Al III
$\mathbf{B}_4$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4a + z_4c\cos\beta)\,\hat{\mathbf{x}} + z_4c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Al IV
$\mathbf{B}_5$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5a + z_5c\cos\beta)\,\hat{\mathbf{x}} + z_5c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Al V
$\mathbf{B}_6$	=	$x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6a + z_6c\cos\beta)\,\hat{\mathbf{x}} + z_6c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Al VI
$\mathbf{B}_7$	=	$x_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7a + z_7c\cos\beta)\,\hat{\mathbf{x}} + z_7c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Al VII
$\mathbf{B}_8$	=	$x_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8a + z_8c\cos\beta)\,\hat{\mathbf{x}} + z_8c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Al VIII
$\mathbf{B}_9$	=	$x_9\,\mathbf{a}_1 + x_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$(x_9a + z_9c\cos\beta)\,\hat{\mathbf{x}} + z_9c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Al IX
$\mathbf{B}_{10}$	=	$x_{10}\,\mathbf{a}_1 + x_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$(x_{10}a + z_{10}c\cos\beta)\,\hat{\mathbf{x}} + z_{10}c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Al X
$\mathbf{B}_{11}$	=	$x_{11}\,\mathbf{a}_1 + x_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$(x_{11}a + z_{11}c\cos\beta)\,\hat{\mathbf{x}} + z_{11}c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Al XI
$\mathbf{B}_{12}$	=	$x_{12}\,\mathbf{a}_1 + x_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3$	=	$(x_{12}a + z_{12}c\cos\beta)\,\hat{\mathbf{x}} + z_{12}c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Al XII
$\mathbf{B}_{13}$	=	$x_{13}\,\mathbf{a}_1 + x_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3$	=	$(x_{13}a + z_{13}c\cos\beta)\,\hat{\mathbf{x}} + z_{13}c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Al XIII
$\mathbf{B}_{14}$	=	$x_{14}\,\mathbf{a}_1 + x_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3$	=	$(x_{14}a + z_{14}c\cos\beta)\,\hat{\mathbf{x}} + z_{14}c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Al XIV
$\mathbf{B}_{15}$	=	$x_{15}\,\mathbf{a}_1 + x_{15}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3$	=	$(x_{15}a + z_{15}c\cos\beta)\,\hat{\mathbf{x}} + z_{15}c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Al XV
$\mathbf{B}_{16}$	=	$x_{16}\,\mathbf{a}_1 + x_{16}\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3$	=	$(x_{16}a + z_{16}c\cos\beta)\,\hat{\mathbf{x}} + z_{16}c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Al XVI
$\mathbf{B}_{17}$	=	$x_{17}\,\mathbf{a}_1 + x_{17}\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3$	=	$(x_{17}a + z_{17}c\cos\beta)\,\hat{\mathbf{x}} + z_{17}c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Al XVII
$\mathbf{B}_{18}$	=	$x_{18}\,\mathbf{a}_1 + x_{18}\,\mathbf{a}_2 + z_{18}\,\mathbf{a}_3$	=	$(x_{18}a + z_{18}c\cos\beta)\,\hat{\mathbf{x}} + z_{18}c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Co I
$\mathbf{B}_{19}$	=	$x_{19}\,\mathbf{a}_1 + x_{19}\,\mathbf{a}_2 + z_{19}\,\mathbf{a}_3$	=	$(x_{19}a + z_{19}c\cos\beta)\,\hat{\mathbf{x}} + z_{19}c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Co II
$\mathbf{B}_{20}$	=	$x_{20}\,\mathbf{a}_1 + x_{20}\,\mathbf{a}_2 + z_{20}\,\mathbf{a}_3$	=	$(x_{20}a + z_{20}c\cos\beta)\,\hat{\mathbf{x}} + z_{20}c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Co III
$\mathbf{B}_{21}$	=	$x_{21}\,\mathbf{a}_1 + x_{21}\,\mathbf{a}_2 + z_{21}\,\mathbf{a}_3$	=	$(x_{21}a + z_{21}c\cos\beta)\,\hat{\mathbf{x}} + z_{21}c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Co IV
$\mathbf{B}_{22}$	=	$x_{22}\,\mathbf{a}_1 + x_{22}\,\mathbf{a}_2 + z_{22}\,\mathbf{a}_3$	=	$(x_{22}a + z_{22}c\cos\beta)\,\hat{\mathbf{x}} + z_{22}c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Co V
$\mathbf{B}_{23}$	=	$x_{23}\,\mathbf{a}_1 + x_{23}\,\mathbf{a}_2 + z_{23}\,\mathbf{a}_3$	=	$(x_{23}a + z_{23}c\cos\beta)\,\hat{\mathbf{x}} + z_{23}c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Co VI
$\mathbf{B}_{24}$	=	$x_{24}\,\mathbf{a}_1 + x_{24}\,\mathbf{a}_2 + z_{24}\,\mathbf{a}_3$	=	$(x_{24}a + z_{24}c\cos\beta)\,\hat{\mathbf{x}} + z_{24}c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Co VII
$\mathbf{B}_{25}$	=	$x_{25}\,\mathbf{a}_1 + x_{25}\,\mathbf{a}_2 + z_{25}\,\mathbf{a}_3$	=	$(x_{25}a + z_{25}c\cos\beta)\,\hat{\mathbf{x}} + z_{25}c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Co VIII
$\mathbf{B}_{26}$	=	$(x_{26} - y_{26})\,\mathbf{a}_1 + (x_{26} + y_{26})\,\mathbf{a}_2 + z_{26}\,\mathbf{a}_3$	=	$(x_{26}a + z_{26}c\cos\beta)\,\hat{\mathbf{x}} + y_{26}b\,\hat{\mathbf{y}} + z_{26}c\sin\beta\,\hat{\mathbf{z}}$	$(4b)$	Al XVIII
$\mathbf{B}_{27}$	=	$(x_{26} + y_{26})\,\mathbf{a}_1 + (x_{26} - y_{26})\,\mathbf{a}_2 + z_{26}\,\mathbf{a}_3$	=	$(x_{26}a + z_{26}c\cos\beta)\,\hat{\mathbf{x}} - y_{26}b\,\hat{\mathbf{y}} + z_{26}c\sin\beta\,\hat{\mathbf{z}}$	$(4b)$	Al XVIII

$$\mathbf{B}_{28} = (x_{27} - y_{27})\,\mathbf{a}_1 + (x_{27} + y_{27})\,\mathbf{a}_2 + z_{27}\,\mathbf{a}_3 = (x_{27}a + z_{27}c\cos\beta)\,\hat{\mathbf{x}} + y_{27}b\,\hat{\mathbf{y}} + z_{27}c\sin\beta\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XIX}$$

$$\mathbf{B}_{29} = (x_{27} + y_{27})\,\mathbf{a}_1 + (x_{27} - y_{27})\,\mathbf{a}_2 + z_{27}\,\mathbf{a}_3 = (x_{27}a + z_{27}c\cos\beta)\,\hat{\mathbf{x}} - y_{27}b\,\hat{\mathbf{y}} + z_{27}c\sin\beta\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XIX}$$

$$\mathbf{B}_{30} = (x_{28} - y_{28})\,\mathbf{a}_1 + (x_{28} + y_{28})\,\mathbf{a}_2 + z_{28}\,\mathbf{a}_3 = (x_{28}a + z_{28}c\cos\beta)\,\hat{\mathbf{x}} + y_{28}b\,\hat{\mathbf{y}} + z_{28}c\sin\beta\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XX}$$

$$\mathbf{B}_{31} = (x_{28} + y_{28})\,\mathbf{a}_1 + (x_{28} - y_{28})\,\mathbf{a}_2 + z_{28}\,\mathbf{a}_3 = (x_{28}a + z_{28}c\cos\beta)\,\hat{\mathbf{x}} - y_{28}b\,\hat{\mathbf{y}} + z_{28}c\sin\beta\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XX}$$

$$\mathbf{B}_{32} = (x_{29} - y_{29})\,\mathbf{a}_1 + (x_{29} + y_{29})\,\mathbf{a}_2 + z_{29}\,\mathbf{a}_3 = (x_{29}a + z_{29}c\cos\beta)\,\hat{\mathbf{x}} + y_{29}b\,\hat{\mathbf{y}} + z_{29}c\sin\beta\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXI}$$

$$\mathbf{B}_{33} = (x_{29} + y_{29})\,\mathbf{a}_1 + (x_{29} - y_{29})\,\mathbf{a}_2 + z_{29}\,\mathbf{a}_3 = (x_{29}a + z_{29}c\cos\beta)\,\hat{\mathbf{x}} - y_{29}b\,\hat{\mathbf{y}} + z_{29}c\sin\beta\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXI}$$

$$\mathbf{B}_{34} = (x_{30} - y_{30})\,\mathbf{a}_1 + (x_{30} + y_{30})\,\mathbf{a}_2 + z_{30}\,\mathbf{a}_3 = (x_{30}a + z_{30}c\cos\beta)\,\hat{\mathbf{x}} + y_{30}b\,\hat{\mathbf{y}} + z_{30}c\sin\beta\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXII}$$

$$\mathbf{B}_{35} = (x_{30} + y_{30})\,\mathbf{a}_1 + (x_{30} - y_{30})\,\mathbf{a}_2 + z_{30}\,\mathbf{a}_3 = (x_{30}a + z_{30}c\cos\beta)\,\hat{\mathbf{x}} - y_{30}b\,\hat{\mathbf{y}} + z_{30}c\sin\beta\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXII}$$

$$\mathbf{B}_{36} = (x_{31} - y_{31})\,\mathbf{a}_1 + (x_{31} + y_{31})\,\mathbf{a}_2 + z_{31}\,\mathbf{a}_3 = (x_{31}a + z_{31}c\cos\beta)\,\hat{\mathbf{x}} + y_{31}b\,\hat{\mathbf{y}} + z_{31}c\sin\beta\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXIII}$$

$$\mathbf{B}_{37} = (x_{31} + y_{31})\,\mathbf{a}_1 + (x_{31} - y_{31})\,\mathbf{a}_2 + z_{31}\,\mathbf{a}_3 = (x_{31}a + z_{31}c\cos\beta)\,\hat{\mathbf{x}} - y_{31}b\,\hat{\mathbf{y}} + z_{31}c\sin\beta\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXIII}$$

$$\mathbf{B}_{38} = (x_{32} - y_{32})\,\mathbf{a}_1 + (x_{32} + y_{32})\,\mathbf{a}_2 + z_{32}\,\mathbf{a}_3 = (x_{32}a + z_{32}c\cos\beta)\,\hat{\mathbf{x}} + y_{32}b\,\hat{\mathbf{y}} + z_{32}c\sin\beta\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXIV}$$

$$\mathbf{B}_{39} = (x_{32} + y_{32})\,\mathbf{a}_1 + (x_{32} - y_{32})\,\mathbf{a}_2 + z_{32}\,\mathbf{a}_3 = (x_{32}a + z_{32}c\cos\beta)\,\hat{\mathbf{x}} - y_{32}b\,\hat{\mathbf{y}} + z_{32}c\sin\beta\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXIV}$$

$$\mathbf{B}_{40} = (x_{33} - y_{33})\,\mathbf{a}_1 + (x_{33} + y_{33})\,\mathbf{a}_2 + z_{33}\,\mathbf{a}_3 = (x_{33}a + z_{33}c\cos\beta)\,\hat{\mathbf{x}} + y_{33}b\,\hat{\mathbf{y}} + z_{33}c\sin\beta\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXV}$$

$$\mathbf{B}_{41} = (x_{33} + y_{33})\,\mathbf{a}_1 + (x_{33} - y_{33})\,\mathbf{a}_2 + z_{33}\,\mathbf{a}_3 = (x_{33}a + z_{33}c\cos\beta)\,\hat{\mathbf{x}} - y_{33}b\,\hat{\mathbf{y}} + z_{33}c\sin\beta\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXV}$$

$$\mathbf{B}_{42} = (x_{34} - y_{34})\,\mathbf{a}_1 + (x_{34} + y_{34})\,\mathbf{a}_2 + z_{34}\,\mathbf{a}_3 = (x_{34}a + z_{34}c\cos\beta)\,\hat{\mathbf{x}} + y_{34}b\,\hat{\mathbf{y}} + z_{34}c\sin\beta\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXVI}$$

$$\mathbf{B}_{43} = (x_{34} + y_{34})\,\mathbf{a}_1 + (x_{34} - y_{34})\,\mathbf{a}_2 + z_{34}\,\mathbf{a}_3 = (x_{34}a + z_{34}c\cos\beta)\,\hat{\mathbf{x}} - y_{34}b\,\hat{\mathbf{y}} + z_{34}c\sin\beta\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXVI}$$

$$\mathbf{B}_{44} = (x_{35} - y_{35})\,\mathbf{a}_1 + (x_{35} + y_{35})\,\mathbf{a}_2 + z_{35}\,\mathbf{a}_3 = (x_{35}a + z_{35}c\cos\beta)\,\hat{\mathbf{x}} + y_{35}b\,\hat{\mathbf{y}} + z_{35}c\sin\beta\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXVII}$$

$$\mathbf{B}_{45} = (x_{35} + y_{35})\,\mathbf{a}_1 + (x_{35} - y_{35})\,\mathbf{a}_2 + z_{35}\,\mathbf{a}_3 = (x_{35}a + z_{35}c\cos\beta)\,\hat{\mathbf{x}} - y_{35}b\,\hat{\mathbf{y}} + z_{35}c\sin\beta\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXVII}$$

$$\mathbf{B}_{46} = (x_{36} - y_{36})\,\mathbf{a}_1 + (x_{36} + y_{36})\,\mathbf{a}_2 + z_{36}\,\mathbf{a}_3 = (x_{36}a + z_{36}c\cos\beta)\,\hat{\mathbf{x}} + y_{36}b\,\hat{\mathbf{y}} + z_{36}c\sin\beta\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXVIII}$$

$$\mathbf{B}_{47} = (x_{36} + y_{36})\,\mathbf{a}_1 + (x_{36} - y_{36})\,\mathbf{a}_2 + z_{36}\,\mathbf{a}_3 = (x_{36}a + z_{36}c\cos\beta)\,\hat{\mathbf{x}} - y_{36}b\,\hat{\mathbf{y}} + z_{36}c\sin\beta\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXVIII}$$

$$\mathbf{B}_{48} = (x_{37} - y_{37})\,\mathbf{a}_1 + (x_{37} + y_{37})\,\mathbf{a}_2 + z_{37}\,\mathbf{a}_3 = (x_{37}a + z_{37}c\cos\beta)\,\hat{\mathbf{x}} + y_{37}b\,\hat{\mathbf{y}} + z_{37}c\sin\beta\,\hat{\mathbf{z}} \quad (4b) \quad \text{Co IX}$$

$$\mathbf{B}_{49} = (x_{37} + y_{37})\,\mathbf{a}_1 + (x_{37} - y_{37})\,\mathbf{a}_2 + z_{37}\,\mathbf{a}_3 = (x_{37}a + z_{37}c\cos\beta)\,\hat{\mathbf{x}} - y_{37}b\,\hat{\mathbf{y}} + z_{37}c\sin\beta\,\hat{\mathbf{z}} \quad (4b) \quad \text{Co IX}$$

$$\mathbf{B}_{50} = (x_{38} - y_{38})\,\mathbf{a}_1 + (x_{38} + y_{38})\,\mathbf{a}_2 + z_{38}\,\mathbf{a}_3 = (x_{38}a + z_{38}c\cos\beta)\,\hat{\mathbf{x}} + y_{38}b\,\hat{\mathbf{y}} + z_{38}c\sin\beta\,\hat{\mathbf{z}} \quad (4b) \quad \text{Co X}$$

$$\mathbf{B}_{51} = (x_{38} + y_{38})\,\mathbf{a}_1 + (x_{38} - y_{38})\,\mathbf{a}_2 + z_{38}\,\mathbf{a}_3 = (x_{38}a + z_{38}c\cos\beta)\,\hat{\mathbf{x}} - y_{38}b\,\hat{\mathbf{y}} + z_{38}c\sin\beta\,\hat{\mathbf{z}} \quad (4b) \quad \text{Co X}$$

References:

- R. C. Hudd and W. H. Taylor, *The Structure of Co_4Al_{13}*, Acta Cryst. **15**, 441–442 (1962), doi:10.1107/S0365110X62001103.

Found in:

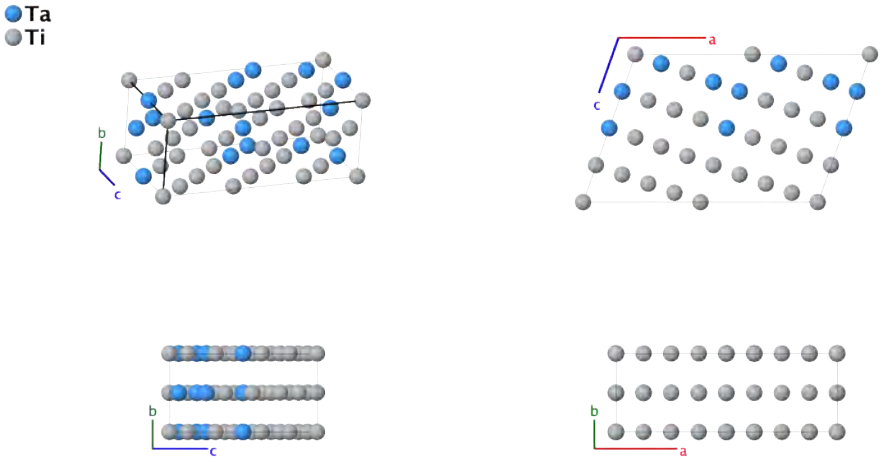
- R. Addou, E. Gaudry, T. Deniozou, M. Heggen, M. Feuerbacher, P. Gille, Y. Grin, R. Widmer, O. Gröning, V. Fournée, J.-M. Dubois, and J. Ledieu, *Structure investigation of the (100) surface of the orthorhombic $Al_{13}Co_4$ crystal*, Phys. Rev. B **80**, 014203 (2009), doi:10.1103/PhysRevB.80.014203.
- T. B. Massalski, H. Okamoto, P. R. Subramanian, and L. Kacprzak, eds., *Binary Alloy Phase Diagrams*, vol. 1 (ASM International, Materials Park, Ohio, USA, 1990), 2^{nd} edn. Ac-Ag to Ca-Zn.

Geometry files:

- CIF: pp. 1515
- POSCAR: pp. 1516

$TaTi_3$ (BCC SQS-16) Structure: AB3_mC32_8_4a_12a

http://aflow.org/prototype-encyclopedia/AB3_mC32_8_4a_12a

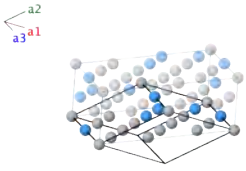


Prototype	:	$TaTi_3$
AFLOW prototype label	:	AB3_mC32_8_4a_12a
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC32
Space group number	:	8
Space group symbol	:	*Cm*
AFLOW prototype command	:	`aflow --proto=AB3_mC32_8_4a_12a --params=`$a, b/a, c/a, \beta, x_1, z_1, x_2, z_2, x_3, z_3, x_4, z_4, x_5, z_5, x_6, z_6, x_7, z_7, x_8, z_8, x_9, z_9, x_{10}, z_{10}, x_{11}, z_{11}, x_{12}, z_{12}, x_{13}, z_{13}, x_{14}, z_{14}, x_{15}, z_{15}, x_{16}, z_{16}$

- This is a special quasirandom structure with 16 atoms per unit cell (SQS-16) for a bcc binary substitutional alloy A_xB_{1-x} (Jiang, 2004). This prototype represents the $x = 0.25$ and $x = 0.75$ structures (change concentration by swapping elements). The $x = 0.5$ structure is given by AB_aP16_2_4i_4i.

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + x_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1a + z_1c\cos\beta)\,\hat{\mathbf{x}} + z_1c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	Ta I
$\mathbf{B}_2$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2a + z_2c\cos\beta)\,\hat{\mathbf{x}} + z_2c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	Ta II
$\mathbf{B}_3$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3a + z_3c\cos\beta)\,\hat{\mathbf{x}} + z_3c\sin\beta\,\hat{\mathbf{z}}$	(2*a*)	Ta III

$\mathbf{B}_4$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Ta IV
$\mathbf{B}_5$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Ti I
$\mathbf{B}_6$	=	$x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Ti II
$\mathbf{B}_7$	=	$x_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Ti III
$\mathbf{B}_8$	=	$x_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Ti IV
$\mathbf{B}_9$	=	$x_9\,\mathbf{a}_1 + x_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$(x_9 a + z_9 c\cos\beta)\,\hat{\mathbf{x}} + z_9 c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Ti V
$\mathbf{B}_{10}$	=	$x_{10}\,\mathbf{a}_1 + x_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$(x_{10} a + z_{10} c\cos\beta)\,\hat{\mathbf{x}} + z_{10} c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Ti VI
$\mathbf{B}_{11}$	=	$x_{11}\,\mathbf{a}_1 + x_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$(x_{11} a + z_{11} c\cos\beta)\,\hat{\mathbf{x}} + z_{11} c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Ti VII
$\mathbf{B}_{12}$	=	$x_{12}\,\mathbf{a}_1 + x_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3$	=	$(x_{12} a + z_{12} c\cos\beta)\,\hat{\mathbf{x}} + z_{12} c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Ti VIII
$\mathbf{B}_{13}$	=	$x_{13}\,\mathbf{a}_1 + x_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3$	=	$(x_{13} a + z_{13} c\cos\beta)\,\hat{\mathbf{x}} + z_{13} c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Ti IX
$\mathbf{B}_{14}$	=	$x_{14}\,\mathbf{a}_1 + x_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3$	=	$(x_{14} a + z_{14} c\cos\beta)\,\hat{\mathbf{x}} + z_{14} c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Ti X
$\mathbf{B}_{15}$	=	$x_{15}\,\mathbf{a}_1 + x_{15}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3$	=	$(x_{15} a + z_{15} c\cos\beta)\,\hat{\mathbf{x}} + z_{15} c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Ti XI
$\mathbf{B}_{16}$	=	$x_{16}\,\mathbf{a}_1 + x_{16}\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3$	=	$(x_{16} a + z_{16} c\cos\beta)\,\hat{\mathbf{x}} + z_{16} c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Ti XII

References:

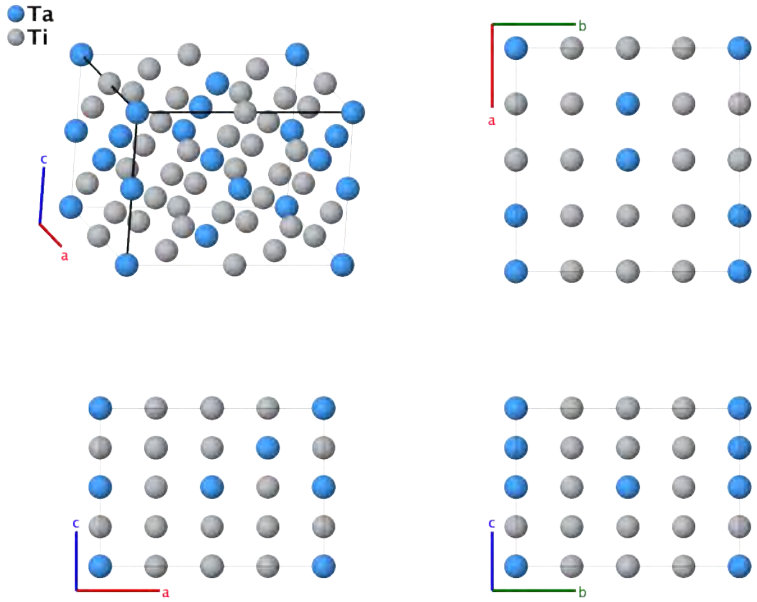
- C. Jiang, C. Wolverton, J. Sofo, L.-Q. Chen, and Z.-K. Liu, *First-principles study of binary bcc alloys using special quasirandom structures*, Phys. Rev. B **69**, 214202 (2004), doi:10.1103/PhysRevB.69.214202.

Geometry files:

- CIF: pp. 1516
- POSCAR: pp. 1517

$TaTi_3$ (BCC SQS-16) Structure: AB3_mC32_8_4a_4a4b

http://aflow.org/prototype-encyclopedia/AB3_mC32_8_4a_4a4b

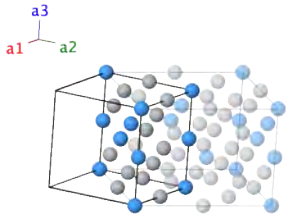


Prototype	:	$TaTi_3$
AFLOW prototype label	:	AB3_mC32_8_4a_4a4b
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC32
Space group number	:	8
Space group symbol	:	*Cm*
AFLOW prototype command	:	`aflow --proto=AB3_mC32_8_4a_4a4b --params=`$a, b/a, c/a, \beta, x_1, z_1, x_2, z_2, x_3, z_3, x_4, z_4, x_5, z_5, x_6, z_6, x_7, z_7, x_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}$

- This is a special quasirandom structure with 16 atoms per unit cell (SQS-16) for the β-phase (high-temperature austenite) bcc substitutional Ti-Ta alloy (Chakraborty, 2016). This protoype contains 25% Ta. Prototypes are listed for other Ta-Ti concentrations: 12.5% Ta (AB7_hR16_166_c_c2h), 18.75% Ta (A3B13_oC32_38_ac_a2bcdef), 31.25% Ta (A5B11_mP16_6_2abc_2a3b3c), and 37.5% Ta (A3B5_oC32_38_abce_abcdf).

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + x_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Ta I
$\mathbf{B}_2$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Ta II
$\mathbf{B}_3$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Ta III
$\mathbf{B}_4$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Ta IV
$\mathbf{B}_5$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Ti I
$\mathbf{B}_6$	=	$x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Ti II
$\mathbf{B}_7$	=	$x_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Ti III
$\mathbf{B}_8$	=	$x_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Ti IV
$\mathbf{B}_9$	=	$(x_9 - y_9)\,\mathbf{a}_1 + (x_9 + y_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$(x_9 a + z_9 c\cos\beta)\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\sin\beta\,\hat{\mathbf{z}}$	$(4b)$	Ti V
$\mathbf{B}_{10}$	=	$(x_9 + y_9)\,\mathbf{a}_1 + (x_9 - y_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$(x_9 a + z_9 c\cos\beta)\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + z_9 c\sin\beta\,\hat{\mathbf{z}}$	$(4b)$	Ti V
$\mathbf{B}_{11}$	=	$(x_{10} - y_{10})\,\mathbf{a}_1 + (x_{10} + y_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$(x_{10} a + z_{10} c\cos\beta)\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\sin\beta\,\hat{\mathbf{z}}$	$(4b)$	Ti VI
$\mathbf{B}_{12}$	=	$(x_{10} + y_{10})\,\mathbf{a}_1 + (x_{10} - y_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$(x_{10} a + z_{10} c\cos\beta)\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} + z_{10} c\sin\beta\,\hat{\mathbf{z}}$	$(4b)$	Ti VI
$\mathbf{B}_{13}$	=	$(x_{11} - y_{11})\,\mathbf{a}_1 + (x_{11} + y_{11})\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$(x_{11} a + z_{11} c\cos\beta)\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + z_{11} c\sin\beta\,\hat{\mathbf{z}}$	$(4b)$	Ti VII
$\mathbf{B}_{14}$	=	$(x_{11} + y_{11})\,\mathbf{a}_1 + (x_{11} - y_{11})\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$(x_{11} a + z_{11} c\cos\beta)\,\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} + z_{11} c\sin\beta\,\hat{\mathbf{z}}$	$(4b)$	Ti VII
$\mathbf{B}_{15}$	=	$(x_{12} - y_{12})\,\mathbf{a}_1 + (x_{12} + y_{12})\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3$	=	$(x_{12} a + z_{12} c\cos\beta)\,\hat{\mathbf{x}} + y_{12} b\,\hat{\mathbf{y}} + z_{12} c\sin\beta\,\hat{\mathbf{z}}$	$(4b)$	Ti VIII
$\mathbf{B}_{16}$	=	$(x_{12} + y_{12})\,\mathbf{a}_1 + (x_{12} - y_{12})\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3$	=	$(x_{12} a + z_{12} c\cos\beta)\,\hat{\mathbf{x}} - y_{12} b\,\hat{\mathbf{y}} + z_{12} c\sin\beta\,\hat{\mathbf{z}}$	$(4b)$	Ti VIII

References:

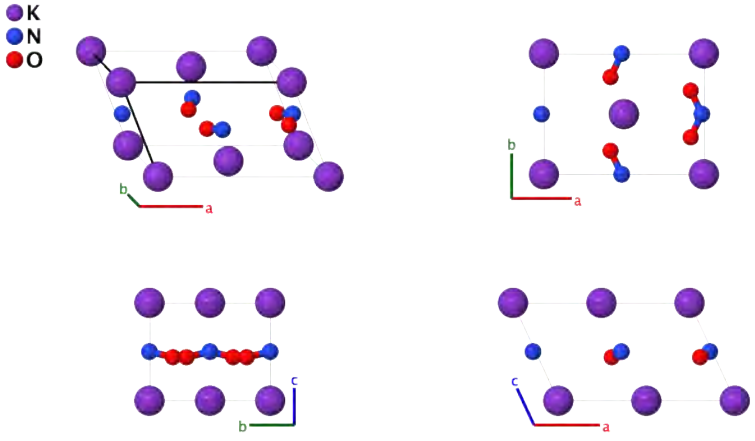
- T. Chakraborty, J. Rogal, and R. Drautz, *Unraveling the composition dependence of the martensitic transformation temperature: A first-principles study of Ti-Ta alloys*, Phys. Rev. B **94**, 224104 (2016), doi:10.1103/PhysRevB.94.224104.

Geometry files:

- CIF: pp. 1517
- POSCAR: pp. 1517

$F5_{11}$ (KNO_2) (*obsolete*) Structure: ABC2_mC8_8_a_a_b

http://aflow.org/prototype-encyclopedia/ABC2_mC8_8_a_a_b

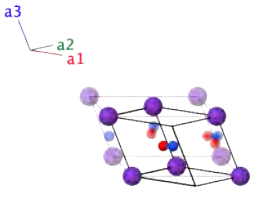


Prototype	:	KNO_2
AFLOW prototype label	:	ABC2_mC8_8_a_a_b
***Strukturbericht* designation**	:	$F5_{11}$
Pearson symbol	:	mC8
Space group number	:	8
Space group symbol	:	*Cm*
AFLOW prototype command	:	`aflow --proto=ABC2_mC8_8_a_a_b --params=`$a, b/a, c/a, \beta, x_1, z_1, x_2, z_2, x_3, y_3, z_3$

- "The room-temperature structure of KNO_2 was first considered to have monoclinic symmetry, ..., but recent studies have established the structure to be rhombohedral ..." (Rao, 1975). The $F5_{11}$ structure is thus neither the ground state structure of KNO_2 nor the room-temperature structure, which is somewhat disordered with space group $R\bar{3}m$. We present this structure as part of the historical record.
- (Ziegler, 1936) gave this structure in the *Am* setting of space group #8. We used FINDSYM to transform it to the standard *Cm* setting, which involved a considerable change in the orientation and length of the primitive lattice vectors.

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$ =	$x_1\,\mathbf{a}_1 + x_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	K

$\mathbf{B}_2$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	N
$\mathbf{B}_3$	=	$(x_3 - y_3)\,\mathbf{a}_1 + (x_3 + y_3)\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(4b)$	O
$\mathbf{B}_4$	=	$(x_3 + y_3)\,\mathbf{a}_1 + (x_3 - y_3)\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(4b)$	O

References:

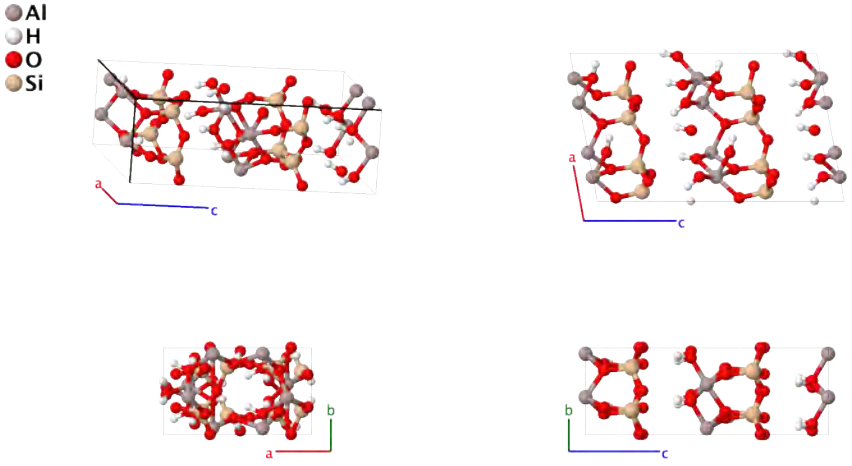
- G. E. Ziegler, *The Crystal Structure of Potassium Nitrite, KNO_2*, Zeitschrift für Kristallographie - Crystalline Materials **94**, 491–499 (1936), doi:10.1524/zkri.1936.94.1.491.
- C. N. R. Rao, B. Prakash, and M. Natarajan, *Crystal Structure Transformations in Inorganic Nitrities, Nitrates, and Carbonates* (National Bureau of Standards, 1975). National Standard Reference Data Series, NSRDS-NBS 53.

Geometry files:

- CIF: pp. 1517
- POSCAR: pp. 1518

Nacrite [$Al_2Si_2O_5(OH)_4$, $S5_4$] Structure: A2B4C9D2_mC68_9_2a_4a_9a_2a

http://aflow.org/prototype-encyclopedia/A2B4C9D2_mC68_9_2a_4a_9a_2a

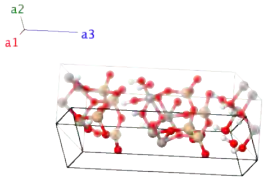


Prototype	:	$Al_2H_4O_9Si_2$
AFLOW prototype label	:	A2B4C9D2_mC68_9_2a_4a_9a_2a
***Strukturbericht* designation**	:	$S5_4$
Pearson symbol	:	mC68
Space group number	:	9
Space group symbol	:	Cc
AFLOW prototype command	:	`aflow --proto=A2B4C9D2_mC68_9_2a_4a_9a_2a --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}, x_{17}, y_{17}, z_{17}$

- (Herrmann, 1943) gave the nacrite structure determined by (Hendricks, 1939) the *Strukturbericht* symbol $S5_4$. This structure was somewhat uncertain, as Hendricks knew that the structure had space group Cc but could only determine the structure in terms of a pseudo-rhombohedral unit cell. This led to a tripling of the unit cell compared to the work of (Zhukhlistov, 2008), who was also able to determine the positions of the hydrogen atoms. We use the modern structure, noting that we can essentially recover the original structure by tripling the unit cell along the c-axis.

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$$\mathbf{B}_1 = (x_1 - y_1)\,\mathbf{a}_1 + (x_1 + y_1)\,\mathbf{a}_2 + z_1\,\mathbf{a}_3 = (x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{Al I}$$

$$\mathbf{B}_2 = (x_1 + y_1)\,\mathbf{a}_1 + (x_1 - y_1)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_1\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_1 a + z_1 c\cos\beta\right)\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_1\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{Al I}$$

$$\mathbf{B}_3 = (x_2 - y_2)\,\mathbf{a}_1 + (x_2 + y_2)\,\mathbf{a}_2 + z_2\,\mathbf{a}_3 = (x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{Al II}$$

$$\mathbf{B}_4 = (x_2 + y_2)\,\mathbf{a}_1 + (x_2 - y_2)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_2\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_2 a + z_2 c\cos\beta\right)\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_2\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{Al II}$$

$$\mathbf{B}_5 = (x_3 - y_3)\,\mathbf{a}_1 + (x_3 + y_3)\,\mathbf{a}_2 + z_3\,\mathbf{a}_3 = (x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{H I}$$

$$\mathbf{B}_6 = (x_3 + y_3)\,\mathbf{a}_1 + (x_3 - y_3)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_3\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{H I}$$

$$\mathbf{B}_7 = (x_4 - y_4)\,\mathbf{a}_1 + (x_4 + y_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3 = (x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{H II}$$

$$\mathbf{B}_8 = (x_4 + y_4)\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_4\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{H II}$$

$$\mathbf{B}_9 = (x_5 - y_5)\,\mathbf{a}_1 + (x_5 + y_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3 = (x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{H III}$$

$$\mathbf{B}_{10} = (x_5 + y_5)\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_5\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{H III}$$

$$\mathbf{B}_{11} = (x_6 - y_6)\,\mathbf{a}_1 + (x_6 + y_6)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 = (x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{H IV}$$

$$\mathbf{B}_{12} = (x_6 + y_6)\,\mathbf{a}_1 + (x_6 - y_6)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{H IV}$$

$$\mathbf{B}_{13} = (x_7 - y_7)\,\mathbf{a}_1 + (x_7 + y_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 = (x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O I}$$

$$\mathbf{B}_{14} = (x_7 + y_7)\,\mathbf{a}_1 + (x_7 - y_7)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_7 a + z_7 c\cos\beta\right)\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_7\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O I}$$

$$\mathbf{B}_{15} = (x_8 - y_8)\,\mathbf{a}_1 + (x_8 + y_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 = (x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O II}$$

$$\mathbf{B}_{16} = (x_8 + y_8)\,\mathbf{a}_1 + (x_8 - y_8)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_8 a + z_8 c\cos\beta\right)\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_8\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O II}$$

$$\mathbf{B}_{17} = (x_9 - y_9)\,\mathbf{a}_1 + (x_9 + y_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 = (x_9 a + z_9 c\cos\beta)\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O III}$$

$$\mathbf{B}_{18} = (x_9 + y_9)\,\mathbf{a}_1 + (x_9 - y_9)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_9 a + z_9 c\cos\beta\right)\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_9\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O III}$$

$$\mathbf{B}_{19} = (x_{10} - y_{10})\,\mathbf{a}_1 + (x_{10} + y_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 = (x_{10} a + z_{10} c\cos\beta)\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O IV}$$

$$\mathbf{B}_{20} = (x_{10} + y_{10})\,\mathbf{a}_1 + (x_{10} - y_{10})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{10} a + z_{10} c\cos\beta\right)\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{10}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O IV}$$

$$\mathbf{B}_{21} = (x_{11} - y_{11})\,\mathbf{a}_1 + (x_{11} + y_{11})\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 = (x_{11} a + z_{11} c\cos\beta)\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + z_{11} c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O V}$$

$$\mathbf{B}_{22} = (x_{11} + y_{11})\,\mathbf{a}_1 + (x_{11} - y_{11})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{11} a + z_{11} c\cos\beta\right)\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{11}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O V}$$

$\mathbf{B}_{23}$	=	$(x_{12} - y_{12})\ \mathbf{a}_1 + (x_{12} + y_{12})\ \mathbf{a}_2 + z_{12}\ \mathbf{a}_3$	=	$(x_{12}a + z_{12}c\cos\beta)\ \hat{\mathbf{x}} + y_{12}b\ \hat{\mathbf{y}} + z_{12}c\sin\beta\ \hat{\mathbf{z}}$	(4*a*)	O VI
$\mathbf{B}_{24}$	=	$(x_{12} + y_{12})\ \mathbf{a}_1 + (x_{12} - y_{12})\ \mathbf{a}_2 + \left(\frac{1}{2} + z_{12}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{12}a + z_{12}c\cos\beta\right)\hat{\mathbf{x}} - y_{12}b\ \hat{\mathbf{y}} + \left(\frac{1}{2} + z_{12}\right)c\sin\beta\ \hat{\mathbf{z}}$	(4*a*)	O VI
$\mathbf{B}_{25}$	=	$(x_{13} - y_{13})\ \mathbf{a}_1 + (x_{13} + y_{13})\ \mathbf{a}_2 + z_{13}\ \mathbf{a}_3$	=	$(x_{13}a + z_{13}c\cos\beta)\ \hat{\mathbf{x}} + y_{13}b\ \hat{\mathbf{y}} + z_{13}c\sin\beta\ \hat{\mathbf{z}}$	(4*a*)	O VII
$\mathbf{B}_{26}$	=	$(x_{13} + y_{13})\ \mathbf{a}_1 + (x_{13} - y_{13})\ \mathbf{a}_2 + \left(\frac{1}{2} + z_{13}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{13}a + z_{13}c\cos\beta\right)\hat{\mathbf{x}} - y_{13}b\ \hat{\mathbf{y}} + \left(\frac{1}{2} + z_{13}\right)c\sin\beta\ \hat{\mathbf{z}}$	(4*a*)	O VII
$\mathbf{B}_{27}$	=	$(x_{14} - y_{14})\ \mathbf{a}_1 + (x_{14} + y_{14})\ \mathbf{a}_2 + z_{14}\ \mathbf{a}_3$	=	$(x_{14}a + z_{14}c\cos\beta)\ \hat{\mathbf{x}} + y_{14}b\ \hat{\mathbf{y}} + z_{14}c\sin\beta\ \hat{\mathbf{z}}$	(4*a*)	O VIII
$\mathbf{B}_{28}$	=	$(x_{14} + y_{14})\ \mathbf{a}_1 + (x_{14} - y_{14})\ \mathbf{a}_2 + \left(\frac{1}{2} + z_{14}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{14}a + z_{14}c\cos\beta\right)\hat{\mathbf{x}} - y_{14}b\ \hat{\mathbf{y}} + \left(\frac{1}{2} + z_{14}\right)c\sin\beta\ \hat{\mathbf{z}}$	(4*a*)	O VIII
$\mathbf{B}_{29}$	=	$(x_{15} - y_{15})\ \mathbf{a}_1 + (x_{15} + y_{15})\ \mathbf{a}_2 + z_{15}\ \mathbf{a}_3$	=	$(x_{15}a + z_{15}c\cos\beta)\ \hat{\mathbf{x}} + y_{15}b\ \hat{\mathbf{y}} + z_{15}c\sin\beta\ \hat{\mathbf{z}}$	(4*a*)	O IX
$\mathbf{B}_{30}$	=	$(x_{15} + y_{15})\ \mathbf{a}_1 + (x_{15} - y_{15})\ \mathbf{a}_2 + \left(\frac{1}{2} + z_{15}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{15}a + z_{15}c\cos\beta\right)\hat{\mathbf{x}} - y_{15}b\ \hat{\mathbf{y}} + \left(\frac{1}{2} + z_{15}\right)c\sin\beta\ \hat{\mathbf{z}}$	(4*a*)	O IX
$\mathbf{B}_{31}$	=	$(x_{16} - y_{16})\ \mathbf{a}_1 + (x_{16} + y_{16})\ \mathbf{a}_2 + z_{16}\ \mathbf{a}_3$	=	$(x_{16}a + z_{16}c\cos\beta)\ \hat{\mathbf{x}} + y_{16}b\ \hat{\mathbf{y}} + z_{16}c\sin\beta\ \hat{\mathbf{z}}$	(4*a*)	Si I
$\mathbf{B}_{32}$	=	$(x_{16} + y_{16})\ \mathbf{a}_1 + (x_{16} - y_{16})\ \mathbf{a}_2 + \left(\frac{1}{2} + z_{16}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{16}a + z_{16}c\cos\beta\right)\hat{\mathbf{x}} - y_{16}b\ \hat{\mathbf{y}} + \left(\frac{1}{2} + z_{16}\right)c\sin\beta\ \hat{\mathbf{z}}$	(4*a*)	Si I
$\mathbf{B}_{33}$	=	$(x_{17} - y_{17})\ \mathbf{a}_1 + (x_{17} + y_{17})\ \mathbf{a}_2 + z_{17}\ \mathbf{a}_3$	=	$(x_{17}a + z_{17}c\cos\beta)\ \hat{\mathbf{x}} + y_{17}b\ \hat{\mathbf{y}} + z_{17}c\sin\beta\ \hat{\mathbf{z}}$	(4*a*)	Si II
$\mathbf{B}_{34}$	=	$(x_{17} + y_{17})\ \mathbf{a}_1 + (x_{17} - y_{17})\ \mathbf{a}_2 + \left(\frac{1}{2} + z_{17}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{17}a + z_{17}c\cos\beta\right)\hat{\mathbf{x}} - y_{17}b\ \hat{\mathbf{y}} + \left(\frac{1}{2} + z_{17}\right)c\sin\beta\ \hat{\mathbf{z}}$	(4*a*)	Si II

References:

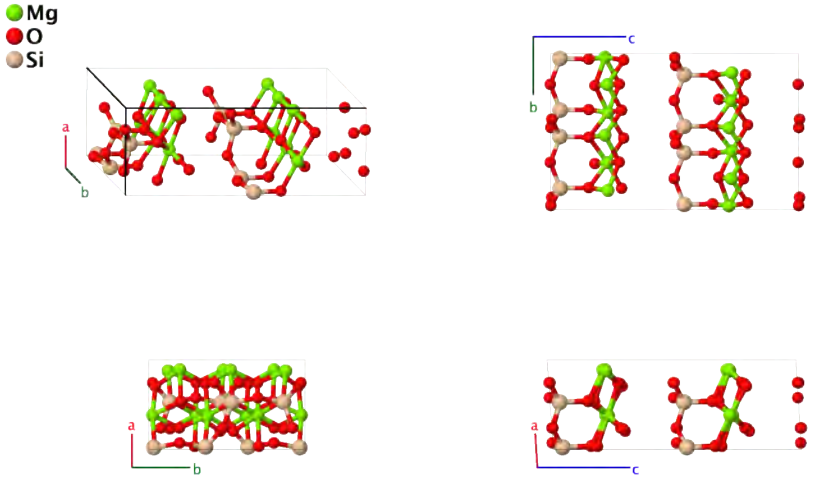
- A. P. Zhukhlistov, *Crystal structure of nacrite from the electron diffraction data*, Crystal. Rep. **53**, 76–82 (2008), doi:10.1134/S1063774508010094.
- S. B. Hendricks, *The Crystal Structure of Nacrite $Al_2O_3{\cdot}2(SiO_2){\cdot}2H_2O$ and the Polymorphism of the Kaolin Minerals*, Zeitschrift für Kristallographie - Crystalline Materials **100**, 509–518 (1939), doi:10.1524/zkri.1939.100.1.509.
- K. Herrmann, ed., *Strukturbericht Band VII 1939* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1943).

Geometry files:

- CIF: pp. 1518
- POSCAR: pp. 1518

Chrysotile ($Mg_3Si_2O_5(OH)_4$) Structure: A3B5C4D2_mC56_9_3a_5a_4a_2a

http://aflow.org/prototype-encyclopedia/A3B5C4D2_mC56_9_3a_5a_4a_2a

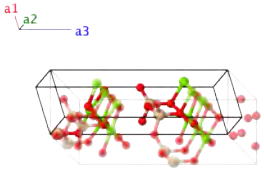


Prototype	:	$Mg_3O_5(OH)_4Si_2$
AFLOW prototype label	:	A3B5C4D2_mC56_9_3a_5a_4a_2a
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC56
Space group number	:	9
Space group symbol	:	Cc
AFLOW prototype command	:	`aflow --proto=A3B5C4D2_mC56_9_3a_5a_4a_2a --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}$

- Chrysotile is more commonly known as "white asbestos." Chrysotile sheets typically curl into tubular fibers and the crystal structure is difficult to determine. (Yada, 1967) has a partial list of the experiments performed to determine this structure. This differs from the early work of (Warren, 1930), $S4_5$, which halves the size of the unit cell and removes the inversion site. For this prototype, the OH molecules are centered on four a Wyckoff positions.

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$\mathbf{B}_1$	=	$(x_1 - y_1)\,\mathbf{a}_1 + (x_1 + y_1)\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\sin\beta\,\hat{\mathbf{z}}$	(4*a*)	Mg I
$\mathbf{B}_2$	=	$(x_1 + y_1)\,\mathbf{a}_1 + (x_1 - y_1)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_1 a + z_1 c\cos\beta\right)\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	(4*a*)	Mg I
$\mathbf{B}_3$	=	$(x_2 - y_2)\,\mathbf{a}_1 + (x_2 + y_2)\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	(4*a*)	Mg II
$\mathbf{B}_4$	=	$(x_2 + y_2)\,\mathbf{a}_1 + (x_2 - y_2)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_2 a + z_2 c\cos\beta\right)\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	(4*a*)	Mg II
$\mathbf{B}_5$	=	$(x_3 - y_3)\,\mathbf{a}_1 + (x_3 + y_3)\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	(4*a*)	Mg III
$\mathbf{B}_6$	=	$(x_3 + y_3)\,\mathbf{a}_1 + (x_3 - y_3)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	(4*a*)	Mg III
$\mathbf{B}_7$	=	$(x_4 - y_4)\,\mathbf{a}_1 + (x_4 + y_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_8$	=	$(x_4 + y_4)\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_9$	=	$(x_5 - y_5)\,\mathbf{a}_1 + (x_5 + y_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	(4*a*)	O II
$\mathbf{B}_{10}$	=	$(x_5 + y_5)\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	(4*a*)	O II
$\mathbf{B}_{11}$	=	$(x_6 - y_6)\,\mathbf{a}_1 + (x_6 + y_6)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	(4*a*)	O III
$\mathbf{B}_{12}$	=	$(x_6 + y_6)\,\mathbf{a}_1 + (x_6 - y_6)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	(4*a*)	O III
$\mathbf{B}_{13}$	=	$(x_7 - y_7)\,\mathbf{a}_1 + (x_7 + y_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	(4*a*)	O IV
$\mathbf{B}_{14}$	=	$(x_7 + y_7)\,\mathbf{a}_1 + (x_7 - y_7)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_7 a + z_7 c\cos\beta\right)\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right)c\sin\beta\,\hat{\mathbf{z}}$	(4*a*)	O IV
$\mathbf{B}_{15}$	=	$(x_8 - y_8)\,\mathbf{a}_1 + (x_8 + y_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}}$	(4*a*)	O V
$\mathbf{B}_{16}$	=	$(x_8 + y_8)\,\mathbf{a}_1 + (x_8 - y_8)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_8 a + z_8 c\cos\beta\right)\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right)c\sin\beta\,\hat{\mathbf{z}}$	(4*a*)	O V
$\mathbf{B}_{17}$	=	$(x_9 - y_9)\,\mathbf{a}_1 + (x_9 + y_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$(x_9 a + z_9 c\cos\beta)\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\sin\beta\,\hat{\mathbf{z}}$	(4*a*)	OH I
$\mathbf{B}_{18}$	=	$(x_9 + y_9)\,\mathbf{a}_1 + (x_9 - y_9)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_9 a + z_9 c\cos\beta\right)\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right)c\sin\beta\,\hat{\mathbf{z}}$	(4*a*)	OH I
$\mathbf{B}_{19}$	=	$(x_{10} - y_{10})\,\mathbf{a}_1 + (x_{10} + y_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$(x_{10} a + z_{10} c\cos\beta)\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\sin\beta\,\hat{\mathbf{z}}$	(4*a*)	OH II
$\mathbf{B}_{20}$	=	$(x_{10} + y_{10})\,\mathbf{a}_1 + (x_{10} - y_{10})\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{10}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{10} a + z_{10} c\cos\beta\right)\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{10}\right)c\sin\beta\,\hat{\mathbf{z}}$	(4*a*)	OH II
$\mathbf{B}_{21}$	=	$(x_{11} - y_{11})\,\mathbf{a}_1 + (x_{11} + y_{11})\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$(x_{11} a + z_{11} c\cos\beta)\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + z_{11} c\sin\beta\,\hat{\mathbf{z}}$	(4*a*)	OH III
$\mathbf{B}_{22}$	=	$(x_{11} + y_{11})\,\mathbf{a}_1 + (x_{11} - y_{11})\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{11}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{11} a + z_{11} c\cos\beta\right)\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{11}\right)c\sin\beta\,\hat{\mathbf{z}}$	(4*a*)	OH III

$\mathbf{B}_{23}$	=	$(x_{12} - y_{12})\ \mathbf{a}_1 + (x_{12} + y_{12})\ \mathbf{a}_2 + z_{12}\ \mathbf{a}_3$	=	$(x_{12}a + z_{12}c\cos\beta)\ \hat{\mathbf{x}} + y_{12}b\,\hat{\mathbf{y}} + z_{12}c\sin\beta\,\hat{\mathbf{z}}$	(4*a*)	OH IV
$\mathbf{B}_{24}$	=	$(x_{12} + y_{12})\ \mathbf{a}_1 + (x_{12} - y_{12})\ \mathbf{a}_2 + \left(\frac{1}{2} + z_{12}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{12}a + z_{12}c\cos\beta\right)\hat{\mathbf{x}} - y_{12}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{12}\right)c\sin\beta\,\hat{\mathbf{z}}$	(4*a*)	OH IV
$\mathbf{B}_{25}$	=	$(x_{13} - y_{13})\ \mathbf{a}_1 + (x_{13} + y_{13})\ \mathbf{a}_2 + z_{13}\ \mathbf{a}_3$	=	$(x_{13}a + z_{13}c\cos\beta)\ \hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} + z_{13}c\sin\beta\,\hat{\mathbf{z}}$	(4*a*)	Si I
$\mathbf{B}_{26}$	=	$(x_{13} + y_{13})\ \mathbf{a}_1 + (x_{13} - y_{13})\ \mathbf{a}_2 + \left(\frac{1}{2} + z_{13}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{13}a + z_{13}c\cos\beta\right)\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{13}\right)c\sin\beta\,\hat{\mathbf{z}}$	(4*a*)	Si I
$\mathbf{B}_{27}$	=	$(x_{14} - y_{14})\ \mathbf{a}_1 + (x_{14} + y_{14})\ \mathbf{a}_2 + z_{14}\ \mathbf{a}_3$	=	$(x_{14}a + z_{14}c\cos\beta)\ \hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} + z_{14}c\sin\beta\,\hat{\mathbf{z}}$	(4*a*)	Si II
$\mathbf{B}_{28}$	=	$(x_{14} + y_{14})\ \mathbf{a}_1 + (x_{14} - y_{14})\ \mathbf{a}_2 + \left(\frac{1}{2} + z_{14}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{14}a + z_{14}c\cos\beta\right)\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{14}\right)c\sin\beta\,\hat{\mathbf{z}}$	(4*a*)	Si II

References:

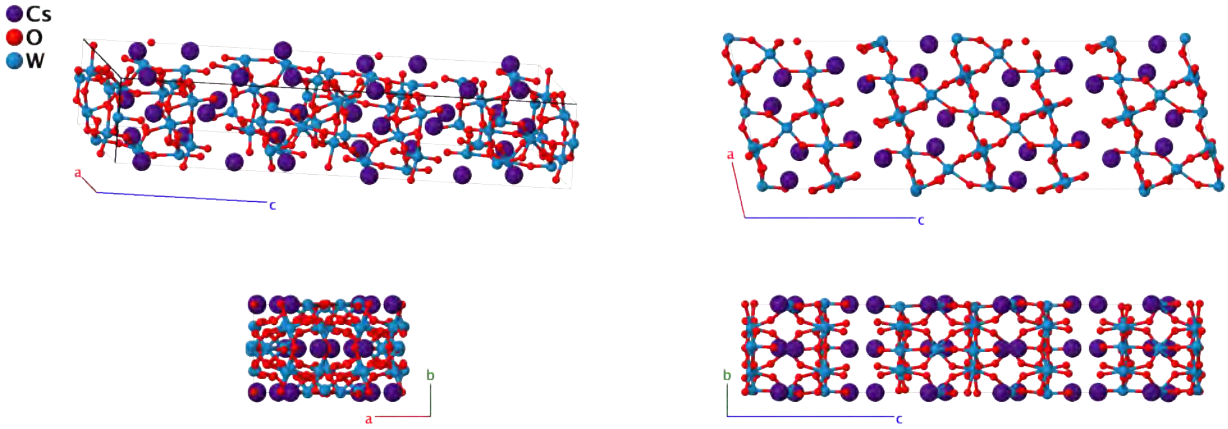
- G. Falini, E. Foresti, M. Gazzano, A. F. Gualtieri, M. Leoni, I. G. Lesci, and N. Roveri, *Tubular-Shaped Stoichiometric Chrysotile Nanocrystals*, Chem.: Eur. J. **10**, 3043–3049 (2004), doi:10.1002/chem.200305685.
- K. Yada, *Study of Chrysotile Asbestos by a High Resolution Electron Microscope*, Acta Cryst. **23**, 704–707 (1967), doi:10.1107/S0365110X67003524.

Geometry files:

- CIF: pp. 1518
- POSCAR: pp. 1519

$Cs_6W_{11}O_{36}$ Structure: A6B36C11_mC212_9_6a_36a_11a

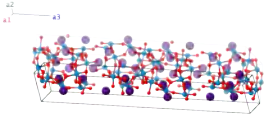
http://aflow.org/prototype-encyclopedia/A6B36C11_mC212_9_6a_36a_11a



Prototype	:	$Cs_6O_{36}W_{11}$
AFLOW prototype label	:	A6B36C11_mC212_9_6a_36a_11a
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC212
Space group number	:	9
Space group symbol	:	Cc
AFLOW prototype command	:	`aflow --proto=A6B36C11_mC212_9_6a_36a_11a --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}, x_{17}, y_{17}, z_{17}, x_{18}, y_{18}, z_{18}, x_{19}, y_{19}, z_{19}, x_{20}, y_{20}, z_{20}, x_{21}, y_{21}, z_{21}, x_{22}, y_{22}, z_{22}, x_{23}, y_{23}, z_{23}, x_{24}, y_{24}, z_{24}, x_{25}, y_{25}, z_{25}, x_{26}, y_{26}, z_{26}, x_{27}, y_{27}, z_{27}, x_{28}, y_{28}, z_{28}, x_{29}, y_{29}, z_{29}, x_{30}, y_{30}, z_{30}, x_{31}, y_{31}, z_{31}, x_{32}, y_{32}, z_{32}, x_{33}, y_{33}, z_{33}, x_{34}, y_{34}, z_{34}, x_{35}, y_{35}, z_{35}, x_{36}, y_{36}, z_{36}, x_{37}, y_{37}, z_{37}, x_{38}, y_{38}, z_{38}, x_{39}, y_{39}, z_{39}, x_{40}, y_{40}, z_{40}, x_{41}, y_{41}, z_{41}, x_{42}, y_{42}, z_{42}, x_{43}, y_{43}, z_{43}, x_{44}, y_{44}, z_{44}, x_{45}, y_{45}, z_{45}, x_{46}, y_{46}, z_{46}, x_{47}, y_{47}, z_{47}, x_{48}, y_{48}, z_{48}, x_{49}, y_{49}, z_{49}, x_{50}, y_{50}, z_{50}, x_{51}, y_{51}, z_{51}, x_{52}, y_{52}, z_{52}, x_{53}, y_{53}, z_{53}$

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$(x_1 - y_1)\,\mathbf{a}_1 + (x_1 + y_1)\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(4a)$	Cs I
$\mathbf{B}_2$	=	$(x_1 + y_1)\,\mathbf{a}_1 + (x_1 - y_1)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_1 a + z_1 c\cos\beta\right)\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4a)$	Cs I
$\mathbf{B}_3$	=	$(x_2 - y_2)\,\mathbf{a}_1 + (x_2 + y_2)\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(4a)$	Cs II

$$\mathbf{B}_4 = (x_2+y_2)\,\mathbf{a}_1+(x_2-y_2)\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_2\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta+x_2a+z_2c\cos\beta\right)\hat{\mathbf{x}}-y_2b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_2\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{Cs II}$$

$$\mathbf{B}_5 = (x_3-y_3)\,\mathbf{a}_1+(x_3+y_3)\,\mathbf{a}_2+z_3\,\mathbf{a}_3 = (x_3a+z_3c\cos\beta)\,\hat{\mathbf{x}}+y_3b\,\hat{\mathbf{y}}+z_3c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{Cs III}$$

$$\mathbf{B}_6 = (x_3+y_3)\,\mathbf{a}_1+(x_3-y_3)\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta+x_3a+z_3c\cos\beta\right)\hat{\mathbf{x}}-y_3b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_3\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{Cs III}$$

$$\mathbf{B}_7 = (x_4-y_4)\,\mathbf{a}_1+(x_4+y_4)\,\mathbf{a}_2+z_4\,\mathbf{a}_3 = (x_4a+z_4c\cos\beta)\,\hat{\mathbf{x}}+y_4b\,\hat{\mathbf{y}}+z_4c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{Cs IV}$$

$$\mathbf{B}_8 = (x_4+y_4)\,\mathbf{a}_1+(x_4-y_4)\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta+x_4a+z_4c\cos\beta\right)\hat{\mathbf{x}}-y_4b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_4\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{Cs IV}$$

$$\mathbf{B}_9 = (x_5-y_5)\,\mathbf{a}_1+(x_5+y_5)\,\mathbf{a}_2+z_5\,\mathbf{a}_3 = (x_5a+z_5c\cos\beta)\,\hat{\mathbf{x}}+y_5b\,\hat{\mathbf{y}}+z_5c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{Cs V}$$

$$\mathbf{B}_{10} = (x_5+y_5)\,\mathbf{a}_1+(x_5-y_5)\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta+x_5a+z_5c\cos\beta\right)\hat{\mathbf{x}}-y_5b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_5\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{Cs V}$$

$$\mathbf{B}_{11} = (x_6-y_6)\,\mathbf{a}_1+(x_6+y_6)\,\mathbf{a}_2+z_6\,\mathbf{a}_3 = (x_6a+z_6c\cos\beta)\,\hat{\mathbf{x}}+y_6b\,\hat{\mathbf{y}}+z_6c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{Cs VI}$$

$$\mathbf{B}_{12} = (x_6+y_6)\,\mathbf{a}_1+(x_6-y_6)\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta+x_6a+z_6c\cos\beta\right)\hat{\mathbf{x}}-y_6b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_6\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{Cs VI}$$

$$\mathbf{B}_{13} = (x_7-y_7)\,\mathbf{a}_1+(x_7+y_7)\,\mathbf{a}_2+z_7\,\mathbf{a}_3 = (x_7a+z_7c\cos\beta)\,\hat{\mathbf{x}}+y_7b\,\hat{\mathbf{y}}+z_7c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O I}$$

$$\mathbf{B}_{14} = (x_7+y_7)\,\mathbf{a}_1+(x_7-y_7)\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta+x_7a+z_7c\cos\beta\right)\hat{\mathbf{x}}-y_7b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_7\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O I}$$

$$\mathbf{B}_{15} = (x_8-y_8)\,\mathbf{a}_1+(x_8+y_8)\,\mathbf{a}_2+z_8\,\mathbf{a}_3 = (x_8a+z_8c\cos\beta)\,\hat{\mathbf{x}}+y_8b\,\hat{\mathbf{y}}+z_8c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O II}$$

$$\mathbf{B}_{16} = (x_8+y_8)\,\mathbf{a}_1+(x_8-y_8)\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta+x_8a+z_8c\cos\beta\right)\hat{\mathbf{x}}-y_8b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_8\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O II}$$

$$\mathbf{B}_{17} = (x_9-y_9)\,\mathbf{a}_1+(x_9+y_9)\,\mathbf{a}_2+z_9\,\mathbf{a}_3 = (x_9a+z_9c\cos\beta)\,\hat{\mathbf{x}}+y_9b\,\hat{\mathbf{y}}+z_9c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O III}$$

$$\mathbf{B}_{18} = (x_9+y_9)\,\mathbf{a}_1+(x_9-y_9)\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta+x_9a+z_9c\cos\beta\right)\hat{\mathbf{x}}-y_9b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_9\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O III}$$

$$\mathbf{B}_{19} = (x_{10}-y_{10})\,\mathbf{a}_1+(x_{10}+y_{10})\,\mathbf{a}_2+z_{10}\,\mathbf{a}_3 = (x_{10}a+z_{10}c\cos\beta)\,\hat{\mathbf{x}}+y_{10}b\,\hat{\mathbf{y}}+z_{10}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O IV}$$

$$\mathbf{B}_{20} = (x_{10}+y_{10})\,\mathbf{a}_1+(x_{10}-y_{10})\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta+x_{10}a+z_{10}c\cos\beta\right)\hat{\mathbf{x}}-y_{10}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{10}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O IV}$$

$$\mathbf{B}_{21} = (x_{11}-y_{11})\,\mathbf{a}_1+(x_{11}+y_{11})\,\mathbf{a}_2+z_{11}\,\mathbf{a}_3 = (x_{11}a+z_{11}c\cos\beta)\,\hat{\mathbf{x}}+y_{11}b\,\hat{\mathbf{y}}+z_{11}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O V}$$

$$\mathbf{B}_{22} = (x_{11}+y_{11})\,\mathbf{a}_1+(x_{11}-y_{11})\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta+x_{11}a+z_{11}c\cos\beta\right)\hat{\mathbf{x}}-y_{11}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{11}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O V}$$

$$\mathbf{B}_{23} = (x_{12}-y_{12})\,\mathbf{a}_1+(x_{12}+y_{12})\,\mathbf{a}_2+z_{12}\,\mathbf{a}_3 = (x_{12}a+z_{12}c\cos\beta)\,\hat{\mathbf{x}}+y_{12}b\,\hat{\mathbf{y}}+z_{12}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O VI}$$

$$\mathbf{B}_{24} = (x_{12}+y_{12})\,\mathbf{a}_1+(x_{12}-y_{12})\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta+x_{12}a+z_{12}c\cos\beta\right)\hat{\mathbf{x}}-y_{12}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{12}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O VI}$$

$$\mathbf{B}_{25} = (x_{13}-y_{13})\,\mathbf{a}_1+(x_{13}+y_{13})\,\mathbf{a}_2+z_{13}\,\mathbf{a}_3 = (x_{13}a+z_{13}c\cos\beta)\,\hat{\mathbf{x}}+y_{13}b\,\hat{\mathbf{y}}+z_{13}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O VII}$$

$$\mathbf{B}_{26} = (x_{13}+y_{13})\,\mathbf{a}_1 + (x_{13}-y_{13})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{13}a + z_{13}c\cos\beta\right)\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{13}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O VII}$$

$$\mathbf{B}_{27} = (x_{14}-y_{14})\,\mathbf{a}_1 + (x_{14}+y_{14})\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3 = (x_{14}a + z_{14}c\cos\beta)\,\hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} + z_{14}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O VIII}$$

$$\mathbf{B}_{28} = (x_{14}+y_{14})\,\mathbf{a}_1 + (x_{14}-y_{14})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{14}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{14}a + z_{14}c\cos\beta\right)\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{14}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O VIII}$$

$$\mathbf{B}_{29} = (x_{15}-y_{15})\,\mathbf{a}_1 + (x_{15}+y_{15})\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 = (x_{15}a + z_{15}c\cos\beta)\,\hat{\mathbf{x}} + y_{15}b\,\hat{\mathbf{y}} + z_{15}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O IX}$$

$$\mathbf{B}_{30} = (x_{15}+y_{15})\,\mathbf{a}_1 + (x_{15}-y_{15})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{15}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{15}a + z_{15}c\cos\beta\right)\hat{\mathbf{x}} - y_{15}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{15}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O IX}$$

$$\mathbf{B}_{31} = (x_{16}-y_{16})\,\mathbf{a}_1 + (x_{16}+y_{16})\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 = (x_{16}a + z_{16}c\cos\beta)\,\hat{\mathbf{x}} + y_{16}b\,\hat{\mathbf{y}} + z_{16}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O X}$$

$$\mathbf{B}_{32} = (x_{16}+y_{16})\,\mathbf{a}_1 + (x_{16}-y_{16})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{16}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{16}a + z_{16}c\cos\beta\right)\hat{\mathbf{x}} - y_{16}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{16}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O X}$$

$$\mathbf{B}_{33} = (x_{17}-y_{17})\,\mathbf{a}_1 + (x_{17}+y_{17})\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 = (x_{17}a + z_{17}c\cos\beta)\,\hat{\mathbf{x}} + y_{17}b\,\hat{\mathbf{y}} + z_{17}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XI}$$

$$\mathbf{B}_{34} = (x_{17}+y_{17})\,\mathbf{a}_1 + (x_{17}-y_{17})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{17}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{17}a + z_{17}c\cos\beta\right)\hat{\mathbf{x}} - y_{17}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{17}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XI}$$

$$\mathbf{B}_{35} = (x_{18}-y_{18})\,\mathbf{a}_1 + (x_{18}+y_{18})\,\mathbf{a}_2 + z_{18}\,\mathbf{a}_3 = (x_{18}a + z_{18}c\cos\beta)\,\hat{\mathbf{x}} + y_{18}b\,\hat{\mathbf{y}} + z_{18}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XII}$$

$$\mathbf{B}_{36} = (x_{18}+y_{18})\,\mathbf{a}_1 + (x_{18}-y_{18})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{18}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{18}a + z_{18}c\cos\beta\right)\hat{\mathbf{x}} - y_{18}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{18}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XII}$$

$$\mathbf{B}_{37} = (x_{19}-y_{19})\,\mathbf{a}_1 + (x_{19}+y_{19})\,\mathbf{a}_2 + z_{19}\,\mathbf{a}_3 = (x_{19}a + z_{19}c\cos\beta)\,\hat{\mathbf{x}} + y_{19}b\,\hat{\mathbf{y}} + z_{19}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XIII}$$

$$\mathbf{B}_{38} = (x_{19}+y_{19})\,\mathbf{a}_1 + (x_{19}-y_{19})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{19}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{19}a + z_{19}c\cos\beta\right)\hat{\mathbf{x}} - y_{19}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{19}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XIII}$$

$$\mathbf{B}_{39} = (x_{20}-y_{20})\,\mathbf{a}_1 + (x_{20}+y_{20})\,\mathbf{a}_2 + z_{20}\,\mathbf{a}_3 = (x_{20}a + z_{20}c\cos\beta)\,\hat{\mathbf{x}} + y_{20}b\,\hat{\mathbf{y}} + z_{20}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XIV}$$

$$\mathbf{B}_{40} = (x_{20}+y_{20})\,\mathbf{a}_1 + (x_{20}-y_{20})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{20}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{20}a + z_{20}c\cos\beta\right)\hat{\mathbf{x}} - y_{20}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{20}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XIV}$$

$$\mathbf{B}_{41} = (x_{21}-y_{21})\,\mathbf{a}_1 + (x_{21}+y_{21})\,\mathbf{a}_2 + z_{21}\,\mathbf{a}_3 = (x_{21}a + z_{21}c\cos\beta)\,\hat{\mathbf{x}} + y_{21}b\,\hat{\mathbf{y}} + z_{21}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XV}$$

$$\mathbf{B}_{42} = (x_{21}+y_{21})\,\mathbf{a}_1 + (x_{21}-y_{21})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{21}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{21}a + z_{21}c\cos\beta\right)\hat{\mathbf{x}} - y_{21}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{21}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XV}$$

$$\mathbf{B}_{43} = (x_{22}-y_{22})\,\mathbf{a}_1 + (x_{22}+y_{22})\,\mathbf{a}_2 + z_{22}\,\mathbf{a}_3 = (x_{22}a + z_{22}c\cos\beta)\,\hat{\mathbf{x}} + y_{22}b\,\hat{\mathbf{y}} + z_{22}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XVI}$$

$$\mathbf{B}_{44} = (x_{22}+y_{22})\,\mathbf{a}_1 + (x_{22}-y_{22})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{22}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{22}a + z_{22}c\cos\beta\right)\hat{\mathbf{x}} - y_{22}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{22}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XVI}$$

$$\mathbf{B}_{45} = (x_{23}-y_{23})\,\mathbf{a}_1 + (x_{23}+y_{23})\,\mathbf{a}_2 + z_{23}\,\mathbf{a}_3 = (x_{23}a + z_{23}c\cos\beta)\,\hat{\mathbf{x}} + y_{23}b\,\hat{\mathbf{y}} + z_{23}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XVII}$$

$$\mathbf{B}_{46} = (x_{23}+y_{23})\,\mathbf{a}_1 + (x_{23}-y_{23})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{23}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{23}a + z_{23}c\cos\beta\right)\hat{\mathbf{x}} - y_{23}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{23}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XVII}$$

$$\mathbf{B}_{47} = (x_{24}-y_{24})\,\mathbf{a}_1 + (x_{24}+y_{24})\,\mathbf{a}_2 + z_{24}\,\mathbf{a}_3 = (x_{24}a + z_{24}c\cos\beta)\,\hat{\mathbf{x}} + y_{24}b\,\hat{\mathbf{y}} + z_{24}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XVIII}$$

$$\mathbf{B}_{48} = (x_{24}+y_{24})\,\mathbf{a}_1 + (x_{24}-y_{24})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{24}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{24}a + z_{24}c\cos\beta\right)\hat{\mathbf{x}} - y_{24}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{24}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XVIII}$$

$$\mathbf{B}_{49} = (x_{25}-y_{25})\,\mathbf{a}_1 + (x_{25}+y_{25})\,\mathbf{a}_2 + z_{25}\,\mathbf{a}_3 = (x_{25}a + z_{25}c\cos\beta)\,\hat{\mathbf{x}} + y_{25}b\,\hat{\mathbf{y}} + z_{25}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XIX}$$

$$\mathbf{B}_{50} = (x_{25}+y_{25})\,\mathbf{a}_1 + (x_{25}-y_{25})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{25}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{25}a + z_{25}c\cos\beta\right)\hat{\mathbf{x}} - y_{25}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{25}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XIX}$$

$$\mathbf{B}_{51} = (x_{26}-y_{26})\,\mathbf{a}_1 + (x_{26}+y_{26})\,\mathbf{a}_2 + z_{26}\,\mathbf{a}_3 = (x_{26}a + z_{26}c\cos\beta)\,\hat{\mathbf{x}} + y_{26}b\,\hat{\mathbf{y}} + z_{26}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XX}$$

$$\mathbf{B}_{52} = (x_{26}+y_{26})\,\mathbf{a}_1 + (x_{26}-y_{26})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{26}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{26}a + z_{26}c\cos\beta\right)\hat{\mathbf{x}} - y_{26}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{26}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XX}$$

$$\mathbf{B}_{53} = (x_{27}-y_{27})\,\mathbf{a}_1 + (x_{27}+y_{27})\,\mathbf{a}_2 + z_{27}\,\mathbf{a}_3 = (x_{27}a + z_{27}c\cos\beta)\,\hat{\mathbf{x}} + y_{27}b\,\hat{\mathbf{y}} + z_{27}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XXI}$$

$$\mathbf{B}_{54} = (x_{27}+y_{27})\,\mathbf{a}_1 + (x_{27}-y_{27})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{27}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{27}a + z_{27}c\cos\beta\right)\hat{\mathbf{x}} - y_{27}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{27}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XXI}$$

$$\mathbf{B}_{55} = (x_{28}-y_{28})\,\mathbf{a}_1 + (x_{28}+y_{28})\,\mathbf{a}_2 + z_{28}\,\mathbf{a}_3 = (x_{28}a + z_{28}c\cos\beta)\,\hat{\mathbf{x}} + y_{28}b\,\hat{\mathbf{y}} + z_{28}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XXII}$$

$$\mathbf{B}_{56} = (x_{28}+y_{28})\,\mathbf{a}_1 + (x_{28}-y_{28})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{28}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{28}a + z_{28}c\cos\beta\right)\hat{\mathbf{x}} - y_{28}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{28}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XXII}$$

$$\mathbf{B}_{57} = (x_{29}-y_{29})\,\mathbf{a}_1 + (x_{29}+y_{29})\,\mathbf{a}_2 + z_{29}\,\mathbf{a}_3 = (x_{29}a + z_{29}c\cos\beta)\,\hat{\mathbf{x}} + y_{29}b\,\hat{\mathbf{y}} + z_{29}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XXIII}$$

$$\mathbf{B}_{58} = (x_{29}+y_{29})\,\mathbf{a}_1 + (x_{29}-y_{29})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{29}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{29}a + z_{29}c\cos\beta\right)\hat{\mathbf{x}} - y_{29}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{29}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XXIII}$$

$$\mathbf{B}_{59} = (x_{30}-y_{30})\,\mathbf{a}_1 + (x_{30}+y_{30})\,\mathbf{a}_2 + z_{30}\,\mathbf{a}_3 = (x_{30}a + z_{30}c\cos\beta)\,\hat{\mathbf{x}} + y_{30}b\,\hat{\mathbf{y}} + z_{30}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XXIV}$$

$$\mathbf{B}_{60} = (x_{30}+y_{30})\,\mathbf{a}_1 + (x_{30}-y_{30})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{30}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{30}a + z_{30}c\cos\beta\right)\hat{\mathbf{x}} - y_{30}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{30}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XXIV}$$

$$\mathbf{B}_{61} = (x_{31}-y_{31})\,\mathbf{a}_1 + (x_{31}+y_{31})\,\mathbf{a}_2 + z_{31}\,\mathbf{a}_3 = (x_{31}a + z_{31}c\cos\beta)\,\hat{\mathbf{x}} + y_{31}b\,\hat{\mathbf{y}} + z_{31}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XXV}$$

$$\mathbf{B}_{62} = (x_{31}+y_{31})\,\mathbf{a}_1 + (x_{31}-y_{31})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{31}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{31}a + z_{31}c\cos\beta\right)\hat{\mathbf{x}} - y_{31}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{31}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XXV}$$

$$\mathbf{B}_{63} = (x_{32}-y_{32})\,\mathbf{a}_1 + (x_{32}+y_{32})\,\mathbf{a}_2 + z_{32}\,\mathbf{a}_3 = (x_{32}a + z_{32}c\cos\beta)\,\hat{\mathbf{x}} + y_{32}b\,\hat{\mathbf{y}} + z_{32}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XXVI}$$

$$\mathbf{B}_{64} = (x_{32}+y_{32})\,\mathbf{a}_1 + (x_{32}-y_{32})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{32}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{32}a + z_{32}c\cos\beta\right)\hat{\mathbf{x}} - y_{32}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{32}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XXVI}$$

$$\mathbf{B}_{65} = (x_{33}-y_{33})\,\mathbf{a}_1 + (x_{33}+y_{33})\,\mathbf{a}_2 + z_{33}\,\mathbf{a}_3 = (x_{33}a + z_{33}c\cos\beta)\,\hat{\mathbf{x}} + y_{33}b\,\hat{\mathbf{y}} + z_{33}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XXVII}$$

$$\mathbf{B}_{66} = (x_{33}+y_{33})\,\mathbf{a}_1 + (x_{33}-y_{33})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{33}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{33}a + z_{33}c\cos\beta\right)\hat{\mathbf{x}} - y_{33}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{33}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XXVII}$$

$$\mathbf{B}_{67} = (x_{34}-y_{34})\,\mathbf{a}_1 + (x_{34}+y_{34})\,\mathbf{a}_2 + z_{34}\,\mathbf{a}_3 = (x_{34}a + z_{34}c\cos\beta)\,\hat{\mathbf{x}} + y_{34}b\,\hat{\mathbf{y}} + z_{34}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XXVIII}$$

$$\mathbf{B}_{68} = (x_{34}+y_{34})\,\mathbf{a}_1 + (x_{34}-y_{34})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{34}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{34}a + z_{34}c\cos\beta\right)\hat{\mathbf{x}} - y_{34}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{34}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XXVIII}$$

$$\mathbf{B}_{69} = (x_{35}-y_{35})\,\mathbf{a}_1 + (x_{35}+y_{35})\,\mathbf{a}_2 + z_{35}\,\mathbf{a}_3 = (x_{35}a + z_{35}c\cos\beta)\,\hat{\mathbf{x}} + y_{35}b\,\hat{\mathbf{y}} + z_{35}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XXIX}$$

$$\mathbf{B}_{70} = (x_{35}+y_{35})\,\mathbf{a}_1 + (x_{35}-y_{35})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{35}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{35}a + z_{35}c\cos\beta\right)\hat{\mathbf{x}} - y_{35}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{35}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XXIX}$$

$$\mathbf{B}_{71} = (x_{36}-y_{36})\,\mathbf{a}_1 + (x_{36}+y_{36})\,\mathbf{a}_2 + z_{36}\,\mathbf{a}_3 = (x_{36}a + z_{36}c\cos\beta)\,\hat{\mathbf{x}} + y_{36}b\,\hat{\mathbf{y}} + z_{36}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XXX}$$

$$\mathbf{B}_{72} = (x_{36}+y_{36})\,\mathbf{a}_1 + (x_{36}-y_{36})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{36}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{36}a + z_{36}c\cos\beta\right)\hat{\mathbf{x}} - y_{36}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{36}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XXX}$$

$$\mathbf{B}_{73} = (x_{37}-y_{37})\,\mathbf{a}_1 + (x_{37}+y_{37})\,\mathbf{a}_2 + z_{37}\,\mathbf{a}_3 = (x_{37}a + z_{37}c\cos\beta)\,\hat{\mathbf{x}} + y_{37}b\,\hat{\mathbf{y}} + z_{37}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XXXI}$$

$$\mathbf{B}_{74} = (x_{37}+y_{37})\,\mathbf{a}_1 + (x_{37}-y_{37})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{37}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{37}a + z_{37}c\cos\beta\right)\hat{\mathbf{x}} - y_{37}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{37}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XXXI}$$

$$\mathbf{B}_{75} = (x_{38}-y_{38})\,\mathbf{a}_1 + (x_{38}+y_{38})\,\mathbf{a}_2 + z_{38}\,\mathbf{a}_3 = (x_{38}a + z_{38}c\cos\beta)\,\hat{\mathbf{x}} + y_{38}b\,\hat{\mathbf{y}} + z_{38}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XXXII}$$

$$\mathbf{B}_{76} = (x_{38}+y_{38})\,\mathbf{a}_1 + (x_{38}-y_{38})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{38}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{38}a + z_{38}c\cos\beta\right)\hat{\mathbf{x}} - y_{38}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{38}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XXXII}$$

$$\mathbf{B}_{77} = (x_{39}-y_{39})\,\mathbf{a}_1 + (x_{39}+y_{39})\,\mathbf{a}_2 + z_{39}\,\mathbf{a}_3 = (x_{39}a + z_{39}c\cos\beta)\,\hat{\mathbf{x}} + y_{39}b\,\hat{\mathbf{y}} + z_{39}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XXXIII}$$

$$\mathbf{B}_{78} = (x_{39}+y_{39})\,\mathbf{a}_1 + (x_{39}-y_{39})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{39}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{39}a + z_{39}c\cos\beta\right)\hat{\mathbf{x}} - y_{39}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{39}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XXXIII}$$

$$\mathbf{B}_{79} = (x_{40}-y_{40})\,\mathbf{a}_1 + (x_{40}+y_{40})\,\mathbf{a}_2 + z_{40}\,\mathbf{a}_3 = (x_{40}a + z_{40}c\cos\beta)\,\hat{\mathbf{x}} + y_{40}b\,\hat{\mathbf{y}} + z_{40}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XXXIV}$$

$$\mathbf{B}_{80} = (x_{40}+y_{40})\,\mathbf{a}_1 + (x_{40}-y_{40})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{40}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{40}a + z_{40}c\cos\beta\right)\hat{\mathbf{x}} - y_{40}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{40}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XXXIV}$$

$$\mathbf{B}_{81} = (x_{41}-y_{41})\,\mathbf{a}_1 + (x_{41}+y_{41})\,\mathbf{a}_2 + z_{41}\,\mathbf{a}_3 = (x_{41}a + z_{41}c\cos\beta)\,\hat{\mathbf{x}} + y_{41}b\,\hat{\mathbf{y}} + z_{41}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XXXV}$$

$$\mathbf{B}_{82} = (x_{41}+y_{41})\,\mathbf{a}_1 + (x_{41}-y_{41})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{41}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{41}a + z_{41}c\cos\beta\right)\hat{\mathbf{x}} - y_{41}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{41}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XXXV}$$

$$\mathbf{B}_{83} = (x_{42}-y_{42})\,\mathbf{a}_1 + (x_{42}+y_{42})\,\mathbf{a}_2 + z_{42}\,\mathbf{a}_3 = (x_{42}a + z_{42}c\cos\beta)\,\hat{\mathbf{x}} + y_{42}b\,\hat{\mathbf{y}} + z_{42}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XXXVI}$$

$$\mathbf{B}_{84} = (x_{42}+y_{42})\,\mathbf{a}_1 + (x_{42}-y_{42})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{42}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{42}a + z_{42}c\cos\beta\right)\hat{\mathbf{x}} - y_{42}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{42}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XXXVI}$$

$$\mathbf{B}_{85} = (x_{43}-y_{43})\,\mathbf{a}_1 + (x_{43}+y_{43})\,\mathbf{a}_2 + z_{43}\,\mathbf{a}_3 = (x_{43}a + z_{43}c\cos\beta)\,\hat{\mathbf{x}} + y_{43}b\,\hat{\mathbf{y}} + z_{43}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{W I}$$

$$\mathbf{B}_{86} = (x_{43}+y_{43})\,\mathbf{a}_1 + (x_{43}-y_{43})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{43}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{43}a + z_{43}c\cos\beta\right)\hat{\mathbf{x}} - y_{43}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{43}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{W I}$$

$$\mathbf{B}_{87} = (x_{44}-y_{44})\,\mathbf{a}_1 + (x_{44}+y_{44})\,\mathbf{a}_2 + z_{44}\,\mathbf{a}_3 = (x_{44}a + z_{44}c\cos\beta)\,\hat{\mathbf{x}} + y_{44}b\,\hat{\mathbf{y}} + z_{44}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{W II}$$

$$\mathbf{B}_{88} = (x_{44}+y_{44})\,\mathbf{a}_1 + (x_{44}-y_{44})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{44}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{44}a + z_{44}c\cos\beta\right)\hat{\mathbf{x}} - y_{44}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{44}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{W II}$$

$$\mathbf{B}_{89} = (x_{45}-y_{45})\,\mathbf{a}_1 + (x_{45}+y_{45})\,\mathbf{a}_2 + z_{45}\,\mathbf{a}_3 = (x_{45}a + z_{45}c\cos\beta)\,\hat{\mathbf{x}} + y_{45}b\,\hat{\mathbf{y}} + z_{45}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{W III}$$

$$\mathbf{B}_{90} = (x_{45}+y_{45})\,\mathbf{a}_1 + (x_{45}-y_{45})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{45}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{45}a + z_{45}c\cos\beta\right)\hat{\mathbf{x}} - y_{45}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{45}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{W III}$$

$$\mathbf{B}_{91} = (x_{46}-y_{46})\,\mathbf{a}_1 + (x_{46}+y_{46})\,\mathbf{a}_2 + z_{46}\,\mathbf{a}_3 = (x_{46}a + z_{46}c\cos\beta)\,\hat{\mathbf{x}} + y_{46}b\,\hat{\mathbf{y}} + z_{46}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{W IV}$$

$$\mathbf{B}_{92} = (x_{46}+y_{46})\,\mathbf{a}_1 + (x_{46}-y_{46})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{46}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{46}a + z_{46}c\cos\beta\right)\hat{\mathbf{x}} - y_{46}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{46}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{W IV}$$

$$\mathbf{B}_{93} = (x_{47}-y_{47})\,\mathbf{a}_1 + (x_{47}+y_{47})\,\mathbf{a}_2 + z_{47}\,\mathbf{a}_3 = (x_{47}a + z_{47}c\cos\beta)\,\hat{\mathbf{x}} + y_{47}b\,\hat{\mathbf{y}} + z_{47}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{W V}$$

$$\mathbf{B}_{94} = (x_{47}+y_{47})\,\mathbf{a}_1 + (x_{47}-y_{47})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{47}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{47}a + z_{47}c\cos\beta\right)\hat{\mathbf{x}} - y_{47}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{47}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{W V}$$

$$\mathbf{B}_{95} = (x_{48}-y_{48})\,\mathbf{a}_1 + (x_{48}+y_{48})\,\mathbf{a}_2 + z_{48}\,\mathbf{a}_3 = (x_{48}a + z_{48}c\cos\beta)\,\hat{\mathbf{x}} + y_{48}b\,\hat{\mathbf{y}} + z_{48}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{W VI}$$

$$\mathbf{B}_{96} = (x_{48}+y_{48})\,\mathbf{a}_1 + (x_{48}-y_{48})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{48}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{48}a + z_{48}c\cos\beta\right)\hat{\mathbf{x}} - y_{48}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{48}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{W VI}$$

$$\mathbf{B}_{97} = (x_{49}-y_{49})\,\mathbf{a}_1 + (x_{49}+y_{49})\,\mathbf{a}_2 + z_{49}\,\mathbf{a}_3 = (x_{49}a + z_{49}c\cos\beta)\,\hat{\mathbf{x}} + y_{49}b\,\hat{\mathbf{y}} + z_{49}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{W VII}$$

$$\mathbf{B}_{98} = (x_{49}+y_{49})\,\mathbf{a}_1 + (x_{49}-y_{49})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{49}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{49}a + z_{49}c\cos\beta\right)\hat{\mathbf{x}} - y_{49}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{49}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{W VII}$$

$$\mathbf{B}_{99} = (x_{50}-y_{50})\,\mathbf{a}_1 + (x_{50}+y_{50})\,\mathbf{a}_2 + z_{50}\,\mathbf{a}_3 = (x_{50}a + z_{50}c\cos\beta)\,\hat{\mathbf{x}} + y_{50}b\,\hat{\mathbf{y}} + z_{50}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{W VIII}$$

$$\mathbf{B}_{100} = (x_{50}+y_{50})\,\mathbf{a}_1 + (x_{50}-y_{50})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{50}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{50}a + z_{50}c\cos\beta\right)\hat{\mathbf{x}} - y_{50}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{50}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{W VIII}$$

$$\mathbf{B}_{101} = (x_{51}-y_{51})\,\mathbf{a}_1 + (x_{51}+y_{51})\,\mathbf{a}_2 + z_{51}\,\mathbf{a}_3 = (x_{51}a + z_{51}c\cos\beta)\,\hat{\mathbf{x}} + y_{51}b\,\hat{\mathbf{y}} + z_{51}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{W IX}$$

$$\mathbf{B}_{102} = (x_{51}+y_{51})\,\mathbf{a}_1 + (x_{51}-y_{51})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{51}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{51}a + z_{51}c\cos\beta\right)\hat{\mathbf{x}} - y_{51}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{51}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{W IX}$$

$$\mathbf{B}_{103} = (x_{52}-y_{52})\,\mathbf{a}_1 + (x_{52}+y_{52})\,\mathbf{a}_2 + z_{52}\,\mathbf{a}_3 = (x_{52}a + z_{52}c\cos\beta)\,\hat{\mathbf{x}} + y_{52}b\,\hat{\mathbf{y}} + z_{52}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{W X}$$

$$\mathbf{B}_{104} = (x_{52}+y_{52})\,\mathbf{a}_1 + (x_{52}-y_{52})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{52}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{52}a + z_{52}c\cos\beta\right)\hat{\mathbf{x}} - y_{52}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{52}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{W X}$$

$$\mathbf{B}_{105} = (x_{53}-y_{53})\,\mathbf{a}_1 + (x_{53}+y_{53})\,\mathbf{a}_2 + z_{53}\,\mathbf{a}_3 = (x_{53}a + z_{53}c\cos\beta)\,\hat{\mathbf{x}} + y_{53}b\,\hat{\mathbf{y}} + z_{53}c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{W XI}$$

$$\mathbf{B}_{106} = (x_{53}+y_{53})\,\mathbf{a}_1 + (x_{53}-y_{53})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{53}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{53}a + z_{53}c\cos\beta\right)\hat{\mathbf{x}} - y_{53}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{53}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4a) \quad \text{W XI}$$

References:

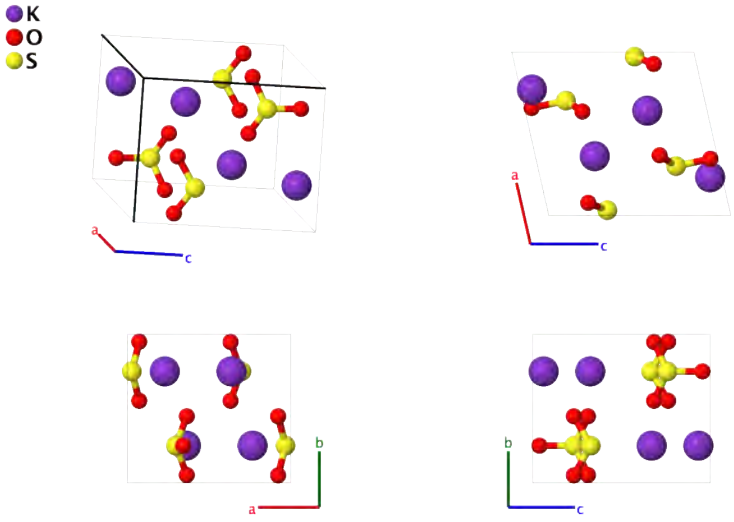
- K. Okada, F. Marumo, and S. Iwai, *The Crystal Structure of* $Cs_6W_{11}O_{36}$, Acta Crystallogr. Sect. B Struct. Sci. **34**, 50–54 (1978), doi:10.1107/S0567740878014934.

Geometry files:

- CIF: pp. 1519
- POSCAR: pp. 1520

$K_2S_2O_5$ ($K0_1$) Structure: A2B5C2_mP18_11_2e_e2f_2e

http://aflow.org/prototype-encyclopedia/A2B5C2_mP18_11_2e_e2f_2e



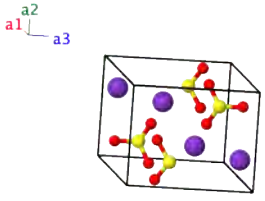
Prototype	:	$K_2O_5S_2$
AFLOW prototype label	:	A2B5C2_mP18_11_2e_e2f_2e
***Strukturbericht* designation**	:	$K0_1$
Pearson symbol	:	mP18
Space group number	:	11
Space group symbol	:	$P2_1/m$
AFLOW prototype command	:	`aflow --proto=A2B5C2_mP18_11_2e_e2f_2e --params=`$a, b/a, c/a, \beta, x_1, z_1, x_2, z_2, x_3, z_3, x_4, z_4, x_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7$

Other compounds with this structure

- $K_2S_2P_5$ and $(NH_4)_2S_2P_5$

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1a + z_1c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_1c\sin\beta\,\hat{\mathbf{z}}$	$(2e)$	K I
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$(-x_1a - z_1c\cos\beta)\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_1c\sin\beta\,\hat{\mathbf{z}}$	$(2e)$	K I
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2a + z_2c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_2c\sin\beta\,\hat{\mathbf{z}}$	$(2e)$	K II

$\mathbf{B}_4$	=	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(2e)$	K II
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(2e)$	O I
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(2e)$	O I
$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(2e)$	S I
$\mathbf{B}_8$	=	$-x_4\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(2e)$	S I
$\mathbf{B}_9$	=	$x_5\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(2e)$	S II
$\mathbf{B}_{10}$	=	$-x_5\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(2e)$	S II
$\mathbf{B}_{11}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(4f)$	O II
$\mathbf{B}_{12}$	=	$-x_6\,\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right) b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(4f)$	O II
$\mathbf{B}_{13}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(4f)$	O II
$\mathbf{B}_{14}$	=	$x_6\,\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right) b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(4f)$	O II
$\mathbf{B}_{15}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(4f)$	O III
$\mathbf{B}_{16}$	=	$-x_7\,\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right) b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(4f)$	O III
$\mathbf{B}_{17}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(4f)$	O III
$\mathbf{B}_{18}$	=	$x_7\,\mathbf{a}_1 + \left(\frac{1}{2} - y_7\right)\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_7\right) b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(4f)$	O III

References:

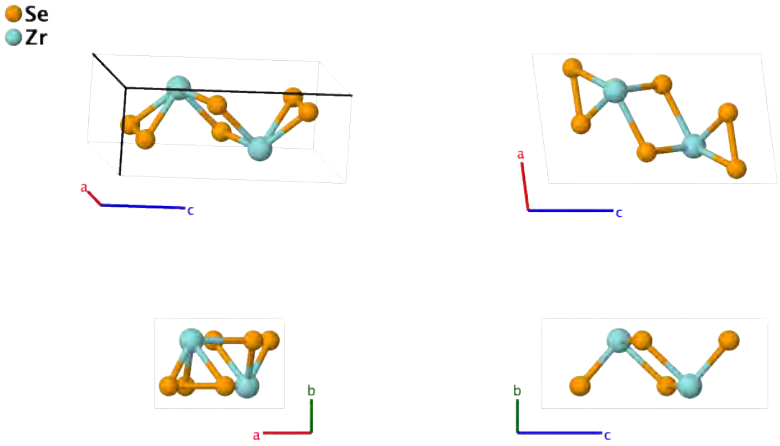
- I.-C. Chen and Y. Wang, *Reinvestigation of potassium pyrosulfite, $K_2S_2O_5$*, Acta Crystallogr. C **40**, 1780–1781 (1984), doi:10.1107/S0108270184009525.

Geometry files:

- CIF: pp. 1521
- POSCAR: pp. 1521

$ZrSe_3$ Structure: A3B_mP8_11_3e_e

http://aflow.org/prototype-encyclopedia/A3B_mP8_11_3e_e



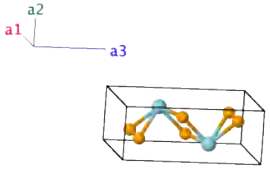
Prototype	:	Se_3Zr
AFLOW prototype label	:	A3B_mP8_11_3e_e
***Strukturbericht* designation**	:	None
Pearson symbol	:	mP8
Space group number	:	11
Space group symbol	:	$P2_1/m$
AFLOW prototype command	:	`aflow --proto=A3B_mP8_11_3e_e --params=`$a, b/a, c/a, \beta, x_1, z_1, x_2, z_2, x_3, z_3, x_4, z_4$

Other compounds with this structure

- HfS_3, $HfSe_3$, $HfTe_3$, $TaSe_3$, TiS_3, ZrS_3, and $ZrTe_3$

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(2e)$	Se I
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$(-x_1 a - z_1 c\cos\beta)\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(2e)$	Se I
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(2e)$	Se II
$\mathbf{B}_4$	=	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(2e)$	Se II
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(2e)$	Se III

$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	(2*e*)	Se III
$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	(2*e*)	Zr
$\mathbf{B}_8$	=	$-x_4\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	(2*e*)	Zr

References:

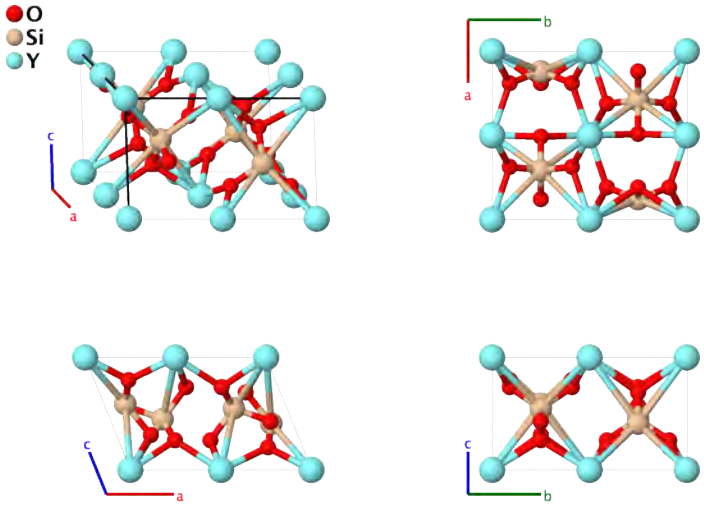
- S. Furuseth, L. Brattås, and A. Kjekshus, *On the Crystal Structures of TiS_3, ZrS_3, $ZrSe_3$, $ZrTe_3$, HfS_3, and $HfSe_3$*, Acta Chem. Scand. **29a**, 623–631 (1975), doi:10.3891/acta.chem.scand.29a-0623.

Geometry files:

- CIF: pp. 1521
- POSCAR: pp. 1521

y-$Y_2Si_2O_7$ Structure: A7B2C2_mP22_11_3e2f_2e_ab

http://aflow.org/prototype-encyclopedia/A7B2C2_mP22_11_3e2f_2e_ab

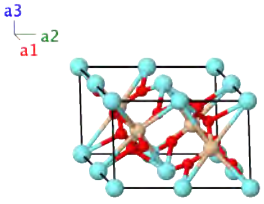


Prototype	:	$O_7Si_2Y_2$
AFLOW prototype label	:	A7B2C2_mP22_11_3e2f_2e_ab
***Strukturbericht* designation**	:	None
Pearson symbol	:	mP22
Space group number	:	11
Space group symbol	:	$P2_1/m$
AFLOW prototype command	:	`aflow --proto=A7B2C2_mP22_11_3e2f_2e_ab --params=`$a, b/a, c/a, \beta, x_3, z_3, x_4, z_4, x_5, z_5, x_6, z_6, x_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9$

- (Becerro, 2003) found that the NMR spectrum of y-$Y_2Si_2O_7$ showed only one peak for yttrium. If that is the case, then this crystal structure is incorrect.

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	Y I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}b\,\hat{\mathbf{y}}$	$(2a)$	Y I
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	$(2b)$	Y II

$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$	(2*b*)	Y II
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3a + z_3c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_3c\sin\beta\,\hat{\mathbf{z}}$	(2*e*)	O I
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3a - z_3c\cos\beta)\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_3c\sin\beta\,\hat{\mathbf{z}}$	(2*e*)	O I
$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4a + z_4c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_4c\sin\beta\,\hat{\mathbf{z}}$	(2*e*)	O II
$\mathbf{B}_8$	=	$-x_4\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4a - z_4c\cos\beta)\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_4c\sin\beta\,\hat{\mathbf{z}}$	(2*e*)	O II
$\mathbf{B}_9$	=	$x_5\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5a + z_5c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_5c\sin\beta\,\hat{\mathbf{z}}$	(2*e*)	O III
$\mathbf{B}_{10}$	=	$-x_5\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5a - z_5c\cos\beta)\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_5c\sin\beta\,\hat{\mathbf{z}}$	(2*e*)	O III
$\mathbf{B}_{11}$	=	$x_6\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6a + z_6c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_6c\sin\beta\,\hat{\mathbf{z}}$	(2*e*)	Si I
$\mathbf{B}_{12}$	=	$-x_6\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6a - z_6c\cos\beta)\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_6c\sin\beta\,\hat{\mathbf{z}}$	(2*e*)	Si I
$\mathbf{B}_{13}$	=	$x_7\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7a + z_7c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_7c\sin\beta\,\hat{\mathbf{z}}$	(2*e*)	Si II
$\mathbf{B}_{14}$	=	$-x_7\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7a - z_7c\cos\beta)\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_7c\sin\beta\,\hat{\mathbf{z}}$	(2*e*)	Si II
$\mathbf{B}_{15}$	=	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8a + z_8c\cos\beta)\,\hat{\mathbf{x}} + y_8b\,\hat{\mathbf{y}} + z_8c\sin\beta\,\hat{\mathbf{z}}$	(4*f*)	O IV
$\mathbf{B}_{16}$	=	$-x_8\,\mathbf{a}_1 + \left(\frac{1}{2} + y_8\right)\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8a - z_8c\cos\beta)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_8\right)b\,\hat{\mathbf{y}} - z_8c\sin\beta\,\hat{\mathbf{z}}$	(4*f*)	O IV
$\mathbf{B}_{17}$	=	$-x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8a - z_8c\cos\beta)\,\hat{\mathbf{x}} - y_8b\,\hat{\mathbf{y}} - z_8c\sin\beta\,\hat{\mathbf{z}}$	(4*f*)	O IV
$\mathbf{B}_{18}$	=	$x_8\,\mathbf{a}_1 + \left(\frac{1}{2} - y_8\right)\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8a + z_8c\cos\beta)\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_8\right)b\,\hat{\mathbf{y}} + z_8c\sin\beta\,\hat{\mathbf{z}}$	(4*f*)	O IV
$\mathbf{B}_{19}$	=	$x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$(x_9a + z_9c\cos\beta)\,\hat{\mathbf{x}} + y_9b\,\hat{\mathbf{y}} + z_9c\sin\beta\,\hat{\mathbf{z}}$	(4*f*)	O V
$\mathbf{B}_{20}$	=	$-x_9\,\mathbf{a}_1 + \left(\frac{1}{2} + y_9\right)\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$(-x_9a - z_9c\cos\beta)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_9\right)b\,\hat{\mathbf{y}} - z_9c\sin\beta\,\hat{\mathbf{z}}$	(4*f*)	O V
$\mathbf{B}_{21}$	=	$-x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$(-x_9a - z_9c\cos\beta)\,\hat{\mathbf{x}} - y_9b\,\hat{\mathbf{y}} - z_9c\sin\beta\,\hat{\mathbf{z}}$	(4*f*)	O V
$\mathbf{B}_{22}$	=	$x_9\,\mathbf{a}_1 + \left(\frac{1}{2} - y_9\right)\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$(x_9a + z_9c\cos\beta)\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_9\right)b\,\hat{\mathbf{y}} + z_9c\sin\beta\,\hat{\mathbf{z}}$	(4*f*)	O V

References:

- N. G. Batalieva and Y. A. Pyatenko, *Artificial yttrialite ("y-phase") - a representative of a new structure type in the rare earth diorthosilicate series*, Sov. Phys. Crystallogr. **16**, 786–789 (1972). Translation of N. G. Batalieva and Yu. A. Pyatenko, Kristallogr. 16, 905 (1971).
- A. I. Becerro, A. Escudero, P. Florian, D. Massiot, and M. D. Alba, *Revisiting $Y_2Si_2O_7$ and Y_2SiO_5 polymorphic structures by ^{89}Y MAS-NMR spectroscopy*, J. Solid State Chem. **177**, 2783–2789 (2004), doi:10.1016/j.jssc.2004.03.047.

Found in:

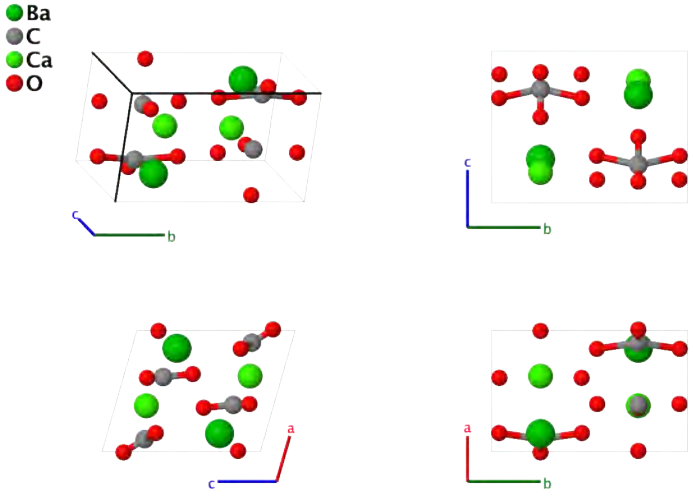
- A. I. Becerro and A. Escudero, *Revision of the crystallographic data of polymorphic $Y_2Si_2O_7$ and Y_2SiO_5 compounds*, Phase Transit. **77**, 1093–1102 (2004), doi:10.1080/01411590412331282814.

Geometry files:

- CIF: pp. 1521
- POSCAR: pp. 1522

Barytocalcite ($BaCa(CO_3)_2$) Structure: AB2CD6_mP20_11_e_2e_e_2e2f

http://aflow.org/prototype-encyclopedia/AB2CD6_mP20_11_e_2e_e_2e2f

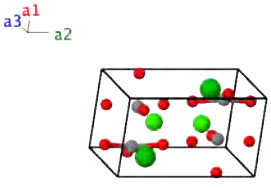


Prototype	:	BaC_2CaO_6
AFLOW prototype label	:	AB2CD6_mP20_11_e_2e_e_2e2f
***Strukturbericht* designation**	:	None
Pearson symbol	:	mP20
Space group number	:	11
Space group symbol	:	$P2_1/m$
AFLOW prototype command	:	`aflow --proto=AB2CD6_mP20_11_e_2e_e_2e2f --params=`$a, b/a, c/a, \beta, x_1, z_1, x_2, z_2, x_3, z_3, x_4, z_4, x_5, z_5, x_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8$

- $BaCa(CO_3)_2$ comes in a variety of crystal structures (Spahr, 2019):
  - monoclinic barytocalcite, space group $P2_1/m$ #11 (the current structure)
  - trigonal paralstonite, space group $P321$ #150,
  - Triclinic alstonite, space group $P1$ #1 or $P\bar{1}$ #2 (Sartori, 1975), and
  - a new monoclinic structure, space group $C2$ #5, synthesized by (Spahr, 2019), and lacking the centrosymmetric character of barytocalcite.

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(2e)$	Ba
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$(-x_1 a - z_1 c\cos\beta)\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(2e)$	Ba
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(2e)$	C I
$\mathbf{B}_4$	=	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(2e)$	C I
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(2e)$	C II
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(2e)$	C II
$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(2e)$	Ca
$\mathbf{B}_8$	=	$-x_4\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(2e)$	Ca
$\mathbf{B}_9$	=	$x_5\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(2e)$	O I
$\mathbf{B}_{10}$	=	$-x_5\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(2e)$	O I
$\mathbf{B}_{11}$	=	$x_6\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(2e)$	O II
$\mathbf{B}_{12}$	=	$-x_6\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(2e)$	O II
$\mathbf{B}_{13}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(4f)$	O III
$\mathbf{B}_{14}$	=	$-x_7\,\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right) b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(4f)$	O III
$\mathbf{B}_{15}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(4f)$	O III
$\mathbf{B}_{16}$	=	$x_7\,\mathbf{a}_1 + \left(\frac{1}{2} - y_7\right)\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_7\right) b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(4f)$	O III
$\mathbf{B}_{17}$	=	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(4f)$	O IV
$\mathbf{B}_{18}$	=	$-x_8\,\mathbf{a}_1 + \left(\frac{1}{2} + y_8\right)\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_8\right) b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(4f)$	O IV
$\mathbf{B}_{19}$	=	$-x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(4f)$	O IV
$\mathbf{B}_{20}$	=	$x_8\,\mathbf{a}_1 + \left(\frac{1}{2} - y_8\right)\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_8\right) b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(4f)$	O IV

References:

- B. Dickens and J. S. Bowen, *The Crystal Structure of $BaCa(CO_3)_2$ (barytocalcite)*, J. Res. Nat. Stand. Sec. A **75**, 197–203 (1971), doi:10.6028/jres.075A.020.
- F. Sartori, *New data on alstonite*, Lithos **8**, 199–207 (1975), doi:10.1016/0024-4937(75)90036-5.

Found in:

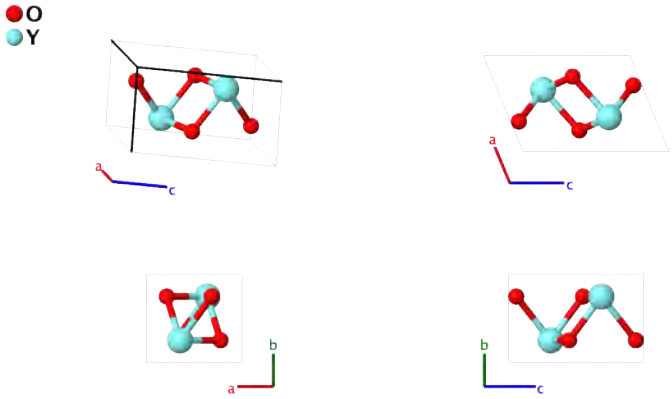
- D. Spahr, L. Bayarjargal, V. Vinograd, R. Luchitskaia, V. Milman, and B. Winkler, *A new $BaCa(CO_3)_2$ polymorph*, Acta Crystallogr. Sect. B Struct. Sci. **75**, 291–300 (2019), doi:10.1107/S2052520619003238.

Geometry files:

- CIF: pp. 1522
- POSCAR: pp. 1522

O(OH)Y Structure: ABC_mP6_11_e_e_e

http://aflow.org/prototype-encyclopedia/ABC_mP6_11_e_e_e



Prototype	:	HO_2Y
AFLOW prototype label	:	ABC_mP6_11_e_e_e
***Strukturbericht* designation**	:	None
Pearson symbol	:	mP6
Space group number	:	11
Space group symbol	:	$P2_1/m$
AFLOW prototype command	:	`aflow --proto=ABC_mP6_11_e_e_e --params=`$a, b/a, c/a, \beta, x_1, z_1, x_2, z_2, x_3, z_3$

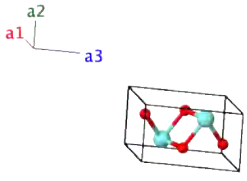
Other compounds with this structure

- HoO(OH), ErO(OH), and YbO(OH)

- For this prototype, the OH molecules are centered on the ($2e$) Wyckoff position.

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	$(x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_1 c\sin\beta\,\hat{\mathbf{z}}$	($2e$)	O
$\mathbf{B}_2$	$-x_1\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	$(-x_1 a - z_1 c\cos\beta)\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_1 c\sin\beta\,\hat{\mathbf{z}}$	($2e$)	O
$\mathbf{B}_3$	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	($2e$)	OH
$\mathbf{B}_4$	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	$(-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_2 c\sin\beta\,\hat{\mathbf{z}}$	($2e$)	OH

$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	(2*e*)	Y
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	(2*e*)	Y

References:

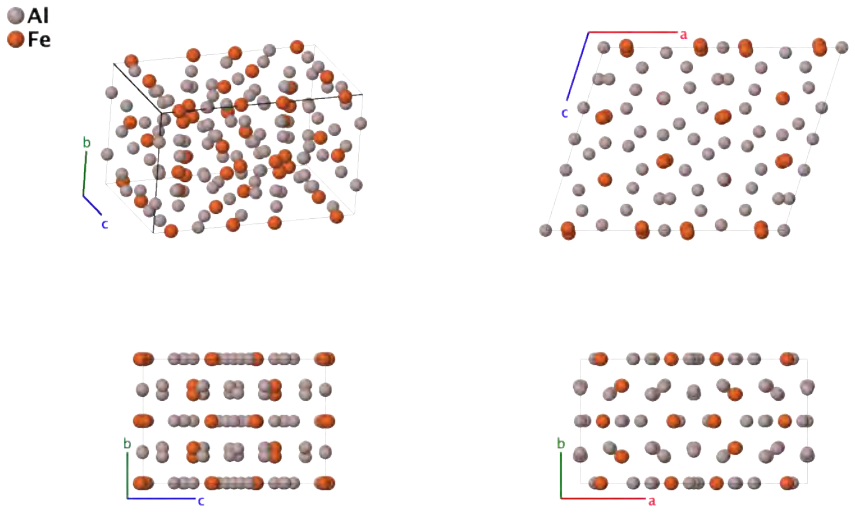
- R. F. Klevtsova and P. V. Klevtsov, *The Crystal Structure of YOOH*, J. Struct. Chem. **5**, 795–797 (1965), doi:10.1007/BF00744232.

Geometry files:

- CIF: pp. 1523
- POSCAR: pp. 1523

$Al_{13}Fe_4$ Structure: A13B4_mC102_12_dg8i5j_4ij

http://aflow.org/prototype-encyclopedia/A13B4_mC102_12_dg8i5j_4ij

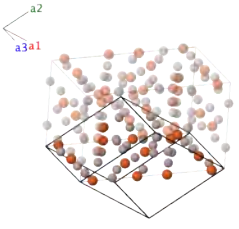


Prototype	:	$Al_{13}Fe_4$
AFLOW prototype label	:	A13B4_mC102_12_dg8i5j_4ij
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC102
Space group number	:	12
Space group symbol	:	$C2/m$
AFLOW prototype command	:	`aflow --proto=A13B4_mC102_12_dg8i5j_4ij --params=`$a, b/a, c/a, \beta, y_2, x_3, z_3, x_4, z_4, x_5, z_5, x_6, z_6, x_7, z_7, x_8, z_8, x_9, z_9, x_{10}, z_{10}, x_{11}, z_{11}, x_{12}, z_{12}, x_{13}, z_{13}, x_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}, x_{17}, y_{17}, z_{17}, x_{18}, y_{18}, z_{18}, x_{19}, y_{19}, z_{19}, x_{20}, y_{20}, z_{20}$

- The Al-IV site is only occupied 70% of the time. This makes the stoichiometry $Al_{12.8}Fe_4$, which (Black I and II, 1955) round to $Al_{13}Fe_4$.

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(a + c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	$(2d)$	Al I
$\mathbf{B}_2$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2$	=	$y_2 b\,\hat{\mathbf{y}}$	$(4g)$	Al II

$\mathbf{B}_3$	=	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2$	=	$-y_2 b\,\hat{\mathbf{y}}$	$(4g)$	Al II
$\mathbf{B}_4$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Al III
$\mathbf{B}_5$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Al III
$\mathbf{B}_6$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Al IV
$\mathbf{B}_7$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Al IV
$\mathbf{B}_8$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Al V
$\mathbf{B}_9$	=	$-x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Al V
$\mathbf{B}_{10}$	=	$x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Al VI
$\mathbf{B}_{11}$	=	$-x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Al VI
$\mathbf{B}_{12}$	=	$x_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Al VII
$\mathbf{B}_{13}$	=	$-x_7\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} - z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Al VII
$\mathbf{B}_{14}$	=	$x_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Al VIII
$\mathbf{B}_{15}$	=	$-x_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} - z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Al VIII
$\mathbf{B}_{16}$	=	$x_9\,\mathbf{a}_1 + x_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$(x_9 a + z_9 c\cos\beta)\,\hat{\mathbf{x}} + z_9 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Al IX
$\mathbf{B}_{17}$	=	$-x_9\,\mathbf{a}_1 - x_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$(-x_9 a - z_9 c\cos\beta)\,\hat{\mathbf{x}} - z_9 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Al IX
$\mathbf{B}_{18}$	=	$x_{10}\,\mathbf{a}_1 + x_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$(x_{10} a + z_{10} c\cos\beta)\,\hat{\mathbf{x}} + z_{10} c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Al X
$\mathbf{B}_{19}$	=	$-x_{10}\,\mathbf{a}_1 - x_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3$	=	$(-x_{10} a - z_{10} c\cos\beta)\,\hat{\mathbf{x}} - z_{10} c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Al X
$\mathbf{B}_{20}$	=	$x_{11}\,\mathbf{a}_1 + x_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$(x_{11} a + z_{11} c\cos\beta)\,\hat{\mathbf{x}} + z_{11} c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Fe I
$\mathbf{B}_{21}$	=	$-x_{11}\,\mathbf{a}_1 - x_{11}\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3$	=	$(-x_{11} a - z_{11} c\cos\beta)\,\hat{\mathbf{x}} - z_{11} c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Fe I
$\mathbf{B}_{22}$	=	$x_{12}\,\mathbf{a}_1 + x_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3$	=	$(x_{12} a + z_{12} c\cos\beta)\,\hat{\mathbf{x}} + z_{12} c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Fe II
$\mathbf{B}_{23}$	=	$-x_{12}\,\mathbf{a}_1 - x_{12}\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3$	=	$(-x_{12} a - z_{12} c\cos\beta)\,\hat{\mathbf{x}} - z_{12} c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Fe II
$\mathbf{B}_{24}$	=	$x_{13}\,\mathbf{a}_1 + x_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3$	=	$(x_{13} a + z_{13} c\cos\beta)\,\hat{\mathbf{x}} + z_{13} c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Fe III
$\mathbf{B}_{25}$	=	$-x_{13}\,\mathbf{a}_1 - x_{13}\,\mathbf{a}_2 - z_{13}\,\mathbf{a}_3$	=	$(-x_{13} a - z_{13} c\cos\beta)\,\hat{\mathbf{x}} - z_{13} c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Fe III
$\mathbf{B}_{26}$	=	$x_{14}\,\mathbf{a}_1 + x_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3$	=	$(x_{14} a + z_{14} c\cos\beta)\,\hat{\mathbf{x}} + z_{14} c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Fe IV
$\mathbf{B}_{27}$	=	$-x_{14}\,\mathbf{a}_1 - x_{14}\,\mathbf{a}_2 - z_{14}\,\mathbf{a}_3$	=	$(-x_{14} a - z_{14} c\cos\beta)\,\hat{\mathbf{x}} - z_{14} c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Fe IV
$\mathbf{B}_{28}$	=	$(x_{15} - y_{15})\,\mathbf{a}_1 + (x_{15} + y_{15})\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3$	=	$(x_{15} a + z_{15} c\cos\beta)\,\hat{\mathbf{x}} + y_{15} b\,\hat{\mathbf{y}} + z_{15} c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Al XI
$\mathbf{B}_{29}$	=	$(-x_{15} - y_{15})\,\mathbf{a}_1 + (-x_{15} + y_{15})\,\mathbf{a}_2 - z_{15}\,\mathbf{a}_3$	=	$(-x_{15} a - z_{15} c\cos\beta)\,\hat{\mathbf{x}} + y_{15} b\,\hat{\mathbf{y}} - z_{15} c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Al XI
$\mathbf{B}_{30}$	=	$(-x_{15} + y_{15})\,\mathbf{a}_1 + (-x_{15} - y_{15})\,\mathbf{a}_2 - z_{15}\,\mathbf{a}_3$	=	$(-x_{15} a - z_{15} c\cos\beta)\,\hat{\mathbf{x}} - y_{15} b\,\hat{\mathbf{y}} - z_{15} c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Al XI
$\mathbf{B}_{31}$	=	$(x_{15} + y_{15})\,\mathbf{a}_1 + (x_{15} - y_{15})\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3$	=	$(x_{15} a + z_{15} c\cos\beta)\,\hat{\mathbf{x}} - y_{15} b\,\hat{\mathbf{y}} + z_{15} c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Al XI
$\mathbf{B}_{32}$	=	$(x_{16} - y_{16})\,\mathbf{a}_1 + (x_{16} + y_{16})\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3$	=	$(x_{16} a + z_{16} c\cos\beta)\,\hat{\mathbf{x}} + y_{16} b\,\hat{\mathbf{y}} + z_{16} c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Al XII
$\mathbf{B}_{33}$	=	$(-x_{16} - y_{16})\,\mathbf{a}_1 + (-x_{16} + y_{16})\,\mathbf{a}_2 - z_{16}\,\mathbf{a}_3$	=	$(-x_{16} a - z_{16} c\cos\beta)\,\hat{\mathbf{x}} + y_{16} b\,\hat{\mathbf{y}} - z_{16} c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Al XII
$\mathbf{B}_{34}$	=	$(-x_{16} + y_{16})\,\mathbf{a}_1 + (-x_{16} - y_{16})\,\mathbf{a}_2 - z_{16}\,\mathbf{a}_3$	=	$(-x_{16} a - z_{16} c\cos\beta)\,\hat{\mathbf{x}} - y_{16} b\,\hat{\mathbf{y}} - z_{16} c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Al XII

$\mathbf{B}_{35}$	=	$(x_{16}+y_{16})\,\mathbf{a}_1+(x_{16}-y_{16})\,\mathbf{a}_2+z_{16}\,\mathbf{a}_3$	=	$(x_{16}a+z_{16}c\cos\beta)\,\hat{\mathbf{x}}-y_{16}b\,\hat{\mathbf{y}}+z_{16}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Al XII
$\mathbf{B}_{36}$	=	$(x_{17}-y_{17})\,\mathbf{a}_1+(x_{17}+y_{17})\,\mathbf{a}_2+z_{17}\,\mathbf{a}_3$	=	$(x_{17}a+z_{17}c\cos\beta)\,\hat{\mathbf{x}}+y_{17}b\,\hat{\mathbf{y}}+z_{17}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Al XIII
$\mathbf{B}_{37}$	=	$(-x_{17}-y_{17})\,\mathbf{a}_1+(-x_{17}+y_{17})\,\mathbf{a}_2-z_{17}\,\mathbf{a}_3$	=	$(-x_{17}a-z_{17}c\cos\beta)\,\hat{\mathbf{x}}+y_{17}b\,\hat{\mathbf{y}}-z_{17}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Al XIII
$\mathbf{B}_{38}$	=	$(-x_{17}+y_{17})\,\mathbf{a}_1+(-x_{17}-y_{17})\,\mathbf{a}_2-z_{17}\,\mathbf{a}_3$	=	$(-x_{17}a-z_{17}c\cos\beta)\,\hat{\mathbf{x}}-y_{17}b\,\hat{\mathbf{y}}-z_{17}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Al XIII
$\mathbf{B}_{39}$	=	$(x_{17}+y_{17})\,\mathbf{a}_1+(x_{17}-y_{17})\,\mathbf{a}_2+z_{17}\,\mathbf{a}_3$	=	$(x_{17}a+z_{17}c\cos\beta)\,\hat{\mathbf{x}}-y_{17}b\,\hat{\mathbf{y}}+z_{17}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Al XIII
$\mathbf{B}_{40}$	=	$(x_{18}-y_{18})\,\mathbf{a}_1+(x_{18}+y_{18})\,\mathbf{a}_2+z_{18}\,\mathbf{a}_3$	=	$(x_{18}a+z_{18}c\cos\beta)\,\hat{\mathbf{x}}+y_{18}b\,\hat{\mathbf{y}}+z_{18}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Al XIV
$\mathbf{B}_{41}$	=	$(-x_{18}-y_{18})\,\mathbf{a}_1+(-x_{18}+y_{18})\,\mathbf{a}_2-z_{18}\,\mathbf{a}_3$	=	$(-x_{18}a-z_{18}c\cos\beta)\,\hat{\mathbf{x}}+y_{18}b\,\hat{\mathbf{y}}-z_{18}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Al XIV
$\mathbf{B}_{42}$	=	$(-x_{18}+y_{18})\,\mathbf{a}_1+(-x_{18}-y_{18})\,\mathbf{a}_2-z_{18}\,\mathbf{a}_3$	=	$(-x_{18}a-z_{18}c\cos\beta)\,\hat{\mathbf{x}}-y_{18}b\,\hat{\mathbf{y}}-z_{18}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Al XIV
$\mathbf{B}_{43}$	=	$(x_{18}+y_{18})\,\mathbf{a}_1+(x_{18}-y_{18})\,\mathbf{a}_2+z_{18}\,\mathbf{a}_3$	=	$(x_{18}a+z_{18}c\cos\beta)\,\hat{\mathbf{x}}-y_{18}b\,\hat{\mathbf{y}}+z_{18}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Al XIV
$\mathbf{B}_{44}$	=	$(x_{19}-y_{19})\,\mathbf{a}_1+(x_{19}+y_{19})\,\mathbf{a}_2+z_{19}\,\mathbf{a}_3$	=	$(x_{19}a+z_{19}c\cos\beta)\,\hat{\mathbf{x}}+y_{19}b\,\hat{\mathbf{y}}+z_{19}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Al XV
$\mathbf{B}_{45}$	=	$(-x_{19}-y_{19})\,\mathbf{a}_1+(-x_{19}+y_{19})\,\mathbf{a}_2-z_{19}\,\mathbf{a}_3$	=	$(-x_{19}a-z_{19}c\cos\beta)\,\hat{\mathbf{x}}+y_{19}b\,\hat{\mathbf{y}}-z_{19}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Al XV
$\mathbf{B}_{46}$	=	$(-x_{19}+y_{19})\,\mathbf{a}_1+(-x_{19}-y_{19})\,\mathbf{a}_2-z_{19}\,\mathbf{a}_3$	=	$(-x_{19}a-z_{19}c\cos\beta)\,\hat{\mathbf{x}}-y_{19}b\,\hat{\mathbf{y}}-z_{19}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Al XV
$\mathbf{B}_{47}$	=	$(x_{19}+y_{19})\,\mathbf{a}_1+(x_{19}-y_{19})\,\mathbf{a}_2+z_{19}\,\mathbf{a}_3$	=	$(x_{19}a+z_{19}c\cos\beta)\,\hat{\mathbf{x}}-y_{19}b\,\hat{\mathbf{y}}+z_{19}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Al XV
$\mathbf{B}_{48}$	=	$(x_{20}-y_{20})\,\mathbf{a}_1+(x_{20}+y_{20})\,\mathbf{a}_2+z_{20}\,\mathbf{a}_3$	=	$(x_{20}a+z_{20}c\cos\beta)\,\hat{\mathbf{x}}+y_{20}b\,\hat{\mathbf{y}}+z_{20}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Fe V
$\mathbf{B}_{49}$	=	$(-x_{20}-y_{20})\,\mathbf{a}_1+(-x_{20}+y_{20})\,\mathbf{a}_2-z_{20}\,\mathbf{a}_3$	=	$(-x_{20}a-z_{20}c\cos\beta)\,\hat{\mathbf{x}}+y_{20}b\,\hat{\mathbf{y}}-z_{20}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Fe V
$\mathbf{B}_{50}$	=	$(-x_{20}+y_{20})\,\mathbf{a}_1+(-x_{20}-y_{20})\,\mathbf{a}_2-z_{20}\,\mathbf{a}_3$	=	$(-x_{20}a-z_{20}c\cos\beta)\,\hat{\mathbf{x}}-y_{20}b\,\hat{\mathbf{y}}-z_{20}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Fe V
$\mathbf{B}_{51}$	=	$(x_{20}+y_{20})\,\mathbf{a}_1+(x_{20}-y_{20})\,\mathbf{a}_2+z_{20}\,\mathbf{a}_3$	=	$(x_{20}a+z_{20}c\cos\beta)\,\hat{\mathbf{x}}-y_{20}b\,\hat{\mathbf{y}}+z_{20}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Fe V

References:

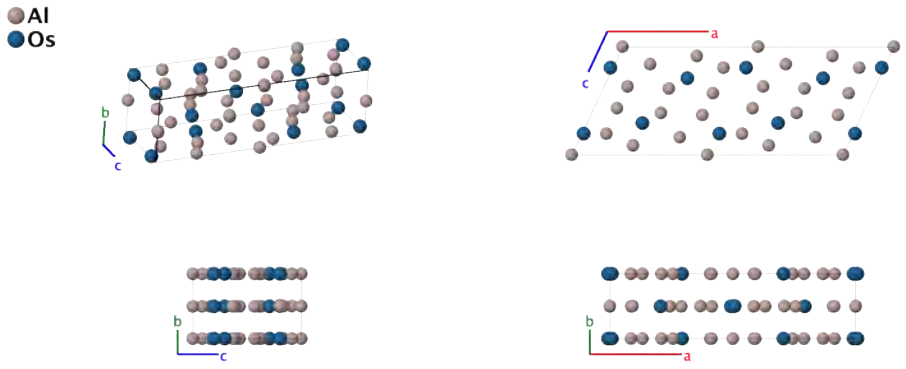
- P. J. Black, *The Structure of $FeAl_3$. I*, Acta Cryst. **8**, 43–48 (1955), doi:10.1107/S0365110X5500011X.
- P. J. Black, *The Structure of $FeAl_3$. II*, Acta Cryst. **8**, 175–182 (1955), doi:10.1107/S0365110X55000637.

Geometry files:

- CIF: pp. 1523
- POSCAR: pp. 1523

Os_4Al_{13} Structure: A13B4_mC34_12_b6i_2i

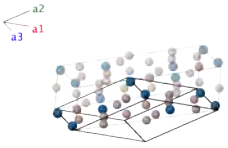
http://aflow.org/prototype-encyclopedia/A13B4_mC34_12_b6i_2i



Prototype	:	$Al_{13}Os_4$
AFLOW prototype label	:	A13B4_mC34_12_b6i_2i
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC34
Space group number	:	12
Space group symbol	:	$C2/m$
AFLOW prototype command	:	`aflow --proto=A13B4_mC34_12_b6i_2i --params=`$a, b/a, c/a, \beta, x_2, z_2, x_3, z_3, x_4, z_4, x_5, z_5, x_6, z_6, x_7, z_7, x_8, z_8, x_9, z_9$

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{2}\mathbf{a}_1 + \frac{1}{2}\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	$(2b)$	Al I
$\mathbf{B}_2$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2a + z_2c\cos\beta)\,\hat{\mathbf{x}} + z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Al II
$\mathbf{B}_3$	=	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2a - z_2c\cos\beta)\,\hat{\mathbf{x}} - z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Al II
$\mathbf{B}_4$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3a + z_3c\cos\beta)\,\hat{\mathbf{x}} + z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Al III
$\mathbf{B}_5$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3a - z_3c\cos\beta)\,\hat{\mathbf{x}} - z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Al III
$\mathbf{B}_6$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4a + z_4c\cos\beta)\,\hat{\mathbf{x}} + z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Al IV
$\mathbf{B}_7$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4a - z_4c\cos\beta)\,\hat{\mathbf{x}} - z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Al IV
$\mathbf{B}_8$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5a + z_5c\cos\beta)\,\hat{\mathbf{x}} + z_5c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Al V
$\mathbf{B}_9$	=	$-x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5a - z_5c\cos\beta)\,\hat{\mathbf{x}} - z_5c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Al V
$\mathbf{B}_{10}$	=	$x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6a + z_6c\cos\beta)\,\hat{\mathbf{x}} + z_6c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Al VI

$\mathbf{B}_{11}$	=	$-x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Al VI
$\mathbf{B}_{12}$	=	$x_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Al VII
$\mathbf{B}_{13}$	=	$-x_7\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} - z_7 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Al VII
$\mathbf{B}_{14}$	=	$x_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + z_8 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Os I
$\mathbf{B}_{15}$	=	$-x_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} - z_8 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Os I
$\mathbf{B}_{16}$	=	$x_9\,\mathbf{a}_1 + x_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$(x_9 a + z_9 c\cos\beta)\,\hat{\mathbf{x}} + z_9 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Os II
$\mathbf{B}_{17}$	=	$-x_9\,\mathbf{a}_1 - x_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$(-x_9 a - z_9 c\cos\beta)\,\hat{\mathbf{x}} - z_9 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Os II

References:

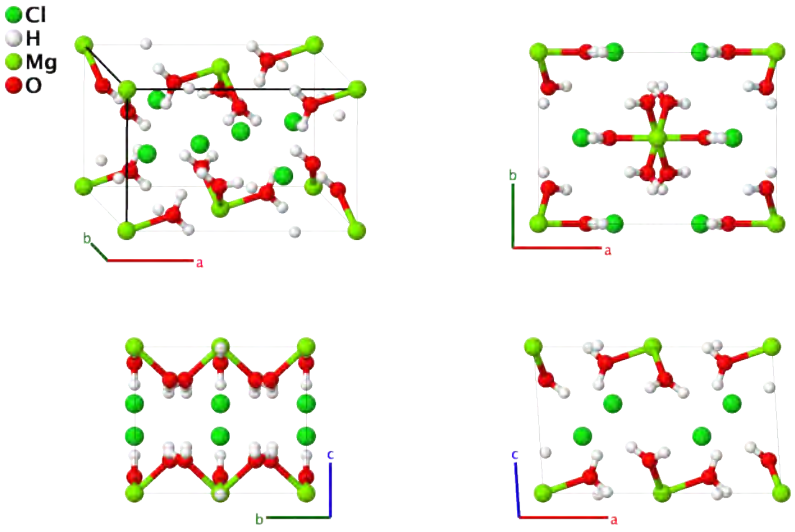
- L.-E. Edshammar, *The Crystal Structure of Os_4Al_{13}*, Acta Chem. Scand. **18**, 2294–2302 (1964), doi:10.3891/acta.chem.scand.18-2294.

Geometry files:

- CIF: pp. 1524
- POSCAR: pp. 1524

Bischofite ($MgCl_2{\cdot}6H_2O$, $J1_7$) Structure: A2B12CD6_mC42_12_i_2i2j_a_ij

http://aflow.org/prototype-encyclopedia/A2B12CD6_mC42_12_i_2i2j_a_ij



Prototype	:	$Cl_2H_{12}MgO_6$
AFLOW prototype label	:	A2B12CD6_mC42_12_i_2i2j_a_ij
***Strukturbericht* designation**	:	$J1_7$
Pearson symbol	:	mC42
Space group number	:	12
Space group symbol	:	$C2/m$
AFLOW prototype command	:	`aflow --proto=A2B12CD6_mC42_12_i_2i2j_a_ij --params=`$a, b/a, c/a, \beta, x_2, z_2, x_3, z_3, x_4, z_4, x_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8$

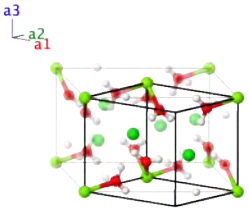
Other compounds with this structure

- $MgBr_2{\cdot}6H_2O$

- This structure is nearly identical to the one presented in (Gottfried, 1937) as $J1_7$, but now includes the positions of the hydrogen atoms.

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	Mg
$\mathbf{B}_2$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Cl
$\mathbf{B}_3$	=	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} - z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Cl
$\mathbf{B}_4$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	H I
$\mathbf{B}_5$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	H I
$\mathbf{B}_6$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	H II
$\mathbf{B}_7$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	H II
$\mathbf{B}_8$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	O I
$\mathbf{B}_9$	=	$-x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	O I
$\mathbf{B}_{10}$	=	$(x_6 - y_6)\,\mathbf{a}_1 + (x_6 + y_6)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	H III
$\mathbf{B}_{11}$	=	$(-x_6 - y_6)\,\mathbf{a}_1 + (-x_6 + y_6)\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	H III
$\mathbf{B}_{12}$	=	$(-x_6 + y_6)\,\mathbf{a}_1 + (-x_6 - y_6)\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	H III
$\mathbf{B}_{13}$	=	$(x_6 + y_6)\,\mathbf{a}_1 + (x_6 - y_6)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	H III
$\mathbf{B}_{14}$	=	$(x_7 - y_7)\,\mathbf{a}_1 + (x_7 + y_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	H IV
$\mathbf{B}_{15}$	=	$(-x_7 - y_7)\,\mathbf{a}_1 + (-x_7 + y_7)\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	H IV
$\mathbf{B}_{16}$	=	$(-x_7 + y_7)\,\mathbf{a}_1 + (-x_7 - y_7)\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	H IV
$\mathbf{B}_{17}$	=	$(x_7 + y_7)\,\mathbf{a}_1 + (x_7 - y_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	H IV
$\mathbf{B}_{18}$	=	$(x_8 - y_8)\,\mathbf{a}_1 + (x_8 + y_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O II
$\mathbf{B}_{19}$	=	$(-x_8 - y_8)\,\mathbf{a}_1 + (-x_8 + y_8)\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O II
$\mathbf{B}_{20}$	=	$(-x_8 + y_8)\,\mathbf{a}_1 + (-x_8 - y_8)\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O II
$\mathbf{B}_{21}$	=	$(x_8 + y_8)\,\mathbf{a}_1 + (x_8 - y_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O II

References:

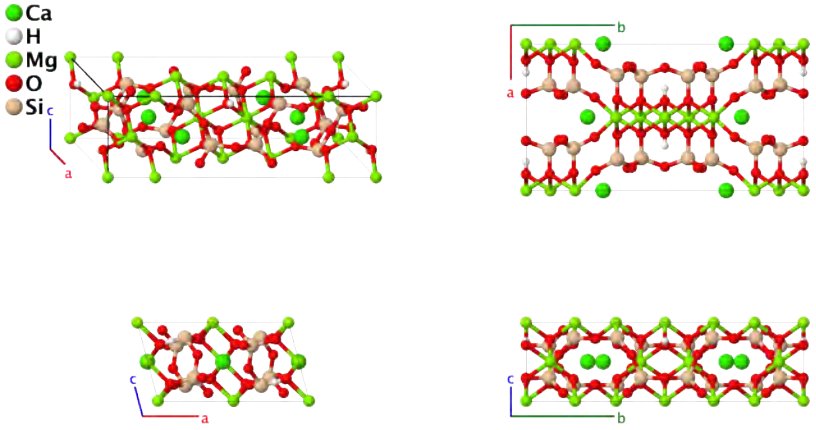
- P. A. Agron and W. R. Busing, *Magnesium dichloride hexahydrate, $MgCl_2{\cdot}6H_2O$, by neutron diffraction*, Acta Crystallogr. C **41**, 8–10 (1985), doi:10.1107/S0108270185002591.
- C. Gottfried and F. Schossberger, eds., *Strukturbericht Band III 1933-1935* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1524
- POSCAR: pp. 1525

Tremolite ($Ca_2Mg_5Si_8O_{22}(OH)_2$, $S4_2$) Structure: A2B2C5D24E8_mC82_12_h_i_agh_2i5j_2j

http://aflow.org/prototype-encyclopedia/A2B2C5D24E8_mC82_12_h_i_agh_2i5j_2j



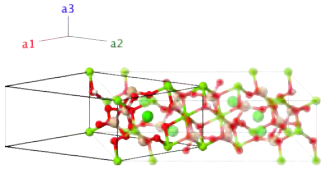
Prototype	:	$Ca_2H_2Mg_5O_{24}Si_8$
AFLOW prototype label	:	A2B2C5D24E8_mC82_12_h_i_agh_2i5j_2j
***Strukturbericht* designation**	:	$S4_2$
Pearson symbol	:	mC82
Space group number	:	12
Space group symbol	:	$C2/m$
AFLOW prototype command	:	`aflow --proto=A2B2C5D24E8_mC82_12_h_i_agh_2i5j_2j --params=`$a, b/a, c/a, \beta, y_2, y_3, y_4, x_5, z_5, x_6, z_6, x_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}$

Other compounds with this structure

- $Ca_2H_2(Mg_{5-x},Fe_x)O_{24}Si_8$ and $Ca_2F_2Mg_5O_{22}Si_8$

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	Mg I
$\mathbf{B}_2$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2$	=	$y_2 b\,\hat{\mathbf{y}}$	$(4g)$	Mg II
$\mathbf{B}_3$	=	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2$	=	$-y_2 b\,\hat{\mathbf{y}}$	$(4g)$	Mg II

$\mathbf{B}_4$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	(4*h*)	Ca
$\mathbf{B}_5$	=	$y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	(4*h*)	Ca
$\mathbf{B}_6$	=	$-y_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	(4*h*)	Mg III
$\mathbf{B}_7$	=	$y_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	(4*h*)	Mg III
$\mathbf{B}_8$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	H
$\mathbf{B}_9$	=	$-x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	H
$\mathbf{B}_{10}$	=	$x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	O I
$\mathbf{B}_{11}$	=	$-x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	O I
$\mathbf{B}_{12}$	=	$x_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	O II
$\mathbf{B}_{13}$	=	$-x_7\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} - z_7 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	O II
$\mathbf{B}_{14}$	=	$(x_8 - y_8)\,\mathbf{a}_1 + (x_8 + y_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	O III
$\mathbf{B}_{15}$	=	$(-x_8 - y_8)\,\mathbf{a}_1 + (-x_8 + y_8)\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	O III
$\mathbf{B}_{16}$	=	$(-x_8 + y_8)\,\mathbf{a}_1 + (-x_8 - y_8)\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	O III
$\mathbf{B}_{17}$	=	$(x_8 + y_8)\,\mathbf{a}_1 + (x_8 - y_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	O III
$\mathbf{B}_{18}$	=	$(x_9 - y_9)\,\mathbf{a}_1 + (x_9 + y_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$(x_9 a + z_9 c\cos\beta)\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	O IV
$\mathbf{B}_{19}$	=	$(-x_9 - y_9)\,\mathbf{a}_1 + (-x_9 + y_9)\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$(-x_9 a - z_9 c\cos\beta)\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} - z_9 c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	O IV
$\mathbf{B}_{20}$	=	$(-x_9 + y_9)\,\mathbf{a}_1 + (-x_9 - y_9)\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$(-x_9 a - z_9 c\cos\beta)\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} - z_9 c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	O IV
$\mathbf{B}_{21}$	=	$(x_9 + y_9)\,\mathbf{a}_1 + (x_9 - y_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$(x_9 a + z_9 c\cos\beta)\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + z_9 c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	O IV
$\mathbf{B}_{22}$	=	$(x_{10} - y_{10})\,\mathbf{a}_1 + (x_{10} + y_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$(x_{10} a + z_{10} c\cos\beta)\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	O V
$\mathbf{B}_{23}$	=	$(-x_{10} - y_{10})\,\mathbf{a}_1 + (-x_{10} + y_{10})\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3$	=	$(-x_{10} a - z_{10} c\cos\beta)\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} - z_{10} c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	O V
$\mathbf{B}_{24}$	=	$(-x_{10} + y_{10})\,\mathbf{a}_1 + (-x_{10} - y_{10})\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3$	=	$(-x_{10} a - z_{10} c\cos\beta)\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} - z_{10} c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	O V
$\mathbf{B}_{25}$	=	$(x_{10} + y_{10})\,\mathbf{a}_1 + (x_{10} - y_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$(x_{10} a + z_{10} c\cos\beta)\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} + z_{10} c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	O V
$\mathbf{B}_{26}$	=	$(x_{11} - y_{11})\,\mathbf{a}_1 + (x_{11} + y_{11})\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$(x_{11} a + z_{11} c\cos\beta)\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + z_{11} c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	O VI
$\mathbf{B}_{27}$	=	$(-x_{11} - y_{11})\,\mathbf{a}_1 + (-x_{11} + y_{11})\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3$	=	$(-x_{11} a - z_{11} c\cos\beta)\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} - z_{11} c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	O VI
$\mathbf{B}_{28}$	=	$(-x_{11} + y_{11})\,\mathbf{a}_1 + (-x_{11} - y_{11})\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3$	=	$(-x_{11} a - z_{11} c\cos\beta)\,\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} - z_{11} c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	O VI
$\mathbf{B}_{29}$	=	$(x_{11} + y_{11})\,\mathbf{a}_1 + (x_{11} - y_{11})\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$(x_{11} a + z_{11} c\cos\beta)\,\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} + z_{11} c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	O VI
$\mathbf{B}_{30}$	=	$(x_{12} - y_{12})\,\mathbf{a}_1 + (x_{12} + y_{12})\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3$	=	$(x_{12} a + z_{12} c\cos\beta)\,\hat{\mathbf{x}} + y_{12} b\,\hat{\mathbf{y}} + z_{12} c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	O VII
$\mathbf{B}_{31}$	=	$(-x_{12} - y_{12})\,\mathbf{a}_1 + (-x_{12} + y_{12})\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3$	=	$(-x_{12} a - z_{12} c\cos\beta)\,\hat{\mathbf{x}} + y_{12} b\,\hat{\mathbf{y}} - z_{12} c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	O VII

$\mathbf{B}_{32}$	=	$(-x_{12}+y_{12})\,\mathbf{a}_1 + (-x_{12}-y_{12})\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3$	=	$(-x_{12}a - z_{12}c\cos\beta)\,\hat{\mathbf{x}} - y_{12}b\,\hat{\mathbf{y}} - z_{12}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O VII
$\mathbf{B}_{33}$	=	$(x_{12}+y_{12})\,\mathbf{a}_1 + (x_{12}-y_{12})\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3$	=	$(x_{12}a + z_{12}c\cos\beta)\,\hat{\mathbf{x}} - y_{12}b\,\hat{\mathbf{y}} + z_{12}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O VII
$\mathbf{B}_{34}$	=	$(x_{13}-y_{13})\,\mathbf{a}_1 + (x_{13}+y_{13})\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3$	=	$(x_{13}a + z_{13}c\cos\beta)\,\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} + z_{13}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Si I
$\mathbf{B}_{35}$	=	$(-x_{13}-y_{13})\,\mathbf{a}_1 + (-x_{13}+y_{13})\,\mathbf{a}_2 - z_{13}\,\mathbf{a}_3$	=	$(-x_{13}a - z_{13}c\cos\beta)\,\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} - z_{13}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Si I
$\mathbf{B}_{36}$	=	$(-x_{13}+y_{13})\,\mathbf{a}_1 + (-x_{13}-y_{13})\,\mathbf{a}_2 - z_{13}\,\mathbf{a}_3$	=	$(-x_{13}a - z_{13}c\cos\beta)\,\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} - z_{13}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Si I
$\mathbf{B}_{37}$	=	$(x_{13}+y_{13})\,\mathbf{a}_1 + (x_{13}-y_{13})\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3$	=	$(x_{13}a + z_{13}c\cos\beta)\,\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} + z_{13}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Si I
$\mathbf{B}_{38}$	=	$(x_{14}-y_{14})\,\mathbf{a}_1 + (x_{14}+y_{14})\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3$	=	$(x_{14}a + z_{14}c\cos\beta)\,\hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} + z_{14}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Si II
$\mathbf{B}_{39}$	=	$(-x_{14}-y_{14})\,\mathbf{a}_1 + (-x_{14}+y_{14})\,\mathbf{a}_2 - z_{14}\,\mathbf{a}_3$	=	$(-x_{14}a - z_{14}c\cos\beta)\,\hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} - z_{14}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Si II
$\mathbf{B}_{40}$	=	$(-x_{14}+y_{14})\,\mathbf{a}_1 + (-x_{14}-y_{14})\,\mathbf{a}_2 - z_{14}\,\mathbf{a}_3$	=	$(-x_{14}a - z_{14}c\cos\beta)\,\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} - z_{14}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Si II
$\mathbf{B}_{41}$	=	$(x_{14}+y_{14})\,\mathbf{a}_1 + (x_{14}-y_{14})\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3$	=	$(x_{14}a + z_{14}c\cos\beta)\,\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} + z_{14}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Si II

References:

- M. Merli, L. Ungaretti, and R. Oberti, *Leverage analysis and structure refinement of minerals*, Am. Mineral. **85**, 532–542 (2000).

Found in:

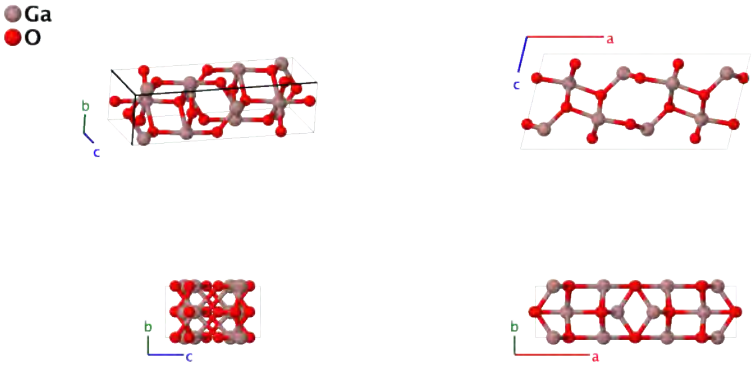
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1525
- POSCAR: pp. 1525

β-Ga_2O_3 Structure: A2B3_mC20_12_2i_3i

http://aflow.org/prototype-encyclopedia/A2B3_mC20_12_2i_3i

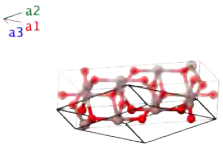


Prototype	:	Ga_2O_3
AFLOW prototype label	:	A2B3_mC20_12_2i_3i
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC20
Space group number	:	12
Space group symbol	:	$C2/m$
AFLOW prototype command	:	`aflow --proto=A2B3_mC20_12_2i_3i --params=`$a, b/a, c/a, \beta, x_1, z_1, x_2, z_2, x_3, z_3, x_4, z_4, x_5, z_5$

- Ga_2O_3 takes on a variety of structures:
  - α-Ga_2O_3, which has the corundum ($D5_1$) structure,
  - β-Ga_2O_3, this structure,
  - γ-Ga_2O_3, and
  - ϵ-Ga_2O_3, a structure with many vacancies which can be approximated by the κ-alumina structure.

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + x_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + z_1 c\sin\beta\,\hat{\mathbf{z}}$	($4i$)	Ga I
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 - x_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$(-x_1 a - z_1 c\cos\beta)\,\hat{\mathbf{x}} - z_1 c\sin\beta\,\hat{\mathbf{z}}$	($4i$)	Ga I
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	($4i$)	Ga II
$\mathbf{B}_4$	=	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} - z_2 c\sin\beta\,\hat{\mathbf{z}}$	($4i$)	Ga II

$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	O I
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	O I
$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	O II
$\mathbf{B}_8$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	O II
$\mathbf{B}_9$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	O III
$\mathbf{B}_{10}$	=	$-x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	O III

References:

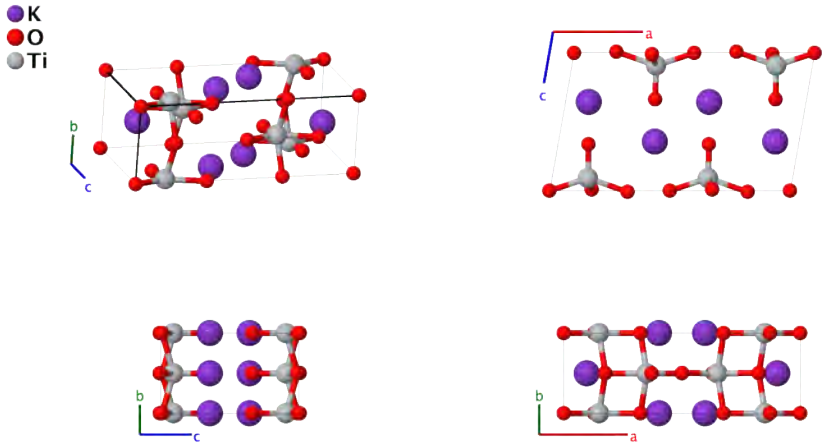
- J. Åhman, G. Svensson, and J. Albertsson, *A Reinvestigation of β-Gallium Oxide*, Acta Crystallogr. C **52**, 1336–1338 (1996), doi:10.1107/S0108270195016404.

Geometry files:

- CIF: pp. 1526
- POSCAR: pp. 1526

$K_2Ti_2O_5$ Structure: A2B5C2_mC18_12_i_a2i_i

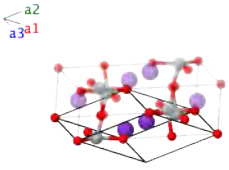
http://aflow.org/prototype-encyclopedia/A2B5C2_mC18_12_i_a2i_i



Prototype	:	$K_2O_5Ti_2$
AFLOW prototype label	:	A2B5C2_mC18_12_i_a2i_i
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC18
Space group number	:	12
Space group symbol	:	$C2/m$
AFLOW prototype command	:	`aflow --proto=A2B5C2_mC18_12_i_a2i_i --params=`$a, b/a, c/a, \beta, x_2, z_2, x_3, z_3, x_4, z_4, x_5, z_5$

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	O I
$\mathbf{B}_2$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2a + z_2c\cos\beta)\,\hat{\mathbf{x}} + z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	K
$\mathbf{B}_3$	=	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2a - z_2c\cos\beta)\,\hat{\mathbf{x}} - z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	K
$\mathbf{B}_4$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3a + z_3c\cos\beta)\,\hat{\mathbf{x}} + z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	O II
$\mathbf{B}_5$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3a - z_3c\cos\beta)\,\hat{\mathbf{x}} - z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	O II
$\mathbf{B}_6$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4a + z_4c\cos\beta)\,\hat{\mathbf{x}} + z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	O III
$\mathbf{B}_7$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4a - z_4c\cos\beta)\,\hat{\mathbf{x}} - z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	O III

$\mathbf{B}_8$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Ti
$\mathbf{B}_9$	=	$-x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Ti

References:

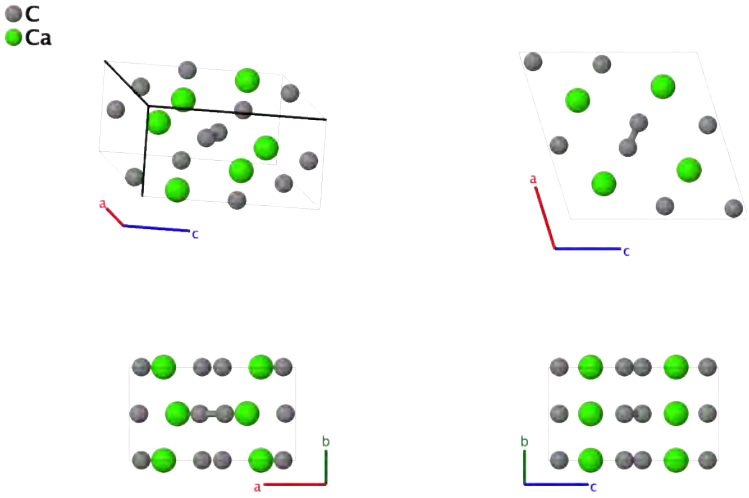
- S. Andersson and A. D. Wadsley, *The Crystal Structure of* $K_2Ti_2O_5$, Acta Chem. Scand. **15**, 663–669 (1961), doi:10.3891/acta.chem.scand.15-0663.

Geometry files:

- CIF: pp. 1526
- POSCAR: pp. 1526

CaC_2-III Structure: A2B_mC12_12_2i_i

http://aflow.org/prototype-encyclopedia/A2B_mC12_12_2i_i

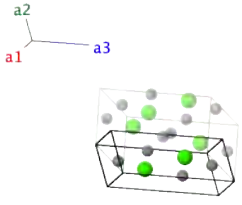


Prototype	:	C_2Ca
AFLOW prototype label	:	A2B_mC12_12_2i_i
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC12
Space group number	:	12
Space group symbol	:	$C2/m$
AFLOW prototype command	:	`aflow --proto=A2B_mC12_12_2i_i --params=`$a, b/a, c/a, \beta, x_1, z_1, x_2, z_2, x_3, z_3$

- This is a meta-stable room temperature structure.
- The stable room temperature configuration has the $C11_a$ structure.
- The low-temperature CaC_2 crystallizes in the ThC_2 (C_g) structure.

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$ =	$x_1\,\mathbf{a}_1 + x_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + z_1 c\sin\beta\,\hat{\mathbf{z}}$	($4i$)	C I

$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 - x_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$(-x_1 a - z_1 c\cos\beta)\,\hat{\mathbf{x}} - z_1 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	C I
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	C II
$\mathbf{B}_4$	=	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} - z_2 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	C II
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Ca
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Ca

References:

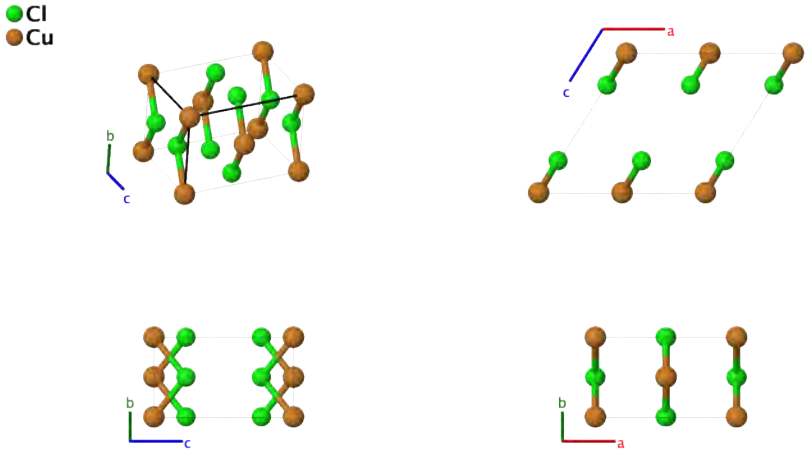
- M. Knapp and U. Ruschewitz, *Structural Phase Transitions in CaC_2*, Chem.: Eur. J. **7**, 874–880 (2001), doi:10.1002/1521-3765(20010216)7:4%3C874::AID-CHEM874%3E3.0.CO;2-V.

Geometry files:

- CIF: pp. 1526
- POSCAR: pp. 1527

Tolbachite ($CuCl_2$) Structure: A2B_mC6_12_i_a

http://aflow.org/prototype-encyclopedia/A2B_mC6_12_i_a



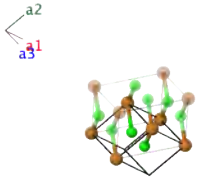
Prototype	:	Cl_2Cu
AFLOW prototype label	:	A2B_mC6_12_i_a
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC6
Space group number	:	12
Space group symbol	:	$C2/m$
AFLOW prototype command	:	`aflow --proto=A2B_mC6_12_i_a --params=`$a, b/a, c/a, \beta, x_2, z_2$

Other compounds with this structure

- $CuBr_2$

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	Cu
$\mathbf{B}_2$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2a + z_2c\cos\beta)\,\hat{\mathbf{x}} + z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Cl
$\mathbf{B}_3$	=	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2a - z_2c\cos\beta)\,\hat{\mathbf{x}} - z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Cl

References:

- P. C. Burns and F. C. Hawthorne, *Tolbachite, $CuCl_2$, the first example of Cu_2^+ octahedrally coordinated by Cl^-*, Am. Mineral. **78**, 187–189 (1993).

Found in:

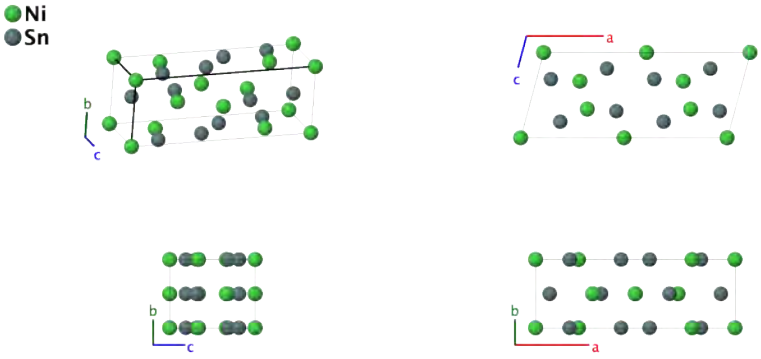
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1527
- POSCAR: pp. 1527

δ-Ni_3Sn_4 ($D7_a$) Structure: A3B4_mC14_12_ai_2i

http://aflow.org/prototype-encyclopedia/A3B4_mC14_12_ai_2i



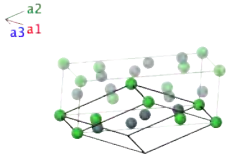
Prototype	:	Ni_3Sn_4
AFLOW prototype label	:	A3B4_mC14_12_ai_2i
***Strukturbericht* designation**	:	$D7_a$
Pearson symbol	:	mC14
Space group number	:	12
Space group symbol	:	$C2/m$
AFLOW prototype command	:	`aflow --proto=A3B4_mC14_12_ai_2i --params=`$a, b/a, c/a, \beta, x_2, z_2, x_3, z_3, x_4, z_4$

Other compounds with this structure

- Cr_3S_4, Cr_3Se_4, Ni_3Sn_4, Fe_3Se_4, Ni_3Se_4, Ti_3Se_4, and V_3Se_4

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($2a$)	Ni I
$\mathbf{B}_2$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	($4i$)	Ni II
$\mathbf{B}_3$	=	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} - z_2 c\sin\beta\,\hat{\mathbf{z}}$	($4i$)	Ni II
$\mathbf{B}_4$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	($4i$)	Sn I
$\mathbf{B}_5$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	($4i$)	Sn I
$\mathbf{B}_6$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	($4i$)	Sn II
$\mathbf{B}_7$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	($4i$)	Sn II

References:

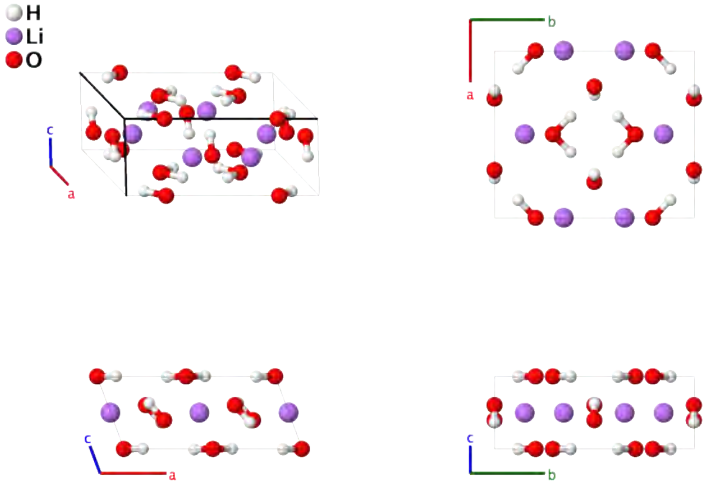
- W. Jeitschko and B. Jaberg, *Structure refinement of Ni_3Sn_4*, Acta Crystallogr. Sect. B Struct. Sci. **38**, 598–600 (1982), doi:10.1107/S056774088200346X.

Geometry files:

- CIF: pp. 1527
- POSCAR: pp. 1528

$LiOH{\cdot}H_2O$ ($B36$) Structure: A3BC2_mC24_12_ij_h_gi

http://aflow.org/prototype-encyclopedia/A3BC2_mC24_12_ij_h_gi

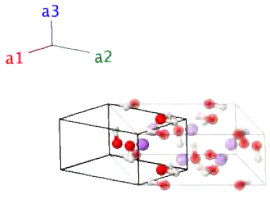


Prototype	:	H_3LiO_2
AFLOW prototype label	:	A3BC2_mC24_12_ij_h_gi
***Strukturbericht* designation**	:	$B36$
Pearson symbol	:	mC24
Space group number	:	12
Space group symbol	:	$C2/m$
AFLOW prototype command	:	`aflow --proto=A3BC2_mC24_12_ij_h_gi --params=`$a, b/a, c/a, \beta, y_1, y_2, x_3, z_3, x_4, z_4, x_5, y_5, z_5$

- The structure given in *Strukturbericht* (Herrmann, 1943) does not give the hydrogen positions, instead giving positions for Li, OH and H_2O.

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2$	=	$y_1 b\,\hat{\mathbf{y}}$	$(4g)$	O I
$\mathbf{B}_2$	=	$y_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2$	=	$-y_1 b\,\hat{\mathbf{y}}$	$(4g)$	O I

$\mathbf{B}_3$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	(4*h*)	Li
$\mathbf{B}_4$	=	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	(4*h*)	Li
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	H I
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	H I
$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	O II
$\mathbf{B}_8$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	O II
$\mathbf{B}_9$	=	$(x_5 - y_5)\,\mathbf{a}_1 + (x_5 + y_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	H II
$\mathbf{B}_{10}$	=	$(-x_5 - y_5)\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	H II
$\mathbf{B}_{11}$	=	$(-x_5 + y_5)\,\mathbf{a}_1 + (-x_5 - y_5)\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	H II
$\mathbf{B}_{12}$	=	$(x_5 + y_5)\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	H II

References:

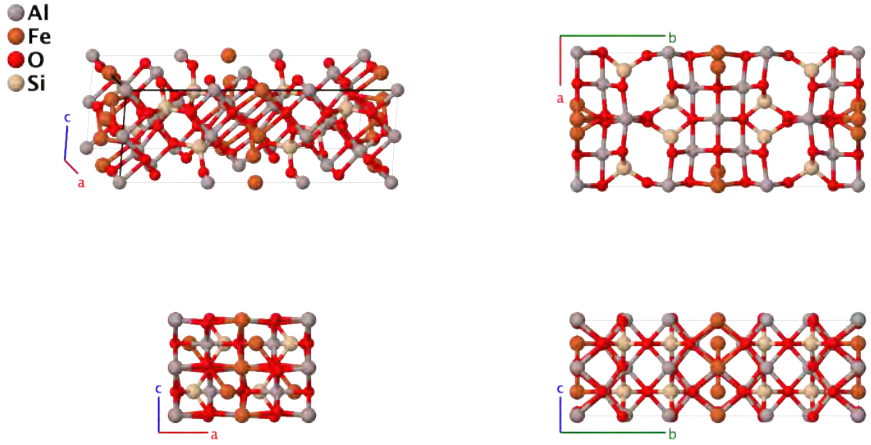
- N. W. Alcock, *Refinement of the crystal structure of lithium hydroxide monohydrate*, Acta Crystallogr. Sect. B Struct. Sci. **27**, 1682–1683 (1971), doi:10.1107/S056774087100459X.
- K. Herrmann, ed., *Strukturbericht Band VII 1939* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1943).

Geometry files:

- CIF: pp. 1528
- POSCAR: pp. 1528

Staurolite ($Al_5Fe_2O_{10}(OH)_2Si_2$) Structure: A5B2C10D2E2_mC84_12_acghj_bdi_5j_2i_j

http://aflow.org/prototype-encyclopedia/A5B2C10D2E2_mC84_12_acghj_bdi_5j_2i_j

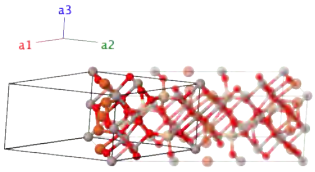


Prototype	:	$Al_5Fe_2O_{10}(OH)_2Si_2$
AFLOW prototype label	:	A5B2C10D2E2_mC84_12_acghj_bdi_5j_2i_j
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC84
Space group number	:	12
Space group symbol	:	$C2/m$
AFLOW prototype command	:	`aflow --proto=A5B2C10D2E2_mC84_12_acghj_bdi_5j_2i_j --params=`$a, b/a, c/a, \beta, y_5, y_6, x_7, z_7, x_8, z_8, x_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}$

- The orthorhombic structure of staurolite determined by (Náray-Szabó, 1929) was given the *Strukturbericht* designation $S0_4$ by (Hermann, 1937). (Smith, 1968) showed that the structure is actually monoclinic with $\beta \approx 90°$, rather than orthorhombic. This paper also corrected the chemical composition of the mineral.
- The hydrogen positions are undetermined, but they are part of a "complex distribution of OH ions," and are "probably" associated with the atoms on the (4*i*) sites (Smith, 1968). We therefore label the (4*i*) sites as OH.
- The metallic sites are actually somewhat disordered. (Smith, 1968) gives the composition of the various sites as
  - Al (2*a*) $Al_{0.67}$ $Fe_{0.33}$
  - Fe (2*b*) $Fe_{0.68}$ $Mn_{0.32}$
  - Al (2*c*) $Al_{0.67}$ $Fe_{0.33}$
  - Fe (2*d*) $Fe_{0.68}$ $Mn_{0.32}$
  - Al (4*g*) $Al_{0.95}$ $Mg_{0.05}$
  - Al (4*h*) $Al_{0.95}$ $Mg_{0.05}$
  - Fe (4*i*) $Fe_{0.64}$ $Al_{0.32}$ $Ti_{0.04}$
  - Si (8*j*) $Si_{0.936}$ $Al_{0.064}$
- (Donnay, 1983) presents a history of the difficulties in determining the staurolite structure.

Base-centered Monoclinic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2}\,b\,\hat{\mathbf{y}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,b\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	Al I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	$(2b)$	Fe I
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	$(2c)$	Al II
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(a + c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	$(2d)$	Fe II
$\mathbf{B}_5$	=	$-y_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2$	=	$y_5 b\,\hat{\mathbf{y}}$	$(4g)$	Al III
$\mathbf{B}_6$	=	$y_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2$	=	$-y_5 b\,\hat{\mathbf{y}}$	$(4g)$	Al III
$\mathbf{B}_7$	=	$-y_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	$(4h)$	Al IV
$\mathbf{B}_8$	=	$y_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	$(4h)$	Al IV
$\mathbf{B}_9$	=	$x_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Fe III
$\mathbf{B}_{10}$	=	$-x_7\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} - z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Fe III
$\mathbf{B}_{11}$	=	$x_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	OH I
$\mathbf{B}_{12}$	=	$-x_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} - z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	OH I
$\mathbf{B}_{13}$	=	$x_9\,\mathbf{a}_1 + x_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$(x_9 a + z_9 c\cos\beta)\,\hat{\mathbf{x}} + z_9 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	OH II
$\mathbf{B}_{14}$	=	$-x_9\,\mathbf{a}_1 - x_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$(-x_9 a - z_9 c\cos\beta)\,\hat{\mathbf{x}} - z_9 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	OH II
$\mathbf{B}_{15}$	=	$(x_{10} - y_{10})\,\mathbf{a}_1 + (x_{10} + y_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$(x_{10} a + z_{10} c\cos\beta)\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Al V
$\mathbf{B}_{16}$	=	$(-x_{10} - y_{10})\,\mathbf{a}_1 + (-x_{10} + y_{10})\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3$	=	$(-x_{10} a - z_{10} c\cos\beta)\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} - z_{10} c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Al V
$\mathbf{B}_{17}$	=	$(-x_{10} + y_{10})\,\mathbf{a}_1 + (-x_{10} - y_{10})\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3$	=	$(-x_{10} a - z_{10} c\cos\beta)\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} - z_{10} c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Al V
$\mathbf{B}_{18}$	=	$(x_{10} + y_{10})\,\mathbf{a}_1 + (x_{10} - y_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$(x_{10} a + z_{10} c\cos\beta)\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} + z_{10} c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Al V
$\mathbf{B}_{19}$	=	$(x_{11} - y_{11})\,\mathbf{a}_1 + (x_{11} + y_{11})\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$(x_{11} a + z_{11} c\cos\beta)\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + z_{11} c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O I
$\mathbf{B}_{20}$	=	$(-x_{11} - y_{11})\,\mathbf{a}_1 + (-x_{11} + y_{11})\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3$	=	$(-x_{11} a - z_{11} c\cos\beta)\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} - z_{11} c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O I
$\mathbf{B}_{21}$	=	$(-x_{11} + y_{11})\,\mathbf{a}_1 + (-x_{11} - y_{11})\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3$	=	$(-x_{11} a - z_{11} c\cos\beta)\,\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} - z_{11} c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O I
$\mathbf{B}_{22}$	=	$(x_{11} + y_{11})\,\mathbf{a}_1 + (x_{11} - y_{11})\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$(x_{11} a + z_{11} c\cos\beta)\,\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} + z_{11} c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O I
$\mathbf{B}_{23}$	=	$(x_{12} - y_{12})\,\mathbf{a}_1 + (x_{12} + y_{12})\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3$	=	$(x_{12} a + z_{12} c\cos\beta)\,\hat{\mathbf{x}} + y_{12} b\,\hat{\mathbf{y}} + z_{12} c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O II
$\mathbf{B}_{24}$	=	$(-x_{12} - y_{12})\,\mathbf{a}_1 + (-x_{12} + y_{12})\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3$	=	$(-x_{12} a - z_{12} c\cos\beta)\,\hat{\mathbf{x}} + y_{12} b\,\hat{\mathbf{y}} - z_{12} c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O II

$\mathbf{B}_{25}$	=	$(-x_{12}+y_{12})\,\mathbf{a}_1+(-x_{12}-y_{12})\,\mathbf{a}_2-z_{12}\,\mathbf{a}_3$	=	$(-x_{12}a-z_{12}c\cos\beta)\,\hat{\mathbf{x}}-y_{12}b\,\hat{\mathbf{y}}-z_{12}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O II
$\mathbf{B}_{26}$	=	$(x_{12}+y_{12})\,\mathbf{a}_1+(x_{12}-y_{12})\,\mathbf{a}_2+z_{12}\,\mathbf{a}_3$	=	$(x_{12}a+z_{12}c\cos\beta)\,\hat{\mathbf{x}}-y_{12}b\,\hat{\mathbf{y}}+z_{12}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O II
$\mathbf{B}_{27}$	=	$(x_{13}-y_{13})\,\mathbf{a}_1+(x_{13}+y_{13})\,\mathbf{a}_2+z_{13}\,\mathbf{a}_3$	=	$(x_{13}a+z_{13}c\cos\beta)\,\hat{\mathbf{x}}+y_{13}b\,\hat{\mathbf{y}}+z_{13}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O III
$\mathbf{B}_{28}$	=	$(-x_{13}-y_{13})\,\mathbf{a}_1+(-x_{13}+y_{13})\,\mathbf{a}_2-z_{13}\,\mathbf{a}_3$	=	$(-x_{13}a-z_{13}c\cos\beta)\,\hat{\mathbf{x}}+y_{13}b\,\hat{\mathbf{y}}-z_{13}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O III
$\mathbf{B}_{29}$	=	$(-x_{13}+y_{13})\,\mathbf{a}_1+(-x_{13}-y_{13})\,\mathbf{a}_2-z_{13}\,\mathbf{a}_3$	=	$(-x_{13}a-z_{13}c\cos\beta)\,\hat{\mathbf{x}}-y_{13}b\,\hat{\mathbf{y}}-z_{13}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O III
$\mathbf{B}_{30}$	=	$(x_{13}+y_{13})\,\mathbf{a}_1+(x_{13}-y_{13})\,\mathbf{a}_2+z_{13}\,\mathbf{a}_3$	=	$(x_{13}a+z_{13}c\cos\beta)\,\hat{\mathbf{x}}-y_{13}b\,\hat{\mathbf{y}}+z_{13}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O III
$\mathbf{B}_{31}$	=	$(x_{14}-y_{14})\,\mathbf{a}_1+(x_{14}+y_{14})\,\mathbf{a}_2+z_{14}\,\mathbf{a}_3$	=	$(x_{14}a+z_{14}c\cos\beta)\,\hat{\mathbf{x}}+y_{14}b\,\hat{\mathbf{y}}+z_{14}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O IV
$\mathbf{B}_{32}$	=	$(-x_{14}-y_{14})\,\mathbf{a}_1+(-x_{14}+y_{14})\,\mathbf{a}_2-z_{14}\,\mathbf{a}_3$	=	$(-x_{14}a-z_{14}c\cos\beta)\,\hat{\mathbf{x}}+y_{14}b\,\hat{\mathbf{y}}-z_{14}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O IV
$\mathbf{B}_{33}$	=	$(-x_{14}+y_{14})\,\mathbf{a}_1+(-x_{14}-y_{14})\,\mathbf{a}_2-z_{14}\,\mathbf{a}_3$	=	$(-x_{14}a-z_{14}c\cos\beta)\,\hat{\mathbf{x}}-y_{14}b\,\hat{\mathbf{y}}-z_{14}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O IV
$\mathbf{B}_{34}$	=	$(x_{14}+y_{14})\,\mathbf{a}_1+(x_{14}-y_{14})\,\mathbf{a}_2+z_{14}\,\mathbf{a}_3$	=	$(x_{14}a+z_{14}c\cos\beta)\,\hat{\mathbf{x}}-y_{14}b\,\hat{\mathbf{y}}+z_{14}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O IV
$\mathbf{B}_{35}$	=	$(x_{15}-y_{15})\,\mathbf{a}_1+(x_{15}+y_{15})\,\mathbf{a}_2+z_{15}\,\mathbf{a}_3$	=	$(x_{15}a+z_{15}c\cos\beta)\,\hat{\mathbf{x}}+y_{15}b\,\hat{\mathbf{y}}+z_{15}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O V
$\mathbf{B}_{36}$	=	$(-x_{15}-y_{15})\,\mathbf{a}_1+(-x_{15}+y_{15})\,\mathbf{a}_2-z_{15}\,\mathbf{a}_3$	=	$(-x_{15}a-z_{15}c\cos\beta)\,\hat{\mathbf{x}}+y_{15}b\,\hat{\mathbf{y}}-z_{15}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O V
$\mathbf{B}_{37}$	=	$(-x_{15}+y_{15})\,\mathbf{a}_1+(-x_{15}-y_{15})\,\mathbf{a}_2-z_{15}\,\mathbf{a}_3$	=	$(-x_{15}a-z_{15}c\cos\beta)\,\hat{\mathbf{x}}-y_{15}b\,\hat{\mathbf{y}}-z_{15}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O V
$\mathbf{B}_{38}$	=	$(x_{15}+y_{15})\,\mathbf{a}_1+(x_{15}-y_{15})\,\mathbf{a}_2+z_{15}\,\mathbf{a}_3$	=	$(x_{15}a+z_{15}c\cos\beta)\,\hat{\mathbf{x}}-y_{15}b\,\hat{\mathbf{y}}+z_{15}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O V
$\mathbf{B}_{39}$	=	$(x_{16}-y_{16})\,\mathbf{a}_1+(x_{16}+y_{16})\,\mathbf{a}_2+z_{16}\,\mathbf{a}_3$	=	$(x_{16}a+z_{16}c\cos\beta)\,\hat{\mathbf{x}}+y_{16}b\,\hat{\mathbf{y}}+z_{16}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Si
$\mathbf{B}_{40}$	=	$(-x_{16}-y_{16})\,\mathbf{a}_1+(-x_{16}+y_{16})\,\mathbf{a}_2-z_{16}\,\mathbf{a}_3$	=	$(-x_{16}a-z_{16}c\cos\beta)\,\hat{\mathbf{x}}+y_{16}b\,\hat{\mathbf{y}}-z_{16}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Si
$\mathbf{B}_{41}$	=	$(-x_{16}+y_{16})\,\mathbf{a}_1+(-x_{16}-y_{16})\,\mathbf{a}_2-z_{16}\,\mathbf{a}_3$	=	$(-x_{16}a-z_{16}c\cos\beta)\,\hat{\mathbf{x}}-y_{16}b\,\hat{\mathbf{y}}-z_{16}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Si
$\mathbf{B}_{42}$	=	$(x_{16}+y_{16})\,\mathbf{a}_1+(x_{16}-y_{16})\,\mathbf{a}_2+z_{16}\,\mathbf{a}_3$	=	$(x_{16}a+z_{16}c\cos\beta)\,\hat{\mathbf{x}}-y_{16}b\,\hat{\mathbf{y}}+z_{16}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Si

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- J. D. H. Donnay and G. Donnay, *The staurolite story*, Tschermaks Min. Petr. Mitt. **31**, 1–15 (1983), doi:10.1007/BF01084757.

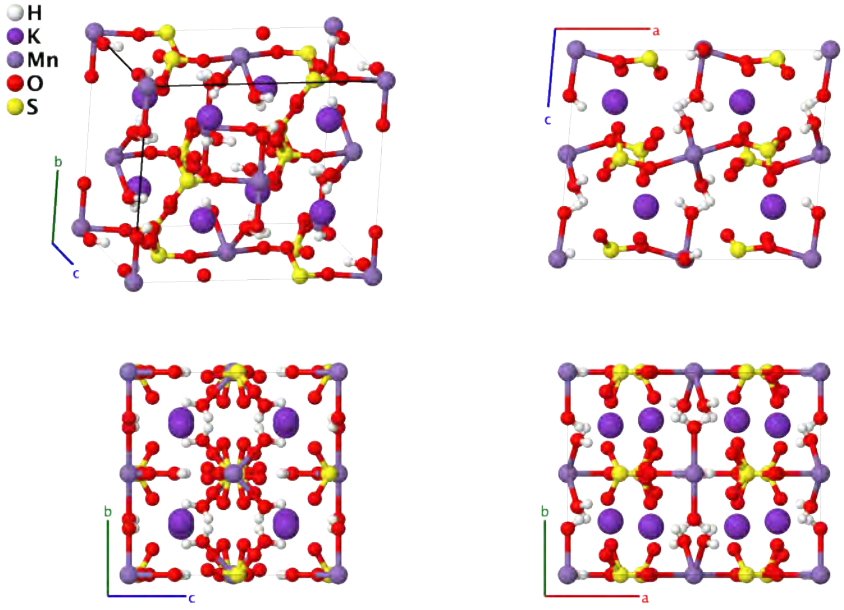
Geometry files:

- CIF: pp. 1528

- POSCAR: pp. 1529

Manganese-leonite [$K_2Mn(SO_4)_2{\cdot}4H_2O$, $H4_{23}$] Structure: A8B2CD15E2_mC112_12_2i3j_j_ad_g4i5j_2i

http://aflow.org/prototype-encyclopedia/A8B2CD15E2_mC112_12_2i3j_j_ad_g4i5j_2i



Prototype	:	$H_8K_2MnO_{12}S_2$
AFLOW prototype label	:	A8B2CD15E2_mC112_12_2i3j_j_ad_g4i5j_2i
***Strukturbericht* designation**	:	$H4_{23}$
Pearson symbol	:	mC112
Space group number	:	12
Space group symbol	:	$C2/m$
AFLOW prototype command	:	`aflow --proto=A8B2CD15E2_mC112_12_2i3j_j_ad_g4i5j_2i --params=`$a, b/a, c/a, \beta, y_3, x_4, z_4, x_5, z_5, x_6, z_6, x_7, z_7, x_8, z_8, x_9, z_9, x_{10}, z_{10}, x_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}, x_{17}, y_{17}, z_{17}, x_{18}, y_{18}, z_{18}, x_{19}, y_{19}, z_{19}, x_{20}, y_{20}, z_{20}$

Other compounds with this structure

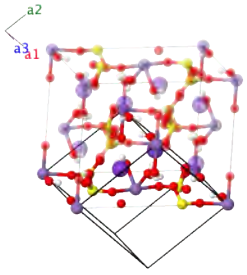
- $(NH_2)_2Co(SO_4)_2{\cdot}4H_2O$, $(NH_2)_2Cu(SO_4)_2{\cdot}4H_2O$, $(NH_2)_2Fe(SO_4)_2{\cdot}4H_2O$, $(NH_2)_2Mg(SO_4)_2{\cdot}4H_2O$, $(NH_2)_2Mn(SO_4)_2{\cdot}4H_2O$, $Cd_2Co(SO_4)_2{\cdot}4H_2O$, $Cd_2Cu(SO_4)_2{\cdot}4H_2O$, $Cd_2Fe(SO_4)_2{\cdot}4H_2O$, $Cd_2Mg(SO_4)_2{\cdot}4H_2O$, $Cd_2Mn(SO_4)_2{\cdot}4H_2O$, $Cs_2Co(SO_4)_2{\cdot}4H_2O$, $Cs_2Cu(SO_4)_2{\cdot}4H_2O$, $Cs_2Fe(SO_4)_2{\cdot}4H_2O$, $Cs_2Mg(SO_4)_2{\cdot}4H_2O$, $Cs_2Mn(SO_4)_2{\cdot}4H_2O$, $K_2Co(SO_4)_2{\cdot}4H_2O$, $K_2Cu(SO_4)_2{\cdot}4H_2O$, $K_2Fe(SO_4)_2{\cdot}4H_2O$ (mereiterite), $K_2Mg(SO_4)_2{\cdot}4H_2O$ (leonite), $Rb_2Co(SO_4)_2{\cdot}4H_2O$, $Rb_2Cu(SO_4)_2{\cdot}4H_2O$, $Rb_2Fe(SO_4)_2{\cdot}4H_2O$, $Rb_2Mg(SO_4)_2{\cdot}4H_2O$, $Rb_2Mn(SO_4)_2{\cdot}4H_2O$, $Tl_2Co(SO_4)_2{\cdot}4H_2O$, $Tl_2Cu(SO_4)_2{\cdot}4H_2O$, $Tl_2Fe(SO_4)_2{\cdot}4H_2O$, $Tl_2Mg(SO_4)_2{\cdot}4H_2O$, and $Tl_2Mn(SO_4)_2{\cdot}4H_2O$

- Properly, the prototype of leonite is $K_2Mg(SO_4)_2{\cdot}4H_2O$, but (Herrmann, 1943) gives manganese-leonite *Strukturbericht* symbol $H4_{23}$, so we will use Mn-leonite as the prototype.
- The room temperature leonite structure, shown here, is in space group $C2/m$ #12. It is characterized by the disorder of the second sulfate group, centered on atom S-II (Hertweck, 2001). The oxygen sites around S-II, labeled O-VII, O-VIII, and O-IX in our notation, are only occupied 50% of the time. (Anspach, 1939), who did the original determination of this structure, was not able to see the disorder, and so gives an ordered structure for both sulfates.

- We use the structure determined by (Hertweck, 2001) at room temperature, 293 K. At lower temperatures, the leonites undergo phase transitions. The exact transition temperature depends on the compound.
- At somewhat lower than room temperature (205 K for Mn-leonite) the sulfate group orders, doubling the size of the unit cell and changing the space group from $C2/m$ to $C2/c$ #15. At still lower temperatures (168 K for Mn-leonite) further ordering takes place and the structure transforms into the $P2_1/c$ #14 space group.

Base-centered Monoclinic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2}\,b\,\hat{\mathbf{y}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,b\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	Mn I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(a + c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	$(2d)$	Mn II
$\mathbf{B}_3$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2$	=	$y_3 b\,\hat{\mathbf{y}}$	$(4g)$	O I
$\mathbf{B}_4$	=	$y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2$	=	$-y_3 b\,\hat{\mathbf{y}}$	$(4g)$	O I
$\mathbf{B}_5$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	H I
$\mathbf{B}_6$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	H I
$\mathbf{B}_7$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	H II
$\mathbf{B}_8$	=	$-x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	H II
$\mathbf{B}_9$	=	$x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	O II
$\mathbf{B}_{10}$	=	$-x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	O II
$\mathbf{B}_{11}$	=	$x_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	O III
$\mathbf{B}_{12}$	=	$-x_7\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} - z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	O III
$\mathbf{B}_{13}$	=	$x_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	O IV
$\mathbf{B}_{14}$	=	$-x_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} - z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	O IV
$\mathbf{B}_{15}$	=	$x_9\,\mathbf{a}_1 + x_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$(x_9 a + z_9 c\cos\beta)\,\hat{\mathbf{x}} + z_9 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	O V
$\mathbf{B}_{16}$	=	$-x_9\,\mathbf{a}_1 - x_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$(-x_9 a - z_9 c\cos\beta)\,\hat{\mathbf{x}} - z_9 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	O V
$\mathbf{B}_{17}$	=	$x_{10}\,\mathbf{a}_1 + x_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$(x_{10} a + z_{10} c\cos\beta)\,\hat{\mathbf{x}} + z_{10} c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	S I
$\mathbf{B}_{18}$	=	$-x_{10}\,\mathbf{a}_1 - x_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3$	=	$(-x_{10} a - z_{10} c\cos\beta)\,\hat{\mathbf{x}} - z_{10} c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	S I
$\mathbf{B}_{19}$	=	$x_{11}\,\mathbf{a}_1 + x_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$(x_{11} a + z_{11} c\cos\beta)\,\hat{\mathbf{x}} + z_{11} c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	S II
$\mathbf{B}_{20}$	=	$-x_{11}\,\mathbf{a}_1 - x_{11}\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3$	=	$(-x_{11} a - z_{11} c\cos\beta)\,\hat{\mathbf{x}} - z_{11} c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	S II
$\mathbf{B}_{21}$	=	$(x_{12} - y_{12})\,\mathbf{a}_1 + (x_{12} + y_{12})\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3$	=	$(x_{12} a + z_{12} c\cos\beta)\,\hat{\mathbf{x}} + y_{12} b\,\hat{\mathbf{y}} + z_{12} c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	H III
$\mathbf{B}_{22}$	=	$(-x_{12} - y_{12})\,\mathbf{a}_1 + (-x_{12} + y_{12})\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3$	=	$(-x_{12} a - z_{12} c\cos\beta)\,\hat{\mathbf{x}} + y_{12} b\,\hat{\mathbf{y}} - z_{12} c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	H III

$$\mathbf{B}_{23} = (-x_{12}+y_{12})\,\mathbf{a}_1 + (-x_{12}-y_{12})\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3 = (-x_{12}a - z_{12}c\cos\beta)\,\hat{\mathbf{x}} - y_{12}b\,\hat{\mathbf{y}} - z_{12}c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{H III}$$

$$\mathbf{B}_{24} = (x_{12}+y_{12})\,\mathbf{a}_1 + (x_{12}-y_{12})\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 = (x_{12}a + z_{12}c\cos\beta)\,\hat{\mathbf{x}} - y_{12}b\,\hat{\mathbf{y}} + z_{12}c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{H III}$$

$$\mathbf{B}_{25} = (x_{13}-y_{13})\,\mathbf{a}_1 + (x_{13}+y_{13})\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3 = (x_{13}a + z_{13}c\cos\beta)\,\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} + z_{13}c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{H IV}$$

$$\mathbf{B}_{26} = (-x_{13}-y_{13})\,\mathbf{a}_1 + (-x_{13}+y_{13})\,\mathbf{a}_2 - z_{13}\,\mathbf{a}_3 = (-x_{13}a - z_{13}c\cos\beta)\,\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} - z_{13}c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{H IV}$$

$$\mathbf{B}_{27} = (-x_{13}+y_{13})\,\mathbf{a}_1 + (-x_{13}-y_{13})\,\mathbf{a}_2 - z_{13}\,\mathbf{a}_3 = (-x_{13}a - z_{13}c\cos\beta)\,\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} - z_{13}c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{H IV}$$

$$\mathbf{B}_{28} = (x_{13}+y_{13})\,\mathbf{a}_1 + (x_{13}-y_{13})\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3 = (x_{13}a + z_{13}c\cos\beta)\,\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} + z_{13}c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{H IV}$$

$$\mathbf{B}_{29} = (x_{14}-y_{14})\,\mathbf{a}_1 + (x_{14}+y_{14})\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3 = (x_{14}a + z_{14}c\cos\beta)\,\hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} + z_{14}c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{H V}$$

$$\mathbf{B}_{30} = (-x_{14}-y_{14})\,\mathbf{a}_1 + (-x_{14}+y_{14})\,\mathbf{a}_2 - z_{14}\,\mathbf{a}_3 = (-x_{14}a - z_{14}c\cos\beta)\,\hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} - z_{14}c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{H V}$$

$$\mathbf{B}_{31} = (-x_{14}+y_{14})\,\mathbf{a}_1 + (-x_{14}-y_{14})\,\mathbf{a}_2 - z_{14}\,\mathbf{a}_3 = (-x_{14}a - z_{14}c\cos\beta)\,\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} - z_{14}c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{H V}$$

$$\mathbf{B}_{32} = (x_{14}+y_{14})\,\mathbf{a}_1 + (x_{14}-y_{14})\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3 = (x_{14}a + z_{14}c\cos\beta)\,\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} + z_{14}c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{H V}$$

$$\mathbf{B}_{33} = (x_{15}-y_{15})\,\mathbf{a}_1 + (x_{15}+y_{15})\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 = (x_{15}a + z_{15}c\cos\beta)\,\hat{\mathbf{x}} + y_{15}b\,\hat{\mathbf{y}} + z_{15}c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{K}$$

$$\mathbf{B}_{34} = (-x_{15}-y_{15})\,\mathbf{a}_1 + (-x_{15}+y_{15})\,\mathbf{a}_2 - z_{15}\,\mathbf{a}_3 = (-x_{15}a - z_{15}c\cos\beta)\,\hat{\mathbf{x}} + y_{15}b\,\hat{\mathbf{y}} - z_{15}c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{K}$$

$$\mathbf{B}_{35} = (-x_{15}+y_{15})\,\mathbf{a}_1 + (-x_{15}-y_{15})\,\mathbf{a}_2 - z_{15}\,\mathbf{a}_3 = (-x_{15}a - z_{15}c\cos\beta)\,\hat{\mathbf{x}} - y_{15}b\,\hat{\mathbf{y}} - z_{15}c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{K}$$

$$\mathbf{B}_{36} = (x_{15}+y_{15})\,\mathbf{a}_1 + (x_{15}-y_{15})\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 = (x_{15}a + z_{15}c\cos\beta)\,\hat{\mathbf{x}} - y_{15}b\,\hat{\mathbf{y}} + z_{15}c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{K}$$

$$\mathbf{B}_{37} = (x_{16}-y_{16})\,\mathbf{a}_1 + (x_{16}+y_{16})\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 = (x_{16}a + z_{16}c\cos\beta)\,\hat{\mathbf{x}} + y_{16}b\,\hat{\mathbf{y}} + z_{16}c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{O VI}$$

$$\mathbf{B}_{38} = (-x_{16}-y_{16})\,\mathbf{a}_1 + (-x_{16}+y_{16})\,\mathbf{a}_2 - z_{16}\,\mathbf{a}_3 = (-x_{16}a - z_{16}c\cos\beta)\,\hat{\mathbf{x}} + y_{16}b\,\hat{\mathbf{y}} - z_{16}c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{O VI}$$

$$\mathbf{B}_{39} = (-x_{16}+y_{16})\,\mathbf{a}_1 + (-x_{16}-y_{16})\,\mathbf{a}_2 - z_{16}\,\mathbf{a}_3 = (-x_{16}a - z_{16}c\cos\beta)\,\hat{\mathbf{x}} - y_{16}b\,\hat{\mathbf{y}} - z_{16}c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{O VI}$$

$$\mathbf{B}_{40} = (x_{16}+y_{16})\,\mathbf{a}_1 + (x_{16}-y_{16})\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 = (x_{16}a + z_{16}c\cos\beta)\,\hat{\mathbf{x}} - y_{16}b\,\hat{\mathbf{y}} + z_{16}c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{O VI}$$

$$\mathbf{B}_{41} = (x_{17}-y_{17})\,\mathbf{a}_1 + (x_{17}+y_{17})\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 = (x_{17}a + z_{17}c\cos\beta)\,\hat{\mathbf{x}} + y_{17}b\,\hat{\mathbf{y}} + z_{17}c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{O VII}$$

$$\mathbf{B}_{42} = (-x_{17}-y_{17})\,\mathbf{a}_1 + (-x_{17}+y_{17})\,\mathbf{a}_2 - z_{17}\,\mathbf{a}_3 = (-x_{17}a - z_{17}c\cos\beta)\,\hat{\mathbf{x}} + y_{17}b\,\hat{\mathbf{y}} - z_{17}c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{O VII}$$

$$\mathbf{B}_{43} = (-x_{17}+y_{17})\,\mathbf{a}_1 + (-x_{17}-y_{17})\,\mathbf{a}_2 - z_{17}\,\mathbf{a}_3 = (-x_{17}a - z_{17}c\cos\beta)\,\hat{\mathbf{x}} - y_{17}b\,\hat{\mathbf{y}} - z_{17}c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{O VII}$$

$$\mathbf{B}_{44} = (x_{17}+y_{17})\,\mathbf{a}_1 + (x_{17}-y_{17})\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 = (x_{17}a + z_{17}c\cos\beta)\,\hat{\mathbf{x}} - y_{17}b\,\hat{\mathbf{y}} + z_{17}c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{O VII}$$

$$\mathbf{B}_{45} = (x_{18}-y_{18})\,\mathbf{a}_1 + (x_{18}+y_{18})\,\mathbf{a}_2 + z_{18}\,\mathbf{a}_3 = (x_{18}a + z_{18}c\cos\beta)\,\hat{\mathbf{x}} + y_{18}b\,\hat{\mathbf{y}} + z_{18}c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{O VIII}$$

$$\mathbf{B}_{46} = (-x_{18}-y_{18})\,\mathbf{a}_1 + (-x_{18}+y_{18})\,\mathbf{a}_2 - z_{18}\,\mathbf{a}_3 = (-x_{18}a - z_{18}c\cos\beta)\,\hat{\mathbf{x}} + y_{18}b\,\hat{\mathbf{y}} - z_{18}c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{O VIII}$$

$\mathbf{B}_{47}$	=	$(-x_{18}+y_{18})\,\mathbf{a}_1 + (-x_{18}-y_{18})\,\mathbf{a}_2 - z_{18}\,\mathbf{a}_3$	=	$(-x_{18}a - z_{18}c\cos\beta)\,\hat{\mathbf{x}} - y_{18}b\,\hat{\mathbf{y}} - z_{18}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O VIII
$\mathbf{B}_{48}$	=	$(x_{18}+y_{18})\,\mathbf{a}_1 + (x_{18}-y_{18})\,\mathbf{a}_2 + z_{18}\,\mathbf{a}_3$	=	$(x_{18}a + z_{18}c\cos\beta)\,\hat{\mathbf{x}} - y_{18}b\,\hat{\mathbf{y}} + z_{18}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O VIII
$\mathbf{B}_{49}$	=	$(x_{19}-y_{19})\,\mathbf{a}_1 + (x_{19}+y_{19})\,\mathbf{a}_2 + z_{19}\,\mathbf{a}_3$	=	$(x_{19}a + z_{19}c\cos\beta)\,\hat{\mathbf{x}} + y_{19}b\,\hat{\mathbf{y}} + z_{19}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O IX
$\mathbf{B}_{50}$	=	$(-x_{19}-y_{19})\,\mathbf{a}_1 + (-x_{19}+y_{19})\,\mathbf{a}_2 - z_{19}\,\mathbf{a}_3$	=	$(-x_{19}a - z_{19}c\cos\beta)\,\hat{\mathbf{x}} + y_{19}b\,\hat{\mathbf{y}} - z_{19}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O IX
$\mathbf{B}_{51}$	=	$(-x_{19}+y_{19})\,\mathbf{a}_1 + (-x_{19}-y_{19})\,\mathbf{a}_2 - z_{19}\,\mathbf{a}_3$	=	$(-x_{19}a - z_{19}c\cos\beta)\,\hat{\mathbf{x}} - y_{19}b\,\hat{\mathbf{y}} - z_{19}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O IX
$\mathbf{B}_{52}$	=	$(x_{19}+y_{19})\,\mathbf{a}_1 + (x_{19}-y_{19})\,\mathbf{a}_2 + z_{19}\,\mathbf{a}_3$	=	$(x_{19}a + z_{19}c\cos\beta)\,\hat{\mathbf{x}} - y_{19}b\,\hat{\mathbf{y}} + z_{19}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O IX
$\mathbf{B}_{53}$	=	$(x_{20}-y_{20})\,\mathbf{a}_1 + (x_{20}+y_{20})\,\mathbf{a}_2 + z_{20}\,\mathbf{a}_3$	=	$(x_{20}a + z_{20}c\cos\beta)\,\hat{\mathbf{x}} + y_{20}b\,\hat{\mathbf{y}} + z_{20}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O X
$\mathbf{B}_{54}$	=	$(-x_{20}-y_{20})\,\mathbf{a}_1 + (-x_{20}+y_{20})\,\mathbf{a}_2 - z_{20}\,\mathbf{a}_3$	=	$(-x_{20}a - z_{20}c\cos\beta)\,\hat{\mathbf{x}} + y_{20}b\,\hat{\mathbf{y}} - z_{20}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O X
$\mathbf{B}_{55}$	=	$(-x_{20}+y_{20})\,\mathbf{a}_1 + (-x_{20}-y_{20})\,\mathbf{a}_2 - z_{20}\,\mathbf{a}_3$	=	$(-x_{20}a - z_{20}c\cos\beta)\,\hat{\mathbf{x}} - y_{20}b\,\hat{\mathbf{y}} - z_{20}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O X
$\mathbf{B}_{56}$	=	$(x_{20}+y_{20})\,\mathbf{a}_1 + (x_{20}-y_{20})\,\mathbf{a}_2 + z_{20}\,\mathbf{a}_3$	=	$(x_{20}a + z_{20}c\cos\beta)\,\hat{\mathbf{x}} - y_{20}b\,\hat{\mathbf{y}} + z_{20}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O X

References:

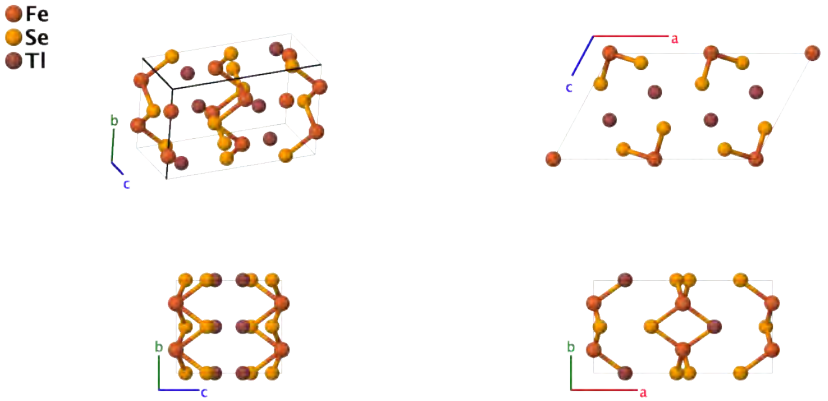
- B. Hertweck, G. Giester, and E. Libowitzky, *The crystal structures of the low-temperature phases of leonite-type compounds, $K_2Me(SO_4)_2{\cdot}4H_2O$ (Me^{2+} = Mg, Mn, Fe)*, Am. Mineral. **86**, 1282–1292 (2001), doi:10.2138/am-2001-1016.
- H. Anspach, *Die Struktur von Mn-Leonit*, Zeitschrift für Kristallographie - Crystalline Materials **101**, 39–77 (1939), doi:10.1524/zkri.1939.101.1.39.
- K. Herrmann, ed., *Strukturbericht Band VII 1939* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1943).

Geometry files:

- CIF: pp. 1529
- POSCAR: pp. 1529

Monoclinic $FeTlSe_2$ Structure: AB2C_mC16_12_g_2i_i

http://aflow.org/prototype-encyclopedia/AB2C_mC16_12_g_2i_i



Prototype	:	$FeSe_2Tl$
AFLOW prototype label	:	AB2C_mC16_12_g_2i_i
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC16
Space group number	:	12
Space group symbol	:	$C2/m$
AFLOW prototype command	:	`aflow --proto=AB2C_mC16_12_g_2i_i --params=`$a, b/a, c/a, \beta, y_1, x_2, z_2, x_3, z_3, x_4, z_4$

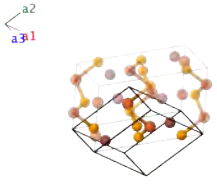
Other compounds with this structure

- $FeTlS_2$

- This is the low temperature structure of $FeTlSe_2$/$FeTlS_2$. At high temperatures, it transforms to the tetragonal $FeTlS_2$ structure.

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2$	=	$y_1 b\,\hat{\mathbf{y}}$	$(4g)$	Fe
$\mathbf{B}_2$	=	$y_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2$	=	$-y_1 b\,\hat{\mathbf{y}}$	$(4g)$	Fe
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Se I
$\mathbf{B}_4$	=	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} - z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Se I

$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Se II
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Se II
$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Tl
$\mathbf{B}_8$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Tl

References:

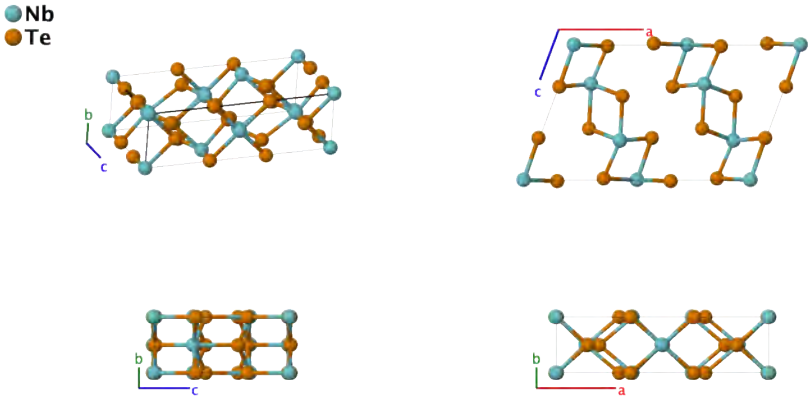
- K. Klepp and H. Boller, *Die Kristallstruktur von TlFeSe$_2$ und T1FeS$_2$*, Monatshefte für Chemie - Chemical Monthly **110**, 1045–1055 (1979), doi:10.1007/BF00910952.

Geometry files:

- CIF: pp. 1530
- POSCAR: pp. 1530

$NbTe_2$ Structure: AB2_mC18_12_ai_3i

http://aflow.org/prototype-encyclopedia/AB2_mC18_12_ai_3i



Prototype	:	$NbTe_2$
AFLOW prototype label	:	AB2_mC18_12_ai_3i
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC18
Space group number	:	12
Space group symbol	:	$C2/m$
AFLOW prototype command	:	`aflow --proto=AB2_mC18_12_ai_3i --params=`$a, b/a, c/a, \beta, x_2, z_2, x_3, z_3, x_4, z_4, x_5, z_5$

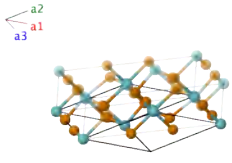
Other compounds with this structure

- $TaTe_2$

- FINDSYM found a somewhat more compact representation of the unit cell (smaller a and β) than that given in (Brown, 1966), so the lattice constants and Wyckoff parameters differ from those given in the reference.

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	Nb I
$\mathbf{B}_2$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Nb II
$\mathbf{B}_3$	=	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} - z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Nb II

$\mathbf{B}_4$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Te I
$\mathbf{B}_5$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Te I
$\mathbf{B}_6$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Te II
$\mathbf{B}_7$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Te II
$\mathbf{B}_8$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Te III
$\mathbf{B}_9$	=	$-x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Te III

References:

- B. E. Brown, *The Crystal Structures of $NbTe_2$ and $TaTe_2$*, Acta Cryst. **20**, 264–267 (1966), doi:10.1107/S0365110X66000501.

Found in:

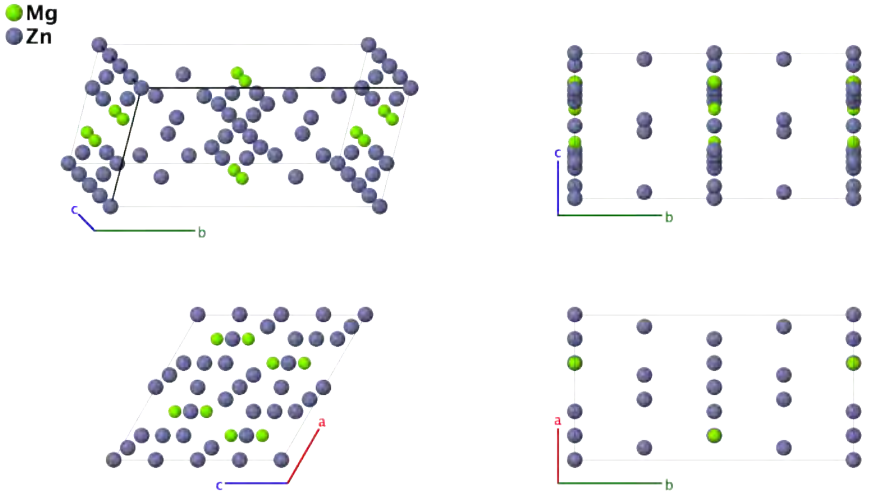
- P. Villars and L. Calvert, *Pearson's Handbook of Crystallographic Data for Intermetallic Phases* (ASM International, Materials Park, OH, 1991), 2nd edn.

Geometry files:

- CIF: pp. 1530
- POSCAR: pp. 1531

$D2_2$ ($MgZn_5$?) (*problematic*) Structure: AB5_mC48_12_2i_ac5i2j

http://aflow.org/prototype-encyclopedia/AB5_mC48_12_2i_ac5i2j

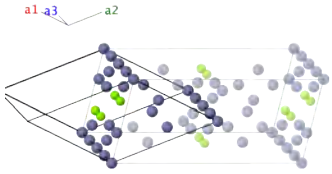


Prototype	:	$MgZn_5$
AFLOW prototype label	:	AB5_mC48_12_2i_ac5i2j
***Strukturbericht* designation**	:	$D2_2$
Pearson symbol	:	mC48
Space group number	:	12
Space group symbol	:	$C2/m$
AFLOW prototype command	:	`aflow --proto=AB5_mC48_12_2i_ac5i2j --params=`$a, b/a, c/a, \beta, x_3, z_3, x_4, z_4, x_5, z_5, x_6, z_6, x_7, z_7, x_8, z_8, x_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}$

- This structure has problems similar to the *B*30 MgZn structure, and for the same reasons. (Hermann, 1937) assigned this the *Strukturbericht* designation $D2_2$, based on the paper of (Tarschish, 1933), who derived it from the hexagonal Laves structure $MgZn_2$ (C14), eventually resulting in a 48 atom cell with composition $MgZn_5$. As with MgZn, he assumed that the space group remained $P6_3/mmc$ #194.
- (McKeehan, 1935) again pointed out that this is impossible. (Hermann, 1937) referenced both papers, giving the space group as $P6_3/mmc$ but listing the atomic coordinates enumerated by McKeehan.
- The McKeehan structure has space group $C2/m$ #12, with 48 atoms in the conventional cell, and 24 atoms in the primitive cell. As with *B*30, this agrees with the structure (Parthé, 1993) designate as $D2_2$.
- It is not clear that any $MgZn_5$ compound actually exists. It does not appear in the assessed Mg-Zn binary phase diagram (Massalski, 1990). It *may* actually be the Mg_2Zn_{11} $D8_c$ structure, but we have found no literature supporting this claim.

Base-centered Monoclinic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2}\,b\,\hat{\mathbf{y}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,b\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	Zn I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	$(2c)$	Zn II
$\mathbf{B}_3$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3a + z_3c\cos\beta)\,\hat{\mathbf{x}} + z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Mg I
$\mathbf{B}_4$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3a - z_3c\cos\beta)\,\hat{\mathbf{x}} - z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Mg I
$\mathbf{B}_5$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4a + z_4c\cos\beta)\,\hat{\mathbf{x}} + z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Mg II
$\mathbf{B}_6$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4a - z_4c\cos\beta)\,\hat{\mathbf{x}} - z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Mg II
$\mathbf{B}_7$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5a + z_5c\cos\beta)\,\hat{\mathbf{x}} + z_5c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Zn III
$\mathbf{B}_8$	=	$-x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5a - z_5c\cos\beta)\,\hat{\mathbf{x}} - z_5c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Zn III
$\mathbf{B}_9$	=	$x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6a + z_6c\cos\beta)\,\hat{\mathbf{x}} + z_6c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Zn IV
$\mathbf{B}_{10}$	=	$-x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6a - z_6c\cos\beta)\,\hat{\mathbf{x}} - z_6c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Zn IV
$\mathbf{B}_{11}$	=	$x_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7a + z_7c\cos\beta)\,\hat{\mathbf{x}} + z_7c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Zn V
$\mathbf{B}_{12}$	=	$-x_7\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7a - z_7c\cos\beta)\,\hat{\mathbf{x}} - z_7c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Zn V
$\mathbf{B}_{13}$	=	$x_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8a + z_8c\cos\beta)\,\hat{\mathbf{x}} + z_8c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Zn VI
$\mathbf{B}_{14}$	=	$-x_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8a - z_8c\cos\beta)\,\hat{\mathbf{x}} - z_8c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Zn VI
$\mathbf{B}_{15}$	=	$x_9\,\mathbf{a}_1 + x_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$(x_9a + z_9c\cos\beta)\,\hat{\mathbf{x}} + z_9c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Zn VII
$\mathbf{B}_{16}$	=	$-x_9\,\mathbf{a}_1 - x_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$(-x_9a - z_9c\cos\beta)\,\hat{\mathbf{x}} - z_9c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Zn VII
$\mathbf{B}_{17}$	=	$(x_{10} - y_{10})\,\mathbf{a}_1 + (x_{10} + y_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$(x_{10}a + z_{10}c\cos\beta)\,\hat{\mathbf{x}} + y_{10}b\,\hat{\mathbf{y}} + z_{10}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Zn VIII
$\mathbf{B}_{18}$	=	$(-x_{10} - y_{10})\,\mathbf{a}_1 + (-x_{10} + y_{10})\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3$	=	$(-x_{10}a - z_{10}c\cos\beta)\,\hat{\mathbf{x}} + y_{10}b\,\hat{\mathbf{y}} - z_{10}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Zn VIII
$\mathbf{B}_{19}$	=	$(-x_{10} + y_{10})\,\mathbf{a}_1 + (-x_{10} - y_{10})\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3$	=	$(-x_{10}a - z_{10}c\cos\beta)\,\hat{\mathbf{x}} - y_{10}b\,\hat{\mathbf{y}} - z_{10}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Zn VIII
$\mathbf{B}_{20}$	=	$(x_{10} + y_{10})\,\mathbf{a}_1 + (x_{10} - y_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$(x_{10}a + z_{10}c\cos\beta)\,\hat{\mathbf{x}} - y_{10}b\,\hat{\mathbf{y}} + z_{10}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Zn VIII
$\mathbf{B}_{21}$	=	$(x_{11} - y_{11})\,\mathbf{a}_1 + (x_{11} + y_{11})\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$(x_{11}a + z_{11}c\cos\beta)\,\hat{\mathbf{x}} + y_{11}b\,\hat{\mathbf{y}} + z_{11}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Zn IX
$\mathbf{B}_{22}$	=	$(-x_{11} - y_{11})\,\mathbf{a}_1 + (-x_{11} + y_{11})\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3$	=	$(-x_{11}a - z_{11}c\cos\beta)\,\hat{\mathbf{x}} + y_{11}b\,\hat{\mathbf{y}} - z_{11}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Zn IX
$\mathbf{B}_{23}$	=	$(-x_{11} + y_{11})\,\mathbf{a}_1 + (-x_{11} - y_{11})\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3$	=	$(-x_{11}a - z_{11}c\cos\beta)\,\hat{\mathbf{x}} - y_{11}b\,\hat{\mathbf{y}} - z_{11}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Zn IX
$\mathbf{B}_{24}$	=	$(x_{11} + y_{11})\,\mathbf{a}_1 + (x_{11} - y_{11})\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$(x_{11}a + z_{11}c\cos\beta)\,\hat{\mathbf{x}} - y_{11}b\,\hat{\mathbf{y}} + z_{11}c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Zn IX

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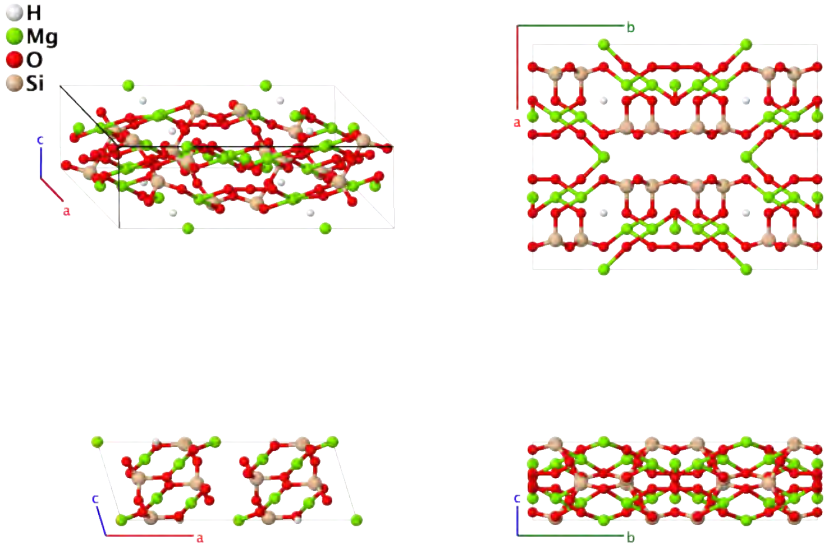
Geometry files:

- CIF: pp. 1531
- POSCAR: pp. 1531

Chrysotile ($H_4Mg_3Si_2O_9$, $S4_5$) Structure: AB6C11D6E4_mC112_12_e_gi2j_i5j_2i2j_2j

http://aflow.org/prototype-encyclopedia/AB6C11D6E4_mC112_12_e_gi2j_i5j_2i2j_2j

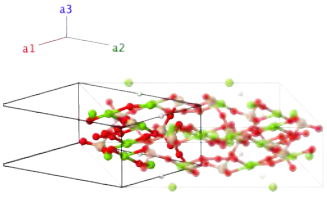


Prototype	:	$(H_2O)(OH)_6O_{11}Mg_6Si_4$
AFLOW prototype label	:	AB6C11D6E4_mC112_12_e_gi2j_i5j_2i2j_2j
***Strukturbericht* designation**	:	$S4_5$
Pearson symbol	:	mC112
Space group number	:	12
Space group symbol	:	$C2/m$
AFLOW prototype command	:	`aflow --proto=AB6C11D6E4_mC112_12_e_gi2j_i5j_2i2j_2j --params=`$a, b/a, c/a, \beta, y_2, x_3, z_3, x_4, z_4, x_5, z_5, x_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}, x_{17}, y_{17}, z_{17}$

- Chrysotile is more commonly known as "white asbestos." Chrysotile sheets typically curl into tubular fibers and the crystal structure is difficult to determine. (Yada, 1967) has a partial list of the experiments performed to determine this structure. The current structure was given the *Strukturbericht* designation $S4_5$ by (Hermann, 1937). It doubles the unit cell of our other chrysotile structure, by (Falini, 2004) and includes an inversion site.
- The structure was originally given in the $I2/m$ setting of space group #12. We used FINDSYM to place it in the standard $C2/m$ setting.

Base-centered Monoclinic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2}\,b\,\hat{\mathbf{y}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,b\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$\frac{1}{2}\,\mathbf{a}_2$	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}}$	$(4e)$	H_2O
$\mathbf{B}_2$	$\frac{1}{2}\,\mathbf{a}_1$	$\frac{1}{4}a\,\hat{\mathbf{x}} - \frac{1}{4}b\,\hat{\mathbf{y}}$	$(4e)$	H_2O
$\mathbf{B}_3$	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2$	$y_2 b\,\hat{\mathbf{y}}$	$(4g)$	Mg I
$\mathbf{B}_4$	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2$	$-y_2 b\,\hat{\mathbf{y}}$	$(4g)$	Mg I
$\mathbf{B}_5$	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Mg II
$\mathbf{B}_6$	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Mg II
$\mathbf{B}_7$	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	O I
$\mathbf{B}_8$	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	O I
$\mathbf{B}_9$	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	OH I
$\mathbf{B}_{10}$	$-x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	OH I
$\mathbf{B}_{11}$	$x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	OH II
$\mathbf{B}_{12}$	$-x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	OH II
$\mathbf{B}_{13}$	$(x_7 - y_7)\,\mathbf{a}_1 + (x_7 + y_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Mg III
$\mathbf{B}_{14}$	$(-x_7 - y_7)\,\mathbf{a}_1 + (-x_7 + y_7)\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	$(-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Mg III
$\mathbf{B}_{15}$	$(-x_7 + y_7)\,\mathbf{a}_1 + (-x_7 - y_7)\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	$(-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Mg III
$\mathbf{B}_{16}$	$(x_7 + y_7)\,\mathbf{a}_1 + (x_7 - y_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Mg III
$\mathbf{B}_{17}$	$(x_8 - y_8)\,\mathbf{a}_1 + (x_8 + y_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	$(x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Mg IV
$\mathbf{B}_{18}$	$(-x_8 - y_8)\,\mathbf{a}_1 + (-x_8 + y_8)\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	$(-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Mg IV
$\mathbf{B}_{19}$	$(-x_8 + y_8)\,\mathbf{a}_1 + (-x_8 - y_8)\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	$(-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Mg IV
$\mathbf{B}_{20}$	$(x_8 + y_8)\,\mathbf{a}_1 + (x_8 - y_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	$(x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Mg IV
$\mathbf{B}_{21}$	$(x_9 - y_9)\,\mathbf{a}_1 + (x_9 + y_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	$(x_9 a + z_9 c\cos\beta)\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O II
$\mathbf{B}_{22}$	$(-x_9 - y_9)\,\mathbf{a}_1 + (-x_9 + y_9)\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	$(-x_9 a - z_9 c\cos\beta)\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} - z_9 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O II

$$
\begin{array}{lcll}
\mathbf{B}_{23} & = & (-x_9 + y_9)\,\mathbf{a}_1 + (-x_9 - y_9)\,\mathbf{a}_2 - z_9\,\mathbf{a}_3 & = (-x_9 a - z_9 c\cos\beta)\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} - z_9 c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{O II} \\
\mathbf{B}_{24} & = & (x_9 + y_9)\,\mathbf{a}_1 + (x_9 - y_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 & = (x_9 a + z_9 c\cos\beta)\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + z_9 c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{O II} \\
\mathbf{B}_{25} & = & (x_{10} - y_{10})\,\mathbf{a}_1 + (x_{10} + y_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 & = (x_{10} a + z_{10} c\cos\beta)\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{O III} \\
\mathbf{B}_{26} & = & (-x_{10} - y_{10})\,\mathbf{a}_1 + (-x_{10} + y_{10})\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 & = (-x_{10} a - z_{10} c\cos\beta)\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} - z_{10} c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{O III} \\
\mathbf{B}_{27} & = & (-x_{10} + y_{10})\,\mathbf{a}_1 + (-x_{10} - y_{10})\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 & = (-x_{10} a - z_{10} c\cos\beta)\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} - z_{10} c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{O III} \\
\mathbf{B}_{28} & = & (x_{10} + y_{10})\,\mathbf{a}_1 + (x_{10} - y_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 & = (x_{10} a + z_{10} c\cos\beta)\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} + z_{10} c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{O III} \\
\mathbf{B}_{29} & = & (x_{11} - y_{11})\,\mathbf{a}_1 + (x_{11} + y_{11})\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 & = (x_{11} a + z_{11} c\cos\beta)\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + z_{11} c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{O IV} \\
\mathbf{B}_{30} & = & (-x_{11} - y_{11})\,\mathbf{a}_1 + (-x_{11} + y_{11})\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3 & = (-x_{11} a - z_{11} c\cos\beta)\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} - z_{11} c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{O IV} \\
\mathbf{B}_{31} & = & (-x_{11} + y_{11})\,\mathbf{a}_1 + (-x_{11} - y_{11})\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3 & = (-x_{11} a - z_{11} c\cos\beta)\,\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} - z_{11} c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{O IV} \\
\mathbf{B}_{32} & = & (x_{11} + y_{11})\,\mathbf{a}_1 + (x_{11} - y_{11})\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 & = (x_{11} a + z_{11} c\cos\beta)\,\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} + z_{11} c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{O IV} \\
\mathbf{B}_{33} & = & (x_{12} - y_{12})\,\mathbf{a}_1 + (x_{12} + y_{12})\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 & = (x_{12} a + z_{12} c\cos\beta)\,\hat{\mathbf{x}} + y_{12} b\,\hat{\mathbf{y}} + z_{12} c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{O V} \\
\mathbf{B}_{34} & = & (-x_{12} - y_{12})\,\mathbf{a}_1 + (-x_{12} + y_{12})\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3 & = (-x_{12} a - z_{12} c\cos\beta)\,\hat{\mathbf{x}} + y_{12} b\,\hat{\mathbf{y}} - z_{12} c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{O V} \\
\mathbf{B}_{35} & = & (-x_{12} + y_{12})\,\mathbf{a}_1 + (-x_{12} - y_{12})\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3 & = (-x_{12} a - z_{12} c\cos\beta)\,\hat{\mathbf{x}} - y_{12} b\,\hat{\mathbf{y}} - z_{12} c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{O V} \\
\mathbf{B}_{36} & = & (x_{12} + y_{12})\,\mathbf{a}_1 + (x_{12} - y_{12})\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 & = (x_{12} a + z_{12} c\cos\beta)\,\hat{\mathbf{x}} - y_{12} b\,\hat{\mathbf{y}} + z_{12} c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{O V} \\
\mathbf{B}_{37} & = & (x_{13} - y_{13})\,\mathbf{a}_1 + (x_{13} + y_{13})\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3 & = (x_{13} a + z_{13} c\cos\beta)\,\hat{\mathbf{x}} + y_{13} b\,\hat{\mathbf{y}} + z_{13} c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{O VI} \\
\mathbf{B}_{38} & = & (-x_{13} - y_{13})\,\mathbf{a}_1 + (-x_{13} + y_{13})\,\mathbf{a}_2 - z_{13}\,\mathbf{a}_3 & = (-x_{13} a - z_{13} c\cos\beta)\,\hat{\mathbf{x}} + y_{13} b\,\hat{\mathbf{y}} - z_{13} c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{O VI} \\
\mathbf{B}_{39} & = & (-x_{13} + y_{13})\,\mathbf{a}_1 + (-x_{13} - y_{13})\,\mathbf{a}_2 - z_{13}\,\mathbf{a}_3 & = (-x_{13} a - z_{13} c\cos\beta)\,\hat{\mathbf{x}} - y_{13} b\,\hat{\mathbf{y}} - z_{13} c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{O VI} \\
\mathbf{B}_{40} & = & (x_{13} + y_{13})\,\mathbf{a}_1 + (x_{13} - y_{13})\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3 & = (x_{13} a + z_{13} c\cos\beta)\,\hat{\mathbf{x}} - y_{13} b\,\hat{\mathbf{y}} + z_{13} c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{O VI} \\
\mathbf{B}_{41} & = & (x_{14} - y_{14})\,\mathbf{a}_1 + (x_{14} + y_{14})\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3 & = (x_{14} a + z_{14} c\cos\beta)\,\hat{\mathbf{x}} + y_{14} b\,\hat{\mathbf{y}} + z_{14} c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{OH III} \\
\mathbf{B}_{42} & = & (-x_{14} - y_{14})\,\mathbf{a}_1 + (-x_{14} + y_{14})\,\mathbf{a}_2 - z_{14}\,\mathbf{a}_3 & = (-x_{14} a - z_{14} c\cos\beta)\,\hat{\mathbf{x}} + y_{14} b\,\hat{\mathbf{y}} - z_{14} c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{OH III} \\
\mathbf{B}_{43} & = & (-x_{14} + y_{14})\,\mathbf{a}_1 + (-x_{14} - y_{14})\,\mathbf{a}_2 - z_{14}\,\mathbf{a}_3 & = (-x_{14} a - z_{14} c\cos\beta)\,\hat{\mathbf{x}} - y_{14} b\,\hat{\mathbf{y}} - z_{14} c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{OH III} \\
\mathbf{B}_{44} & = & (x_{14} + y_{14})\,\mathbf{a}_1 + (x_{14} - y_{14})\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3 & = (x_{14} a + z_{14} c\cos\beta)\,\hat{\mathbf{x}} - y_{14} b\,\hat{\mathbf{y}} + z_{14} c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{OH III} \\
\mathbf{B}_{45} & = & (x_{15} - y_{15})\,\mathbf{a}_1 + (x_{15} + y_{15})\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 & = (x_{15} a + z_{15} c\cos\beta)\,\hat{\mathbf{x}} + y_{15} b\,\hat{\mathbf{y}} + z_{15} c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{OH IV} \\
\mathbf{B}_{46} & = & (-x_{15} - y_{15})\,\mathbf{a}_1 + (-x_{15} + y_{15})\,\mathbf{a}_2 - z_{15}\,\mathbf{a}_3 & = (-x_{15} a - z_{15} c\cos\beta)\,\hat{\mathbf{x}} + y_{15} b\,\hat{\mathbf{y}} - z_{15} c\sin\beta\,\hat{\mathbf{z}} \quad (8j) \quad \text{OH IV}
\end{array}
$$

$\mathbf{B}_{47}$	=	$(-x_{15}+y_{15})\,\mathbf{a}_1+(-x_{15}-y_{15})\,\mathbf{a}_2-z_{15}\,\mathbf{a}_3$	=	$(-x_{15}a-z_{15}c\cos\beta)\,\hat{\mathbf{x}}-y_{15}b\,\hat{\mathbf{y}}-z_{15}c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	OH IV
$\mathbf{B}_{48}$	=	$(x_{15}+y_{15})\,\mathbf{a}_1+(x_{15}-y_{15})\,\mathbf{a}_2+z_{15}\,\mathbf{a}_3$	=	$(x_{15}a+z_{15}c\cos\beta)\,\hat{\mathbf{x}}-y_{15}b\,\hat{\mathbf{y}}+z_{15}c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	OH IV
$\mathbf{B}_{49}$	=	$(x_{16}-y_{16})\,\mathbf{a}_1+(x_{16}+y_{16})\,\mathbf{a}_2+z_{16}\,\mathbf{a}_3$	=	$(x_{16}a+z_{16}c\cos\beta)\,\hat{\mathbf{x}}+y_{16}b\,\hat{\mathbf{y}}+z_{16}c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	Si I
$\mathbf{B}_{50}$	=	$(-x_{16}-y_{16})\,\mathbf{a}_1+(-x_{16}+y_{16})\,\mathbf{a}_2-z_{16}\,\mathbf{a}_3$	=	$(-x_{16}a-z_{16}c\cos\beta)\,\hat{\mathbf{x}}+y_{16}b\,\hat{\mathbf{y}}-z_{16}c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	Si I
$\mathbf{B}_{51}$	=	$(-x_{16}+y_{16})\,\mathbf{a}_1+(-x_{16}-y_{16})\,\mathbf{a}_2-z_{16}\,\mathbf{a}_3$	=	$(-x_{16}a-z_{16}c\cos\beta)\,\hat{\mathbf{x}}-y_{16}b\,\hat{\mathbf{y}}-z_{16}c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	Si I
$\mathbf{B}_{52}$	=	$(x_{16}+y_{16})\,\mathbf{a}_1+(x_{16}-y_{16})\,\mathbf{a}_2+z_{16}\,\mathbf{a}_3$	=	$(x_{16}a+z_{16}c\cos\beta)\,\hat{\mathbf{x}}-y_{16}b\,\hat{\mathbf{y}}+z_{16}c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	Si I
$\mathbf{B}_{53}$	=	$(x_{17}-y_{17})\,\mathbf{a}_1+(x_{17}+y_{17})\,\mathbf{a}_2+z_{17}\,\mathbf{a}_3$	=	$(x_{17}a+z_{17}c\cos\beta)\,\hat{\mathbf{x}}+y_{17}b\,\hat{\mathbf{y}}+z_{17}c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	Si II
$\mathbf{B}_{54}$	=	$(-x_{17}-y_{17})\,\mathbf{a}_1+(-x_{17}+y_{17})\,\mathbf{a}_2-z_{17}\,\mathbf{a}_3$	=	$(-x_{17}a-z_{17}c\cos\beta)\,\hat{\mathbf{x}}+y_{17}b\,\hat{\mathbf{y}}-z_{17}c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	Si II
$\mathbf{B}_{55}$	=	$(-x_{17}+y_{17})\,\mathbf{a}_1+(-x_{17}-y_{17})\,\mathbf{a}_2-z_{17}\,\mathbf{a}_3$	=	$(-x_{17}a-z_{17}c\cos\beta)\,\hat{\mathbf{x}}-y_{17}b\,\hat{\mathbf{y}}-z_{17}c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	Si II
$\mathbf{B}_{56}$	=	$(x_{17}+y_{17})\,\mathbf{a}_1+(x_{17}-y_{17})\,\mathbf{a}_2+z_{17}\,\mathbf{a}_3$	=	$(x_{17}a+z_{17}c\cos\beta)\,\hat{\mathbf{x}}-y_{17}b\,\hat{\mathbf{y}}+z_{17}c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	Si II

References:

- B. E. Warren and W. L. Bragg, *The Structure of Chrysotile $H_4Mg_3Si_2O_9$*, Zeitschrift für Kristallographie - Crystalline Materials **76**, 201–210 (1931), doi:10.1524/zkri.1931.76.1.201.
- K. Yada, *Study of Chrysotile Asbestos by a High Resolution Electron Microscope*, Acta Cryst. **23**, 704–707 (1967), doi:10.1107/S0365110X67003524.

Found in:

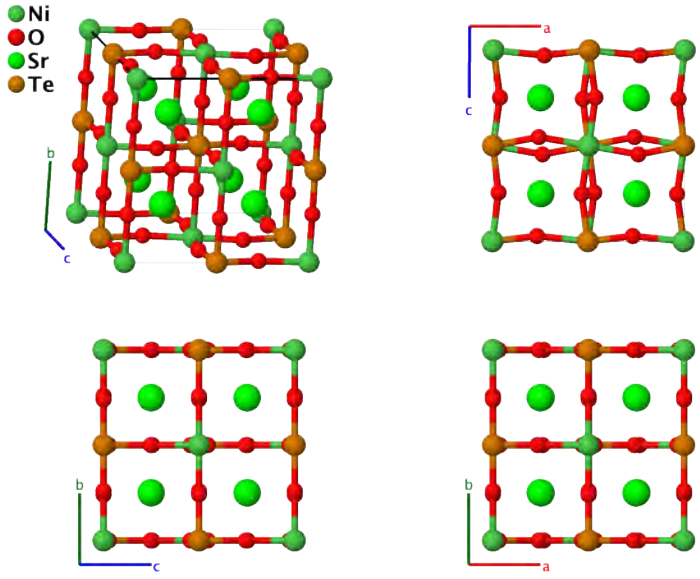
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1531
- POSCAR: pp. 1532

Sr_2NiTeO_6 Structure: AB6C2D_mC40_12_ad_gh4i_j_bc

http://aflow.org/prototype-encyclopedia/AB6C2D_mC40_12_ad_gh4i_j_bc



Prototype	:	NiO_6Sr_2Te
AFLOW prototype label	:	AB6C2D_mC40_12_ad_gh4i_j_bc
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC40
Space group number	:	12
Space group symbol	:	$C2/m$
AFLOW prototype command	:	`aflow --proto=AB6C2D_mC40_12_ad_gh4i_j_bc --params=`$a, b/a, c/a, \beta, y_5, y_6, x_7, z_7, x_8, z_8, x_9, z_9, x_{10}, z_{10}, x_{11}, y_{11}, z_{11}$

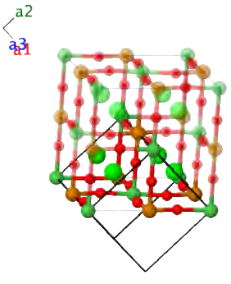
Other compounds with this structure

- Sr_2NiTeO_6 and Cs_2RbDy_6

- At high temperatures Sr_2NiTeO_6 transforms into the cubic perovskite $E2_1$ structure.

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(2*a*)	Ni I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	(2*b*)	Te I
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	(2*c*)	Te II
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(a + c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	(2*d*)	Ni II
$\mathbf{B}_5$	=	$-y_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2$	=	$y_5 b\,\hat{\mathbf{y}}$	(4*g*)	O I
$\mathbf{B}_6$	=	$y_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2$	=	$-y_5 b\,\hat{\mathbf{y}}$	(4*g*)	O I
$\mathbf{B}_7$	=	$-y_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	(4*h*)	O II
$\mathbf{B}_8$	=	$y_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	(4*h*)	O II
$\mathbf{B}_9$	=	$x_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	O III
$\mathbf{B}_{10}$	=	$-x_7\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} - z_7 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	O III
$\mathbf{B}_{11}$	=	$x_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + z_8 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	O IV
$\mathbf{B}_{12}$	=	$-x_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} - z_8 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	O IV
$\mathbf{B}_{13}$	=	$x_9\,\mathbf{a}_1 + x_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$(x_9 a + z_9 c\cos\beta)\,\hat{\mathbf{x}} + z_9 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	O V
$\mathbf{B}_{14}$	=	$-x_9\,\mathbf{a}_1 - x_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$(-x_9 a - z_9 c\cos\beta)\,\hat{\mathbf{x}} - z_9 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	O V
$\mathbf{B}_{15}$	=	$x_{10}\,\mathbf{a}_1 + x_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$(x_{10} a + z_{10} c\cos\beta)\,\hat{\mathbf{x}} + z_{10} c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	O VI
$\mathbf{B}_{16}$	=	$-x_{10}\,\mathbf{a}_1 - x_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3$	=	$(-x_{10} a - z_{10} c\cos\beta)\,\hat{\mathbf{x}} - z_{10} c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	O VI
$\mathbf{B}_{17}$	=	$(x_{11} - y_{11})\,\mathbf{a}_1 + (x_{11} + y_{11})\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$(x_{11} a + z_{11} c\cos\beta)\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + z_{11} c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	Sr
$\mathbf{B}_{18}$	=	$(-x_{11} - y_{11})\,\mathbf{a}_1 + (-x_{11} + y_{11})\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3$	=	$(-x_{11} a - z_{11} c\cos\beta)\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} - z_{11} c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	Sr
$\mathbf{B}_{19}$	=	$(-x_{11} + y_{11})\,\mathbf{a}_1 + (-x_{11} - y_{11})\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3$	=	$(-x_{11} a - z_{11} c\cos\beta)\,\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} - z_{11} c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	Sr
$\mathbf{B}_{20}$	=	$(x_{11} + y_{11})\,\mathbf{a}_1 + (x_{11} - y_{11})\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$(x_{11} a + z_{11} c\cos\beta)\,\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} + z_{11} c\sin\beta\,\hat{\mathbf{z}}$	(8*j*)	Sr

References:

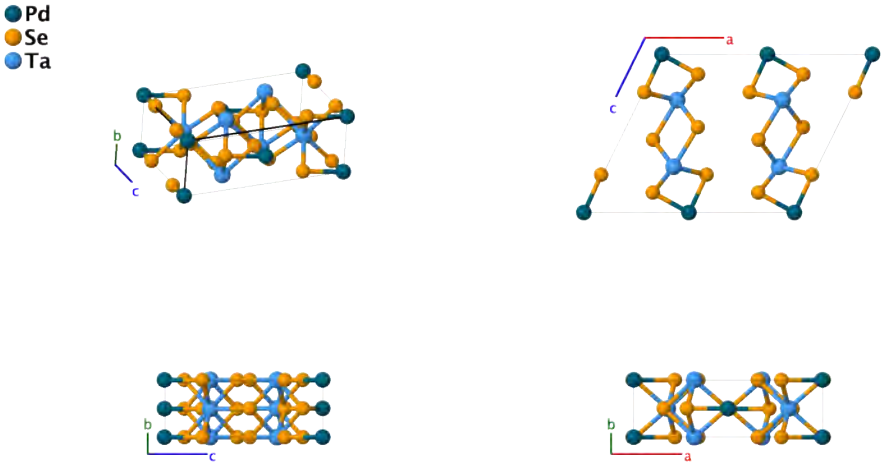
- D. Iwanaga, Y. Inaguma, and M. Itoh, *Structure and Magnetic Properties of Sr_2NiAO_6 (A = W, Te)*, Mater. Res. Bull. **35**, 449–457 (2000), doi:10.1016/S0025-5408(00)00222-1.

Geometry files:

- CIF: pp. 1532
- POSCAR: pp. 1532

Ta_2PdSe_6 Structure: AB6C2_mC18_12_a_3i_i

http://aflow.org/prototype-encyclopedia/AB6C2_mC18_12_a_3i_i



Prototype	:	$PdSe_6Ta_2$
AFLOW prototype label	:	AB6C2_mC18_12_a_3i_i
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC18
Space group number	:	12
Space group symbol	:	$C2/m$
AFLOW prototype command	:	`aflow --proto=AB6C2_mC18_12_a_3i_i --params=`$a, b/a, c/a, \beta, x_2, z_2, x_3, z_3, x_4, z_4, x_5, z_5$

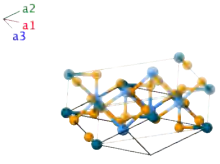
Other compounds with this structure

- Nb_2PdS_6, Nb_2PdSe_6, and Ta_2PdS_6

- (Keszler, 1985) gave the structure in the $I2/m$ setting of space group #12. We used FINDSYM to change this to the standard $C2/m$ setting. Because of this change, our primitive vectors are linear combinations of the original ones, and the lattice has been rotated.

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$ =	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	Pd

$\mathbf{B}_2$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Se I
$\mathbf{B}_3$	=	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} - z_2 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Se I
$\mathbf{B}_4$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Se II
$\mathbf{B}_5$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Se II
$\mathbf{B}_6$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Se III
$\mathbf{B}_7$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Se III
$\mathbf{B}_8$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Ta
$\mathbf{B}_9$	=	$-x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Ta

References:

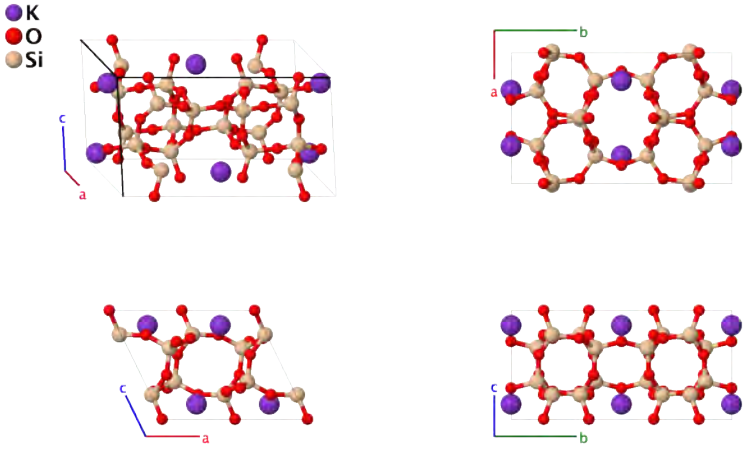
- D. A. Keszler, P. J. Squattrito, N. E. Brese, J. A. Ibers, M. Shang, and J. Lu, *New layered ternary chalcogenides: tantalum palladium sulfide (Ta_2PdS_6), tantalum palladium selenide (Ta_2PdSe_6), niobium palladium sulfide (Nb_2PdS_6), niobium palladium selenide (Nb_2PdSe_6)*, Inorg. Chem. pp. 3063–3067 (1985), doi:10.1021/ic00213a038.

Geometry files:

- CIF: pp. 1533
- POSCAR: pp. 1533

Sanidine ($KAlSi_3O_8$, $S6_7$) Structure: AB8C4_mC52_12_i_gi3j_2j

http://aflow.org/prototype-encyclopedia/AB8C4_mC52_12_i_gi3j_2j



Prototype	:	$AlKO_8Si_3$
AFLOW prototype label	:	AB8C4_mC52_12_i_gi3j_2j
***Strukturbericht* designation**	:	$S6_7$
Pearson symbol	:	mC52
Space group number	:	12
Space group symbol	:	$C2/m$
AFLOW prototype command	:	`aflow --proto=AB8C4_mC52_12_i_gi3j_2j --params=`$a, b/a, c/a, \beta, y_1, x_2, z_2, x_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8$

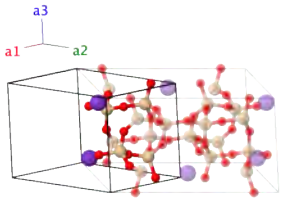
Other compounds with this structure

- $NaAlSi_3O_8$ and $(BaKNa)AlSi_3O_8$ (orthoclase)

- Begining with the $S6_7$ defining paper of (Taylor, 1933), all researchers agree that the sites Si-I and Si-II are actually a random mixture of silicon and aluminum with overall stoichiometry $AlSi_3$. Most results found in (Downs, 2003) assume that this composition holds independently for both sites, but (Scambos, 1987) states that the Si-I site is $Al_{1.064}Si_{2.936}$ while the Si-II site is $Al_{0.936}Si_{3.064}$.
- (Scambos, 1987) list the coordinate $y_8 = 0.18813$ for the Si-II site in Table 2. This position does not result in the expected SiO_4 tetrahedra, and the interatomic distances and angles found do not agree with their Table 5. (Downs, 2003) correct this to $y_8 = 0.11813$, and this value results in good tetrahedra and reproduces the distances and angles reported by (Scambos, 1987). We use this later value on this page.

Base-centered Monoclinic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2}\,b\,\hat{\mathbf{y}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,b\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2$	=	$y_1 b\,\hat{\mathbf{y}}$	$(4g)$	O I
$\mathbf{B}_2$	=	$y_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2$	=	$-y_1 b\,\hat{\mathbf{y}}$	$(4g)$	O I
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	K
$\mathbf{B}_4$	=	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} - z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	K
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	O II
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	O II
$\mathbf{B}_7$	=	$(x_4 - y_4)\,\mathbf{a}_1 + (x_4 + y_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O III
$\mathbf{B}_8$	=	$(-x_4 - y_4)\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O III
$\mathbf{B}_9$	=	$(-x_4 + y_4)\,\mathbf{a}_1 + (-x_4 - y_4)\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O III
$\mathbf{B}_{10}$	=	$(x_4 + y_4)\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O III
$\mathbf{B}_{11}$	=	$(x_5 - y_5)\,\mathbf{a}_1 + (x_5 + y_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O IV
$\mathbf{B}_{12}$	=	$(-x_5 - y_5)\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O IV
$\mathbf{B}_{13}$	=	$(-x_5 + y_5)\,\mathbf{a}_1 + (-x_5 - y_5)\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O IV
$\mathbf{B}_{14}$	=	$(x_5 + y_5)\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O IV
$\mathbf{B}_{15}$	=	$(x_6 - y_6)\,\mathbf{a}_1 + (x_6 + y_6)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O V
$\mathbf{B}_{16}$	=	$(-x_6 - y_6)\,\mathbf{a}_1 + (-x_6 + y_6)\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O V
$\mathbf{B}_{17}$	=	$(-x_6 + y_6)\,\mathbf{a}_1 + (-x_6 - y_6)\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O V
$\mathbf{B}_{18}$	=	$(x_6 + y_6)\,\mathbf{a}_1 + (x_6 - y_6)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	O V
$\mathbf{B}_{19}$	=	$(x_7 - y_7)\,\mathbf{a}_1 + (x_7 + y_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Si I
$\mathbf{B}_{20}$	=	$(-x_7 - y_7)\,\mathbf{a}_1 + (-x_7 + y_7)\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Si I
$\mathbf{B}_{21}$	=	$(-x_7 + y_7)\,\mathbf{a}_1 + (-x_7 - y_7)\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Si I

$\mathbf{B}_{22}$	=	$(x_7 + y_7)\,\mathbf{a}_1 + (x_7 - y_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Si I
$\mathbf{B}_{23}$	=	$(x_8 - y_8)\,\mathbf{a}_1 + (x_8 + y_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Si II
$\mathbf{B}_{24}$	=	$(-x_8 - y_8)\,\mathbf{a}_1 + (-x_8 + y_8)\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Si II
$\mathbf{B}_{25}$	=	$(-x_8 + y_8)\,\mathbf{a}_1 + (-x_8 - y_8)\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Si II
$\mathbf{B}_{26}$	=	$(x_8 + y_8)\,\mathbf{a}_1 + (x_8 - y_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	Si II

References:

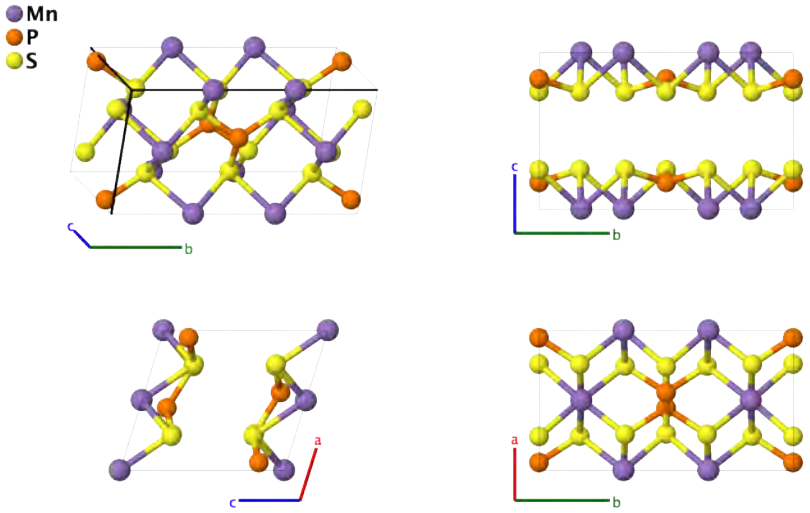
- T. A. Scambos, J. R. Smyth, and T. C. McCormick, *Crystal-structure refinement of high sanidine from the upper mantle*, Am. Mineral. **72**, 973–978 (1987).
- W. H. Taylor, *The Structure of Sanidine and Other Felspars*, Zeitschrift für Kristallographie - Crystalline Materials **85**, 425–442 (1933), doi:10.1524/zkri.1933.85.1.425.
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1533
- POSCAR: pp. 1533

$MnPS_3$ Structure: ABC3_mC20_12_g_i_ij

http://aflow.org/prototype-encyclopedia/ABC3_mC20_12_g_i_ij



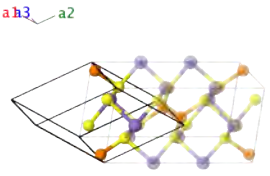
Prototype	:	$MnPS_3$
AFLOW prototype label	:	ABC3_mC20_12_g_i_ij
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC20
Space group number	:	12
Space group symbol	:	$C2/m$
AFLOW prototype command	:	`aflow --proto=ABC3_mC20_12_g_i_ij --params=`$a, b/a, c/a, \beta, y_1, x_2, z_2, x_3, z_3, x_4, y_4, z_4$

Other compounds with this structure

- $CdPS_3$, $CdPSe_3$, $CoPS_3$, $CoPSe_3$, $FePS_3$, $FePSe_3$, $MnPSe_3$, $NiPS_3$, and $NiPSe_3$

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2$	=	$y_1 b\,\hat{\mathbf{y}}$	$(4g)$	Mn
$\mathbf{B}_2$	=	$y_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2$	=	$-y_1 b\,\hat{\mathbf{y}}$	$(4g)$	Mn
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	P
$\mathbf{B}_4$	=	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} - z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	P

$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	S I
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	S I
$\mathbf{B}_7$	=	$(x_4 - y_4)\,\mathbf{a}_1 + (x_4 + y_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	S II
$\mathbf{B}_8$	=	$(-x_4 - y_4)\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	S II
$\mathbf{B}_9$	=	$(-x_4 + y_4)\,\mathbf{a}_1 + (-x_4 - y_4)\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	S II
$\mathbf{B}_{10}$	=	$(x_4 + y_4)\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(8j)$	S II

References:

- G. Ouvrard, R. Brec, and J. Rouxel, *Structural determination of some MPS_3 layered phases (M = Mn, Fe, Co, Ni and Cd)*, Mater. Res. Bull. **20**, 1181–1189 (1985), doi:10.1016/0025-5408(85)90092-3.

Found in:

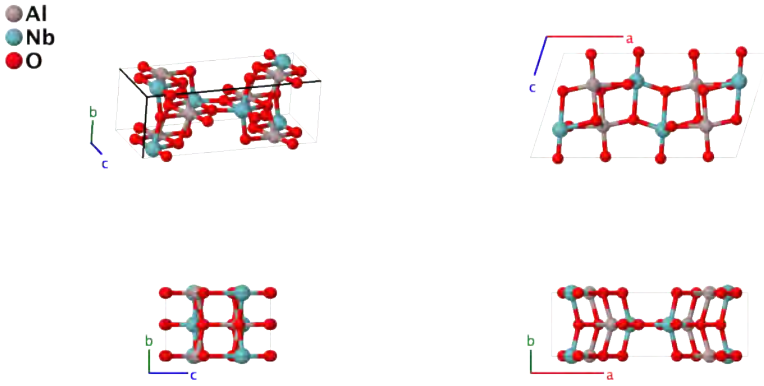
- V. Zhukov, S. Alvarez, and D. Novikov, *Electronic band structure of the magnetic layered semiconductors MPS_3 (M = Mn, Fe and Ni)*, J. Phys. Chem. Solids **57**, 647–652 (1996), doi:10.1016/0022-3697(95)00203-0.

Geometry files:

- CIF: pp. 1534
- POSCAR: pp. 1534

$AlNbO_4$ Structure: ABC4_mC24_12_i_i_4i

http://aflow.org/prototype-encyclopedia/ABC4_mC24_12_i_i_4i

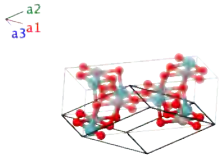


Prototype	:	$AlNbO_4$
AFLOW prototype label	:	ABC4_mC24_12_i_i_4i
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC24
Space group number	:	12
Space group symbol	:	$C2/m$
AFLOW prototype command	:	`aflow --proto=ABC4_mC24_12_i_i_4i --params=`$a, b/a, c/a, \beta, x_1, z_1, x_2, z_2, x_3, z_3, x_4, z_4, x_5, z_5, x_6, z_6$

- (Ardit, 2012) looked at samples with obvious disordering on the sites we label Al and Nb. For their sample N00, the occupation of our Al site was 80% aluminum and 20% niobium, with the reverse concentrations on our Nb site.
- (Pederson, 1962) analyzed a sample of $AlNbO_4$ assuming that these sites were fully ordered, however the atomic positions listed there have some errors, so we used the data found by Ardit *et al.*
- The atomic positions for this structure are not found in the main text. They are filed as item #AM-12-035 in the American Mineralogist repository.

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + x_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Al
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 - x_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$(-x_1 a - z_1 c\cos\beta)\,\hat{\mathbf{x}} - z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Al
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	Nb

$\mathbf{B}_4$	=	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2a - z_2c\cos\beta)\,\hat{\mathbf{x}} - z_2c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Nb
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3a + z_3c\cos\beta)\,\hat{\mathbf{x}} + z_3c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	O I
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3a - z_3c\cos\beta)\,\hat{\mathbf{x}} - z_3c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	O I
$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4a + z_4c\cos\beta)\,\hat{\mathbf{x}} + z_4c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	O II
$\mathbf{B}_8$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4a - z_4c\cos\beta)\,\hat{\mathbf{x}} - z_4c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	O II
$\mathbf{B}_9$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5a + z_5c\cos\beta)\,\hat{\mathbf{x}} + z_5c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	O III
$\mathbf{B}_{10}$	=	$-x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5a - z_5c\cos\beta)\,\hat{\mathbf{x}} - z_5c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	O III
$\mathbf{B}_{11}$	=	$x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6a + z_6c\cos\beta)\,\hat{\mathbf{x}} + z_6c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	O IV
$\mathbf{B}_{12}$	=	$-x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6a - z_6c\cos\beta)\,\hat{\mathbf{x}} - z_6c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	O IV

References:

- M. Ardit, M. Dondi, and G. Cruciani, *Structural stability, cation ordering, and local relaxation along the $AlNbO_4$-$Al_{0.5}Cr_{0.5}NbO_4$ join*, Am. Mineral. **97**, 910–917 (2012), doi:10.2138/am.2012.3977. Structural data from http://www.minsocam.org/MSA/AmMin/TOC/2012/MJ12_Data/Ardit_p910_12.zip.
- B. F. Pedersen, *The Crystal Structure of Aluminium Niobium Oxide ($AlNbO_4$)*, Acta Chem. Scand. **16**, 421–430 (1962), doi:10.3891/acta.chem.scand.16-0421.

Found in:

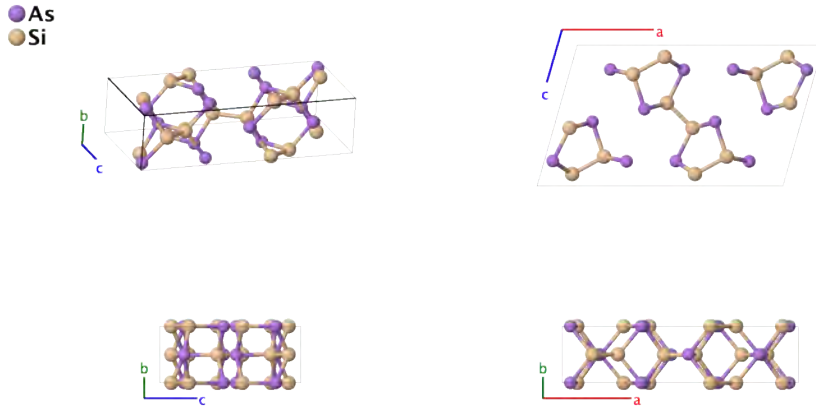
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1534
- POSCAR: pp. 1535

SiAs Structure: AB_mC24_12_3i_3i

http://aflow.org/prototype-encyclopedia/AB_mC24_12_3i_3i



Prototype	:	AsSi
AFLOW prototype label	:	AB_mC24_12_3i_3i
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC24
Space group number	:	12
Space group symbol	:	$C2/m$
AFLOW prototype command	:	`aflow --proto=AB_mC24_12_3i_3i --params=`$a, b/a, c/a, \beta, x_1, z_1, x_2, z_2, x_3, z_3, x_4, z_4, x_5, z_5, x_6, z_6$

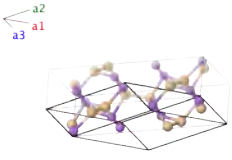
Other compounds with this structure

- GaTe and GeAs

- The structures of GaTe and GeAs were apparently determined by (Bryden, 1965) before (Wadsten, 1965) found the structure of SiAs, but they were never published (Pearson, 1964; Wadsten, 1965; Mead, 1982). As SiAs was the first published determination of this structure, we use it as the prototype.

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + x_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	As I
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 - x_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$(-x_1 a - z_1 c\cos\beta)\,\hat{\mathbf{x}} - z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	As I
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(4i)$	As II

$\mathbf{B}_4$	=	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2a - z_2c\cos\beta)\,\hat{\mathbf{x}} - z_2c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	As II
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3a + z_3c\cos\beta)\,\hat{\mathbf{x}} + z_3c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	As III
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3a - z_3c\cos\beta)\,\hat{\mathbf{x}} - z_3c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	As III
$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4a + z_4c\cos\beta)\,\hat{\mathbf{x}} + z_4c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Si I
$\mathbf{B}_8$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4a - z_4c\cos\beta)\,\hat{\mathbf{x}} - z_4c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Si I
$\mathbf{B}_9$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5a + z_5c\cos\beta)\,\hat{\mathbf{x}} + z_5c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Si II
$\mathbf{B}_{10}$	=	$-x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5a - z_5c\cos\beta)\,\hat{\mathbf{x}} - z_5c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Si II
$\mathbf{B}_{11}$	=	$x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6a + z_6c\cos\beta)\,\hat{\mathbf{x}} + z_6c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Si III
$\mathbf{B}_{12}$	=	$-x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6a - z_6c\cos\beta)\,\hat{\mathbf{x}} - z_6c\sin\beta\,\hat{\mathbf{z}}$	(4*i*)	Si III

References:

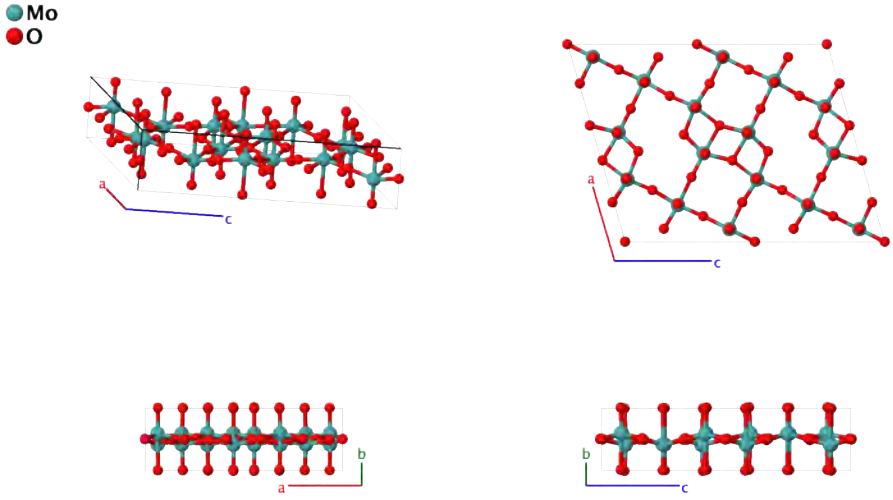
- T. Wadsten, *The Crystal Structure of SiAs*, Acta Chem. Scand. **19**, 1232–1238 (1965), doi:10.3891/acta.chem.scand.19-1232.
- W. B. Pearson, *The crystal structures of semiconductors and a general valence rule*, Acta Cryst. **17**, 1–15 (1964), doi:10.1107/S0365110X64000019.
- D. G. Mead, *Long wavelength study of semiconducting germanium arsenide, GeAs*, Infrared Phys. **22**, 209–213 (1982), doi:10.1016/0020-0891(82)90045-8.
- J. H. Bryden, *Private Communication* (1965). To T. Wadsten.

Geometry files:

- CIF: pp. 1535
- POSCAR: pp. 1535

High-Temperature Mo_8O_{23} Structure: A8B23_mP62_13_4g_c11g

http://aflow.org/prototype-encyclopedia/A8B23_mP62_13_4g_c11g

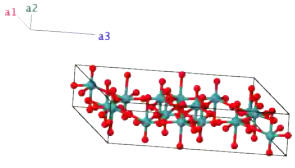


Prototype	:	Mo_8O_{23}
AFLOW prototype label	:	A8B23_mP62_13_4g_c11g
***Strukturbericht* designation**	:	None
Pearson symbol	:	mP62
Space group number	:	13
Space group symbol	:	$P2/c$
AFLOW prototype command	:	`aflow --proto=A8B23_mP62_13_4g_c11g --params=`$a, b/a, c/a, \beta, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}$

- Above 285 K, the structure exhibits an incommensurate charge density wave. This data was taken at 370 K, so the structure given here is only approximate. Below 285 K the CDW is locked in, leading to the low-temperature Mo_8O_{23} structure.

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$\mathbf{B}_1$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}b\,\hat{\mathbf{y}}$	$(2c)$	O I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	$(2c)$	O I
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2a + z_2c\cos\beta)\,\hat{\mathbf{x}} + y_2b\,\hat{\mathbf{y}} + z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4g)$	Mo I
$\mathbf{B}_4$	=	$-x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_2a - z_2c\cos\beta\right)\hat{\mathbf{x}} + y_2b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4g)$	Mo I
$\mathbf{B}_5$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2a - z_2c\cos\beta)\,\hat{\mathbf{x}} - y_2b\,\hat{\mathbf{y}} - z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4g)$	Mo I
$\mathbf{B}_6$	=	$x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_2a + z_2c\cos\beta\right)\hat{\mathbf{x}} - y_2b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4g)$	Mo I
$\mathbf{B}_7$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3a + z_3c\cos\beta)\,\hat{\mathbf{x}} + y_3b\,\hat{\mathbf{y}} + z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4g)$	Mo II
$\mathbf{B}_8$	=	$-x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_3a - z_3c\cos\beta\right)\hat{\mathbf{x}} + y_3b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4g)$	Mo II
$\mathbf{B}_9$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3a - z_3c\cos\beta)\,\hat{\mathbf{x}} - y_3b\,\hat{\mathbf{y}} - z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4g)$	Mo II
$\mathbf{B}_{10}$	=	$x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_3a + z_3c\cos\beta\right)\hat{\mathbf{x}} - y_3b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4g)$	Mo II
$\mathbf{B}_{11}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4a + z_4c\cos\beta)\,\hat{\mathbf{x}} + y_4b\,\hat{\mathbf{y}} + z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4g)$	Mo III
$\mathbf{B}_{12}$	=	$-x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_4a - z_4c\cos\beta\right)\hat{\mathbf{x}} + y_4b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4g)$	Mo III
$\mathbf{B}_{13}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4a - z_4c\cos\beta)\,\hat{\mathbf{x}} - y_4b\,\hat{\mathbf{y}} - z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4g)$	Mo III
$\mathbf{B}_{14}$	=	$x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_4a + z_4c\cos\beta\right)\hat{\mathbf{x}} - y_4b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4g)$	Mo III
$\mathbf{B}_{15}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5a + z_5c\cos\beta)\,\hat{\mathbf{x}} + y_5b\,\hat{\mathbf{y}} + z_5c\sin\beta\,\hat{\mathbf{z}}$	$(4g)$	Mo IV
$\mathbf{B}_{16}$	=	$-x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_5a - z_5c\cos\beta\right)\hat{\mathbf{x}} + y_5b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4g)$	Mo IV
$\mathbf{B}_{17}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5a - z_5c\cos\beta)\,\hat{\mathbf{x}} - y_5b\,\hat{\mathbf{y}} - z_5c\sin\beta\,\hat{\mathbf{z}}$	$(4g)$	Mo IV
$\mathbf{B}_{18}$	=	$x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_5a + z_5c\cos\beta\right)\hat{\mathbf{x}} - y_5b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4g)$	Mo IV
$\mathbf{B}_{19}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6a + z_6c\cos\beta)\,\hat{\mathbf{x}} + y_6b\,\hat{\mathbf{y}} + z_6c\sin\beta\,\hat{\mathbf{z}}$	$(4g)$	O II
$\mathbf{B}_{20}$	=	$-x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_6a - z_6c\cos\beta\right)\hat{\mathbf{x}} + y_6b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4g)$	O II
$\mathbf{B}_{21}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6a - z_6c\cos\beta)\,\hat{\mathbf{x}} - y_6b\,\hat{\mathbf{y}} - z_6c\sin\beta\,\hat{\mathbf{z}}$	$(4g)$	O II
$\mathbf{B}_{22}$	=	$x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_6a + z_6c\cos\beta\right)\hat{\mathbf{x}} - y_6b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4g)$	O II
$\mathbf{B}_{23}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7a + z_7c\cos\beta)\,\hat{\mathbf{x}} + y_7b\,\hat{\mathbf{y}} + z_7c\sin\beta\,\hat{\mathbf{z}}$	$(4g)$	O III
$\mathbf{B}_{24}$	=	$-x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_7a - z_7c\cos\beta\right)\hat{\mathbf{x}} + y_7b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4g)$	O III
$\mathbf{B}_{25}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7a - z_7c\cos\beta)\,\hat{\mathbf{x}} - y_7b\,\hat{\mathbf{y}} - z_7c\sin\beta\,\hat{\mathbf{z}}$	$(4g)$	O III
$\mathbf{B}_{26}$	=	$x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_7a + z_7c\cos\beta\right)\hat{\mathbf{x}} - y_7b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4g)$	O III
$\mathbf{B}_{27}$	=	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8a + z_8c\cos\beta)\,\hat{\mathbf{x}} + y_8b\,\hat{\mathbf{y}} + z_8c\sin\beta\,\hat{\mathbf{z}}$	$(4g)$	O IV

$$\mathbf{B}_{28} = -x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_8 a - z_8 c\cos\beta\right)\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_8\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4g) \quad \text{O IV}$$

$$\mathbf{B}_{29} = -x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3 = (-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}} \quad (4g) \quad \text{O IV}$$

$$\mathbf{B}_{30} = x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_8 a + z_8 c\cos\beta\right)\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_8\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4g) \quad \text{O IV}$$

$$\mathbf{B}_{31} = x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 = (x_9 a + z_9 c\cos\beta)\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\sin\beta\,\hat{\mathbf{z}} \quad (4g) \quad \text{O V}$$

$$\mathbf{B}_{32} = -x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_9 a - z_9 c\cos\beta\right)\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_9\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4g) \quad \text{O V}$$

$$\mathbf{B}_{33} = -x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3 = (-x_9 a - z_9 c\cos\beta)\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} - z_9 c\sin\beta\,\hat{\mathbf{z}} \quad (4g) \quad \text{O V}$$

$$\mathbf{B}_{34} = x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_9 a + z_9 c\cos\beta\right)\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_9\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4g) \quad \text{O V}$$

$$\mathbf{B}_{35} = x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 = (x_{10} a + z_{10} c\cos\beta)\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\sin\beta\,\hat{\mathbf{z}} \quad (4g) \quad \text{O VI}$$

$$\mathbf{B}_{36} = -x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{10} a - z_{10} c\cos\beta\right)\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{10}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4g) \quad \text{O VI}$$

$$\mathbf{B}_{37} = -x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 = (-x_{10} a - z_{10} c\cos\beta)\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} - z_{10} c\sin\beta\,\hat{\mathbf{z}} \quad (4g) \quad \text{O VI}$$

$$\mathbf{B}_{38} = x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{10} a + z_{10} c\cos\beta\right)\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{10}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4g) \quad \text{O VI}$$

$$\mathbf{B}_{39} = x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 = (x_{11} a + z_{11} c\cos\beta)\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + z_{11} c\sin\beta\,\hat{\mathbf{z}} \quad (4g) \quad \text{O VII}$$

$$\mathbf{B}_{40} = -x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{11} a - z_{11} c\cos\beta\right)\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{11}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4g) \quad \text{O VII}$$

$$\mathbf{B}_{41} = -x_{11}\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3 = (-x_{11} a - z_{11} c\cos\beta)\,\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} - z_{11} c\sin\beta\,\hat{\mathbf{z}} \quad (4g) \quad \text{O VII}$$

$$\mathbf{B}_{42} = x_{11}\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{11} a + z_{11} c\cos\beta\right)\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{11}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4g) \quad \text{O VII}$$

$$\mathbf{B}_{43} = x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 = (x_{12} a + z_{12} c\cos\beta)\,\hat{\mathbf{x}} + y_{12} b\,\hat{\mathbf{y}} + z_{12} c\sin\beta\,\hat{\mathbf{z}} \quad (4g) \quad \text{O VIII}$$

$$\mathbf{B}_{44} = -x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{12} a - z_{12} c\cos\beta\right)\hat{\mathbf{x}} + y_{12} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{12}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4g) \quad \text{O VIII}$$

$$\mathbf{B}_{45} = -x_{12}\,\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3 = (-x_{12} a - z_{12} c\cos\beta)\,\hat{\mathbf{x}} - y_{12} b\,\hat{\mathbf{y}} - z_{12} c\sin\beta\,\hat{\mathbf{z}} \quad (4g) \quad \text{O VIII}$$

$$\mathbf{B}_{46} = x_{12}\,\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{12} a + z_{12} c\cos\beta\right)\hat{\mathbf{x}} - y_{12} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{12}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4g) \quad \text{O VIII}$$

$$\mathbf{B}_{47} = x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3 = (x_{13} a + z_{13} c\cos\beta)\,\hat{\mathbf{x}} + y_{13} b\,\hat{\mathbf{y}} + z_{13} c\sin\beta\,\hat{\mathbf{z}} \quad (4g) \quad \text{O IX}$$

$$\mathbf{B}_{48} = -x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{13} a - z_{13} c\cos\beta\right)\hat{\mathbf{x}} + y_{13} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{13}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4g) \quad \text{O IX}$$

$$\mathbf{B}_{49} = -x_{13}\,\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 - z_{13}\,\mathbf{a}_3 = (-x_{13} a - z_{13} c\cos\beta)\,\hat{\mathbf{x}} - y_{13} b\,\hat{\mathbf{y}} - z_{13} c\sin\beta\,\hat{\mathbf{z}} \quad (4g) \quad \text{O IX}$$

$$\mathbf{B}_{50} = x_{13}\,\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{13} a + z_{13} c\cos\beta\right)\hat{\mathbf{x}} - y_{13} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{13}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4g) \quad \text{O IX}$$

$$\mathbf{B}_{51} = x_{14}\,\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3 = (x_{14}a + z_{14}c\cos\beta)\,\hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} + z_{14}c\sin\beta\,\hat{\mathbf{z}} \quad (4g) \quad \text{O X}$$

$$\mathbf{B}_{52} = -x_{14}\,\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{14}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{14}a - z_{14}c\cos\beta\right)\hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{14}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4g) \quad \text{O X}$$

$$\mathbf{B}_{53} = -x_{14}\,\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 - z_{14}\,\mathbf{a}_3 = (-x_{14}a - z_{14}c\cos\beta)\,\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} - z_{14}c\sin\beta\,\hat{\mathbf{z}} \quad (4g) \quad \text{O X}$$

$$\mathbf{B}_{54} = x_{14}\,\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{14}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{14}a + z_{14}c\cos\beta\right)\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{14}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4g) \quad \text{O X}$$

$$\mathbf{B}_{55} = x_{15}\,\mathbf{a}_1 + y_{15}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 = (x_{15}a + z_{15}c\cos\beta)\,\hat{\mathbf{x}} + y_{15}b\,\hat{\mathbf{y}} + z_{15}c\sin\beta\,\hat{\mathbf{z}} \quad (4g) \quad \text{O XI}$$

$$\mathbf{B}_{56} = -x_{15}\,\mathbf{a}_1 + y_{15}\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{15}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{15}a - z_{15}c\cos\beta\right)\hat{\mathbf{x}} + y_{15}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{15}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4g) \quad \text{O XI}$$

$$\mathbf{B}_{57} = -x_{15}\,\mathbf{a}_1 - y_{15}\,\mathbf{a}_2 - z_{15}\,\mathbf{a}_3 = (-x_{15}a - z_{15}c\cos\beta)\,\hat{\mathbf{x}} - y_{15}b\,\hat{\mathbf{y}} - z_{15}c\sin\beta\,\hat{\mathbf{z}} \quad (4g) \quad \text{O XI}$$

$$\mathbf{B}_{58} = x_{15}\,\mathbf{a}_1 - y_{15}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{15}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{15}a + z_{15}c\cos\beta\right)\hat{\mathbf{x}} - y_{15}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{15}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4g) \quad \text{O XI}$$

$$\mathbf{B}_{59} = x_{16}\,\mathbf{a}_1 + y_{16}\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 = (x_{16}a + z_{16}c\cos\beta)\,\hat{\mathbf{x}} + y_{16}b\,\hat{\mathbf{y}} + z_{16}c\sin\beta\,\hat{\mathbf{z}} \quad (4g) \quad \text{O XII}$$

$$\mathbf{B}_{60} = -x_{16}\,\mathbf{a}_1 + y_{16}\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{16}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{16}a - z_{16}c\cos\beta\right)\hat{\mathbf{x}} + y_{16}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{16}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4g) \quad \text{O XII}$$

$$\mathbf{B}_{61} = -x_{16}\,\mathbf{a}_1 - y_{16}\,\mathbf{a}_2 - z_{16}\,\mathbf{a}_3 = (-x_{16}a - z_{16}c\cos\beta)\,\hat{\mathbf{x}} - y_{16}b\,\hat{\mathbf{y}} - z_{16}c\sin\beta\,\hat{\mathbf{z}} \quad (4g) \quad \text{O XII}$$

$$\mathbf{B}_{62} = x_{16}\,\mathbf{a}_1 - y_{16}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{16}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{16}a + z_{16}c\cos\beta\right)\hat{\mathbf{x}} - y_{16}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{16}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4g) \quad \text{O XII}$$

References:

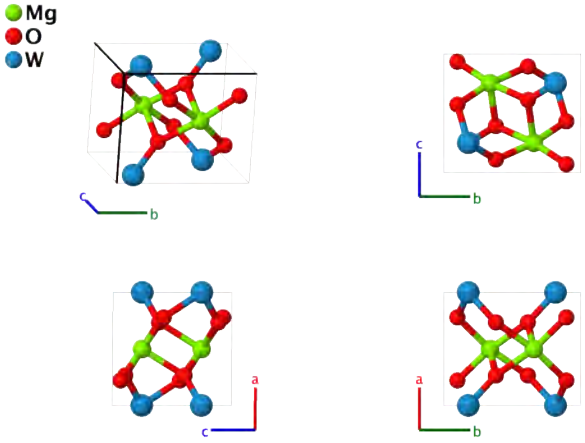
- H. Fujishita, M. Sato, S. Sato, and S. Hoshino, *Structure Determination of low-dimensional conductor Mo_8O_{23}*, J. Solid State Chem. **66**, 40–46 (1987), doi:10.1016/0022-4596(87)90218-0.

Geometry files:

- CIF: pp. 1535
- POSCAR: pp. 1536

Huanzalaite ($MgWO_4$, $H0_6$) Structure: AB4C_mP12_13_f_2g_e

http://aflow.org/prototype-encyclopedia/AB4C_mP12_13_f_2g_e



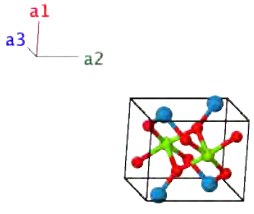
Prototype	:	MgO_4W
AFLOW prototype label	:	AB4C_mP12_13_f_2g_e
***Strukturbericht* designation**	:	$H0_6$
Pearson symbol	:	mP12
Space group number	:	13
Space group symbol	:	$P2/c$
AFLOW prototype command	:	`aflow --proto=AB4C_mP12_13_f_2g_e --params=`$a, b/a, c/a, \beta, y_1, y_2, x_3, y_3, z_3, x_4, y_4, z_4$

Other compounds with this structure

- $CoWO_4$, $FeWO_4$ (ferberite), $MnWO_4$ (hübernite), $NiWO_4$, and $ZnWO_4$

- Most authors refer to this as the wolframite structure, but that name properly belongs to the solid solution $(Mn,Fe)WO_4$. (Herman, 1937) chose this "magnesium wolframite" compound to represent the $H0_6$ structures.

Simple Monoclinic primitive vectors:



$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$

Basis vectors:

	Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$ =	$y_1\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(2e)$	W

$\mathbf{B}_2$	=	$-y_1\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(2e)$	W
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}a + \frac{1}{4}c\cos\beta\right)\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(2f)$	Mg
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}a + \frac{3}{4}c\cos\beta\right)\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(2f)$	Mg
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(4g)$	O I
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_3 a - z_3 c\cos\beta\right)\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4g)$	O I
$\mathbf{B}_7$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(4g)$	O I
$\mathbf{B}_8$	=	$x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4g)$	O I
$\mathbf{B}_9$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(4g)$	O II
$\mathbf{B}_{10}$	=	$-x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_4 a - z_4 c\cos\beta\right)\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4g)$	O II
$\mathbf{B}_{11}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(4g)$	O II
$\mathbf{B}_{12}$	=	$x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4g)$	O II

References:

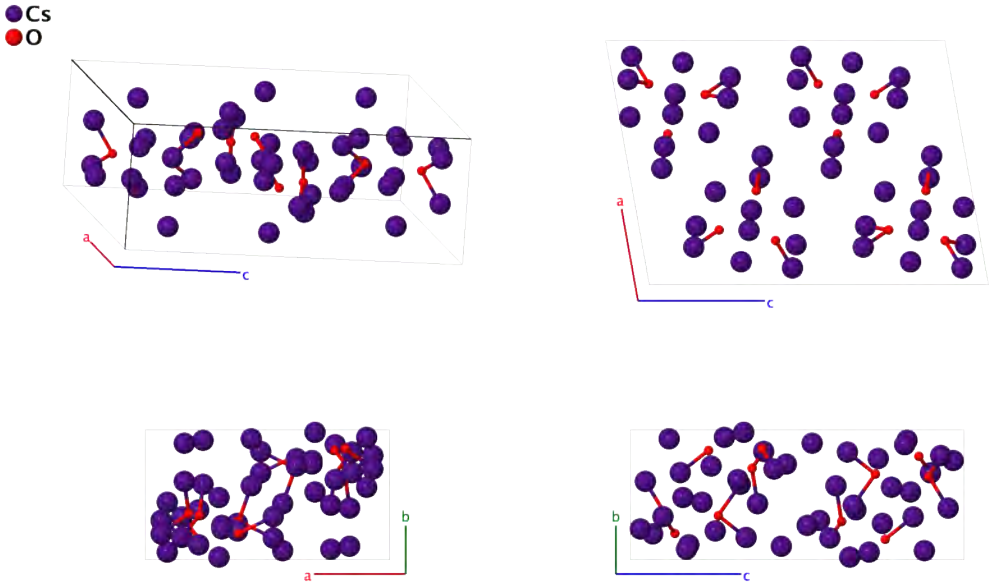
- V. B. Kravchenko, *Crystal structure of the monoclinic form of magnesium tungstate $MgWO_4$*, J. Struct. Chem. **10**, 139–140 (1969), doi:10.1007/BF00751974.
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1536
- POSCAR: pp. 1536

$Cs_{11}O_3$ Structure: A11B3_mP56_14_11e_3e

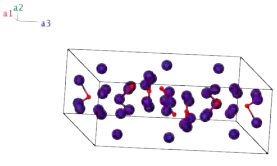
http://aflow.org/prototype-encyclopedia/A11B3_mP56_14_11e_3e



Prototype	:	$Cs_{11}O_3$
AFLOW prototype label	:	A11B3_mP56_14_11e_3e
***Strukturbericht* designation**	:	None
Pearson symbol	:	mP56
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=A11B3_mP56_14_11e_3e --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}$

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cs I

$$\mathbf{B}_2 = -x_1\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_1\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_1\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_1 a - z_1 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_1\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_1\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs I}$$

$$\mathbf{B}_3 = -x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3 = (-x_1 a - z_1 c\cos\beta)\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} - z_1 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs I}$$

$$\mathbf{B}_4 = x_1\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_1\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_1\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_1 a + z_1 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_1\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_1\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs I}$$

$$\mathbf{B}_5 = x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3 = (x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs II}$$

$$\mathbf{B}_6 = -x_2\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_2\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_2\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_2 a - z_2 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_2\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_2\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs II}$$

$$\mathbf{B}_7 = -x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3 = (-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} - z_2 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs II}$$

$$\mathbf{B}_8 = x_2\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_2\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_2\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_2 a + z_2 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_2\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_2\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs II}$$

$$\mathbf{B}_9 = x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3 = (x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs III}$$

$$\mathbf{B}_{10} = -x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_3 a - z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_3\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_3\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs III}$$

$$\mathbf{B}_{11} = -x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3 = (-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs III}$$

$$\mathbf{B}_{12} = x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_3\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_3\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs III}$$

$$\mathbf{B}_{13} = x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3 = (x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs IV}$$

$$\mathbf{B}_{14} = -x_4\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_4 a - z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_4\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_4\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs IV}$$

$$\mathbf{B}_{15} = -x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3 = (-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs IV}$$

$$\mathbf{B}_{16} = x_4\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_4\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_4\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs IV}$$

$$\mathbf{B}_{17} = x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3 = (x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs V}$$

$$\mathbf{B}_{18} = -x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_5 a - z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_5\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_5\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs V}$$

$$\mathbf{B}_{19} = -x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3 = (-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs V}$$

$$\mathbf{B}_{20} = x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_5\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_5\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs V}$$

$$\mathbf{B}_{21} = x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 = (x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs VI}$$

$$\mathbf{B}_{22} = -x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_6 a - z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_6\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_6\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs VI}$$

$$\mathbf{B}_{23} = -x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3 = (-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs VI}$$

$$\mathbf{B}_{24} = x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_6\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs VI}$$

$$\mathbf{B}_{25} = x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 = (x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs VII}$$

$$\mathbf{B}_{26} = -x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_7 a - z_7 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_7\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_7\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs VII}$$

$$\mathbf{B}_{27} = -x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3 = (-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs VII}$$

$$\mathbf{B}_{28} = x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_7 a + z_7 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_7\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_7\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs VII}$$

$$\mathbf{B}_{29} = x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 = (x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs VIII}$$

$$\mathbf{B}_{30} = -x_8\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_8 a - z_8 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_8\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_8\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs VIII}$$

$$\mathbf{B}_{31} = -x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3 = (-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs VIII}$$

$$\mathbf{B}_{32} = x_8\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_8 a + z_8 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_8\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_8\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs VIII}$$

$$\mathbf{B}_{33} = x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 = (x_9 a + z_9 c\cos\beta)\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs IX}$$

$$\mathbf{B}_{34} = -x_9\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_9 a - z_9 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_9\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_9\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs IX}$$

$$\mathbf{B}_{35} = -x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3 = (-x_9 a - z_9 c\cos\beta)\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} - z_9 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs IX}$$

$$\mathbf{B}_{36} = x_9\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_9 a + z_9 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_9\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_9\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs IX}$$

$$\mathbf{B}_{37} = x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 = (x_{10} a + z_{10} c\cos\beta)\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs X}$$

$$\mathbf{B}_{38} = -x_{10}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{10}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{10} a - z_{10} c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{10}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{10}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs X}$$

$$\mathbf{B}_{39} = -x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 = (-x_{10} a - z_{10} c\cos\beta)\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} - z_{10} c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs X}$$

$$\mathbf{B}_{40} = x_{10}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{10}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{10} a + z_{10} c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{10}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{10}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs X}$$

$$\mathbf{B}_{41} = x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 = (x_{11} a + z_{11} c\cos\beta)\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + z_{11} c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs XI}$$

$$\mathbf{B}_{42} = -x_{11}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{11}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{11} a - z_{11} c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{11}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{11}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs XI}$$

$$\mathbf{B}_{43} = -x_{11}\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3 = (-x_{11} a - z_{11} c\cos\beta)\,\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} - z_{11} c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs XI}$$

$$\mathbf{B}_{44} = x_{11}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{11}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{11} a + z_{11} c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{11}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{11}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cs XI}$$

$$\mathbf{B}_{45} = x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 = (x_{12} a + z_{12} c\cos\beta)\,\hat{\mathbf{x}} + y_{12} b\,\hat{\mathbf{y}} + z_{12} c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O I}$$

$\mathbf{B}_{46}$	=	$-x_{12}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{12}\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{12}\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_{12}a - z_{12}c\cos\beta\right)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{12}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{12}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_{47}$	=	$-x_{12}\,\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3$	=	$(-x_{12}a - z_{12}c\cos\beta)\,\hat{\mathbf{x}} - y_{12}b\,\hat{\mathbf{y}} - z_{12}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_{48}$	=	$x_{12}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{12}\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{12}\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{12}a + z_{12}c\cos\beta\right)\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{12}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{12}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_{49}$	=	$x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3$	=	$(x_{13}a + z_{13}c\cos\beta)\,\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} + z_{13}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_{50}$	=	$-x_{13}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{13}\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{13}\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_{13}a - z_{13}c\cos\beta\right)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{13}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{13}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_{51}$	=	$-x_{13}\,\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 - z_{13}\,\mathbf{a}_3$	=	$(-x_{13}a - z_{13}c\cos\beta)\,\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} - z_{13}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_{52}$	=	$x_{13}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{13}\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{13}\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{13}a + z_{13}c\cos\beta\right)\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{13}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{13}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_{53}$	=	$x_{14}\,\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3$	=	$(x_{14}a + z_{14}c\cos\beta)\,\hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} + z_{14}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O III
$\mathbf{B}_{54}$	=	$-x_{14}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{14}\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{14}\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_{14}a - z_{14}c\cos\beta\right)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{14}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{14}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O III
$\mathbf{B}_{55}$	=	$-x_{14}\,\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 - z_{14}\,\mathbf{a}_3$	=	$(-x_{14}a - z_{14}c\cos\beta)\,\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} - z_{14}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O III
$\mathbf{B}_{56}$	=	$x_{14}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{14}\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{14}\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{14}a + z_{14}c\cos\beta\right)\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{14}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{14}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O III

References:

- A. Simon and E. Westerbeck, *Über Suboxide der Alkalimetalle. 10. Das "komplexe Metall" $Cs_{11}O_3$*, Z. Anorg. Allg. Chem. **428**, 187–198 (1977), doi:10.1002/zaac.19774280123.

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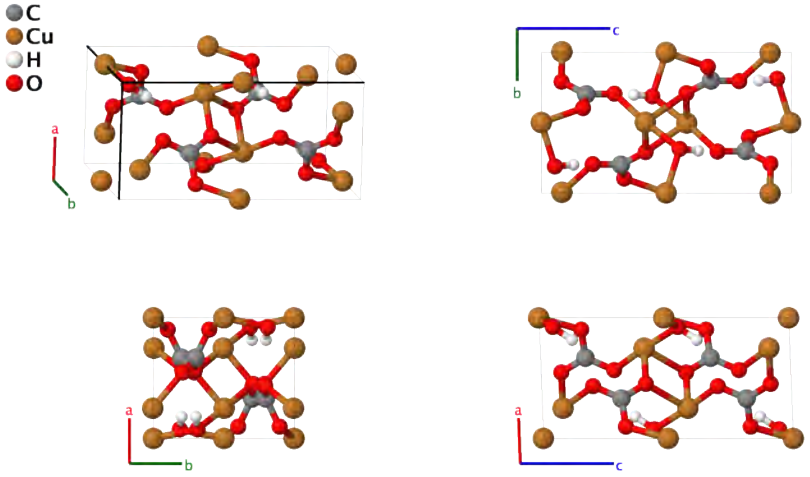
- T. B. Massalski, H. Okamoto, P. R. Subramanian, and L. Kacprzak, eds., *Binary Alloy Phase Diagrams*, vol. 2 (ASM International, Materials Park, Ohio, USA, 1990), 2^{nd} edn. Cd-Ce to Hf-Rb.

Geometry files:

- CIF: pp. 1536
- POSCAR: pp. 1537

Azurite [$Cu_3(CO_3)_2(OH)_2$, $G7_4$] Structure: A2B3C2D8_mP30_14_e_ce_e_4e

http://aflow.org/prototype-encyclopedia/A2B3C2D8_mP30_14_e_ce_e_4e

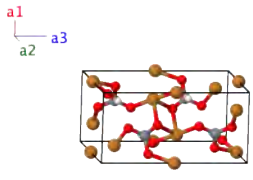


Prototype	:	$C_2Cu_3H_2O_8$
AFLOW prototype label	:	A2B3C2D8_mP30_14_e_ce_e_4e
***Strukturbericht* designation**	:	$G7_4$
Pearson symbol	:	mP30
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=A2B3C2D8_mP30_14_e_ce_e_4e --params=`$a, b/a, c/a, \beta, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8$

- (Rule, 2011) actually present two possible structures for azurite: the current structure, with space group $P2_1/c$ #14, which is identical to the structure designated $G7_4$ by (Hermann, 1937), except that the OH radical is now separated into the O(IV) and H atom sites; and a structure with the same unit cell dimensions but without an inversion site, leading to space group $P2_1$ #4. The actual difference between the two structures is extremely small, and we choose to use the traditional structure.
- Structural information was taken from powder diffraction data at 1.28 K.

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	(2*c*)	Cu I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}b\,\hat{\mathbf{y}}$	(2*c*)	Cu I
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2a + z_2c\cos\beta)\,\hat{\mathbf{x}} + y_2b\,\hat{\mathbf{y}} + z_2c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	C
$\mathbf{B}_4$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2}+y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_2a - z_2c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	C
$\mathbf{B}_5$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2a - z_2c\cos\beta)\,\hat{\mathbf{x}} - y_2b\,\hat{\mathbf{y}} - z_2c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	C
$\mathbf{B}_6$	=	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2}-y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_2a + z_2c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	C
$\mathbf{B}_7$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3a + z_3c\cos\beta)\,\hat{\mathbf{x}} + y_3b\,\hat{\mathbf{y}} + z_3c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Cu II
$\mathbf{B}_8$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2}+y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_3a - z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Cu II
$\mathbf{B}_9$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3a - z_3c\cos\beta)\,\hat{\mathbf{x}} - y_3b\,\hat{\mathbf{y}} - z_3c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Cu II
$\mathbf{B}_{10}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2}-y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_3a + z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Cu II
$\mathbf{B}_{11}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4a + z_4c\cos\beta)\,\hat{\mathbf{x}} + y_4b\,\hat{\mathbf{y}} + z_4c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	H
$\mathbf{B}_{12}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2}+y_4\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_4a - z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	H
$\mathbf{B}_{13}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4a - z_4c\cos\beta)\,\hat{\mathbf{x}} - y_4b\,\hat{\mathbf{y}} - z_4c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	H
$\mathbf{B}_{14}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2}-y_4\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_4a + z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	H
$\mathbf{B}_{15}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5a + z_5c\cos\beta)\,\hat{\mathbf{x}} + y_5b\,\hat{\mathbf{y}} + z_5c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	O I
$\mathbf{B}_{16}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2}+y_5\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_5a - z_5c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	O I
$\mathbf{B}_{17}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5a - z_5c\cos\beta)\,\hat{\mathbf{x}} - y_5b\,\hat{\mathbf{y}} - z_5c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	O I
$\mathbf{B}_{18}$	=	$x_5\,\mathbf{a}_1 + \left(\frac{1}{2}-y_5\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_5a + z_5c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	O I
$\mathbf{B}_{19}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6a + z_6c\cos\beta)\,\hat{\mathbf{x}} + y_6b\,\hat{\mathbf{y}} + z_6c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	O II
$\mathbf{B}_{20}$	=	$-x_6\,\mathbf{a}_1 + \left(\frac{1}{2}+y_6\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_6a - z_6c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_6\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	O II
$\mathbf{B}_{21}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6a - z_6c\cos\beta)\,\hat{\mathbf{x}} - y_6b\,\hat{\mathbf{y}} - z_6c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	O II
$\mathbf{B}_{22}$	=	$x_6\,\mathbf{a}_1 + \left(\frac{1}{2}-y_6\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_6a + z_6c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_6\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	O II

$$\mathbf{B}_{23} = x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 = (x_7a + z_7c\cos\beta)\,\hat{\mathbf{x}} + y_7b\,\hat{\mathbf{y}} + z_7c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O III}$$

$$\mathbf{B}_{24} = -x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_7a - z_7c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_7\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_7\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O III}$$

$$\mathbf{B}_{25} = -x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3 = (-x_7a - z_7c\cos\beta)\,\hat{\mathbf{x}} - y_7b\,\hat{\mathbf{y}} - z_7c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O III}$$

$$\mathbf{B}_{26} = x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_7a + z_7c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_7\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_7\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O III}$$

$$\mathbf{B}_{27} = x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 = (x_8a + z_8c\cos\beta)\,\hat{\mathbf{x}} + y_8b\,\hat{\mathbf{y}} + z_8c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O IV}$$

$$\mathbf{B}_{28} = -x_8\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_8a - z_8c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_8\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_8\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O IV}$$

$$\mathbf{B}_{29} = -x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3 = (-x_8a - z_8c\cos\beta)\,\hat{\mathbf{x}} - y_8b\,\hat{\mathbf{y}} - z_8c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O IV}$$

$$\mathbf{B}_{30} = x_8\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_8a + z_8c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_8\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_8\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O IV}$$

References:

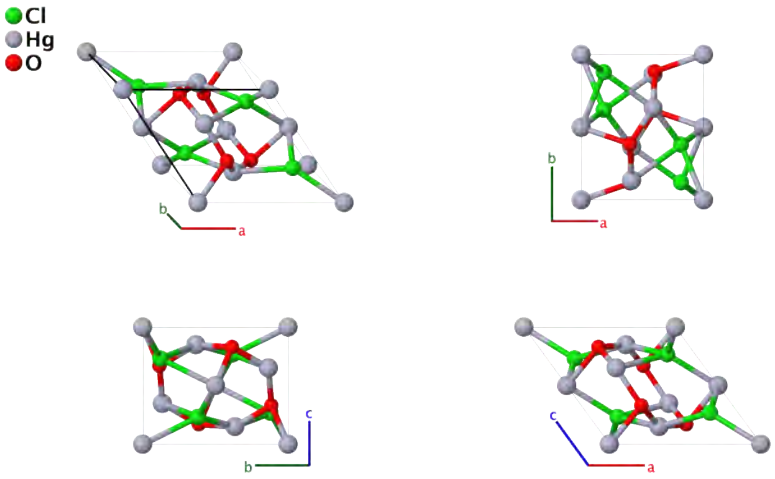
- K. C. Rule, M. Reehuis, M. C. R. Gibson, B. Ouladdiaf, M. J. Gutmann, J.-U. Hoffmann, S. Gerischer, D. A. Tennant, S. Süllow, and M. Lang, *Magnetic and crystal structure of azurite $Cu_3(CO_3)_2(OH)_2$ as determined by neutron diffraction*, Phys. Rev. B **83**, 104401 (2011), doi:10.1103/PhysRevB.83.104401.
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1537
- POSCAR: pp. 1538

$HgCl_2 \cdot 2HgO$ Structure: A2B3C2_mP14_14_e_ae_e

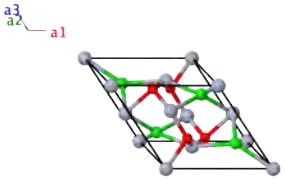
http://aflow.org/prototype-encyclopedia/A2B3C2_mP14_14_e_ae_e



Prototype	:	$Cl_2Hg_3O_2$
AFLOW prototype label	:	A2B3C2_mP14_14_e_ae_e
***Strukturbericht* designation**	:	None
Pearson symbol	:	mP14
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=A2B3C2_mP14_14_e_ae_e --params=`$a, b/a, c/a, \beta, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4$

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	Hg I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Hg I
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2a + z_2c\cos\beta)\,\hat{\mathbf{x}} + y_2b\,\hat{\mathbf{y}} + z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cl
$\mathbf{B}_4$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_2a - z_2c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cl
$\mathbf{B}_5$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2a - z_2c\cos\beta)\,\hat{\mathbf{x}} - y_2b\,\hat{\mathbf{y}} - z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cl

$\mathbf{B}_6$	=	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_2 a + z_2 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Cl
$\mathbf{B}_7$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Hg II
$\mathbf{B}_8$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_3 a - z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Hg II
$\mathbf{B}_9$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Hg II
$\mathbf{B}_{10}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Hg II
$\mathbf{B}_{11}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	O
$\mathbf{B}_{12}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_4 a - z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	O
$\mathbf{B}_{13}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	O
$\mathbf{B}_{14}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	O

References:

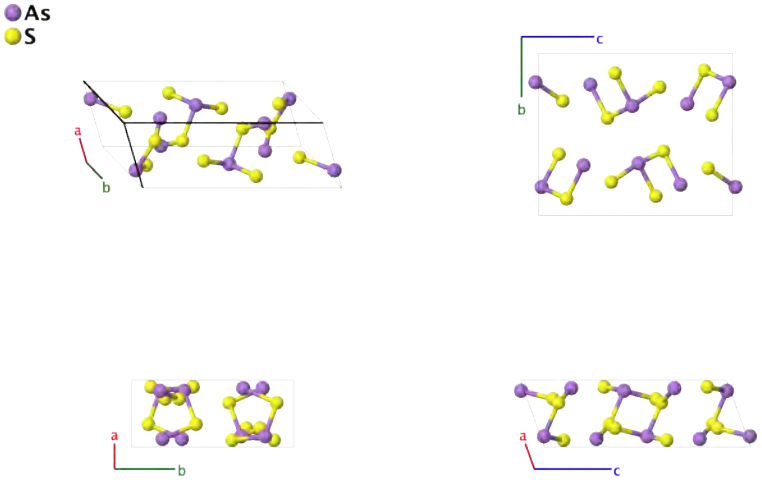
- S. Šćavnićar, *The crystal structure of trimercuric oxychloride, $HgCl_2{\cdot}2HgO$*, Acta Cryst. **8**, 379–383 (1955), doi:10.1107/S0365110X55001278.

Geometry files:

- CIF: pp. 1538
- POSCAR: pp. 1538

Orpiment (As_2S_3, $D5_f$) Structure: A2B3_mP20_14_2e_3e

http://aflow.org/prototype-encyclopedia/A2B3_mP20_14_2e_3e



Prototype	:	As_2S_3
AFLOW prototype label	:	A2B3_mP20_14_2e_3e
***Strukturbericht* designation**	:	$D5_f$
Pearson symbol	:	mP20
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=A2B3_mP20_14_2e_3e --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5$

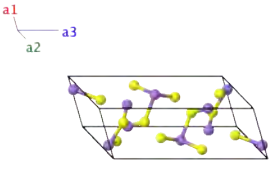
Other compounds with this structure

- As_2Se_3 and As_2Te_3

- (Morimoto, 1954) gives the structure in setting $P2_1/n$ of space group #14. We used FINDSYM to translate the structure to the standard $P2_1/c$ setting. This included a redefinition of the primitive vectors of the lattice.

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

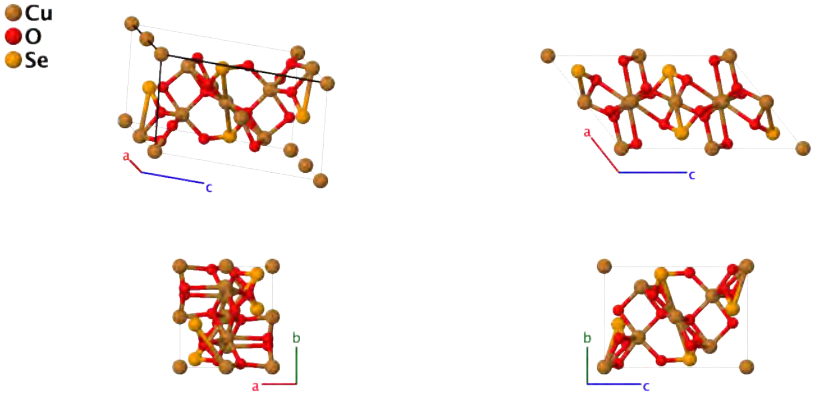
		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1a + z_1c\cos\beta)\,\hat{\mathbf{x}} + y_1b\,\hat{\mathbf{y}} + z_1c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As I
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2}+y_1\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_1a - z_1c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As I
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$(-x_1a - z_1c\cos\beta)\,\hat{\mathbf{x}} - y_1b\,\hat{\mathbf{y}} - z_1c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As I
$\mathbf{B}_4$	=	$x_1\,\mathbf{a}_1 + \left(\frac{1}{2}-y_1\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_1a + z_1c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As I
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2a + z_2c\cos\beta)\,\hat{\mathbf{x}} + y_2b\,\hat{\mathbf{y}} + z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As II
$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2}+y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_2a - z_2c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As II
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2a - z_2c\cos\beta)\,\hat{\mathbf{x}} - y_2b\,\hat{\mathbf{y}} - z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As II
$\mathbf{B}_8$	=	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2}-y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_2a + z_2c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As II
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3a + z_3c\cos\beta)\,\hat{\mathbf{x}} + y_3b\,\hat{\mathbf{y}} + z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S I
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2}+y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_3a - z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S I
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3a - z_3c\cos\beta)\,\hat{\mathbf{x}} - y_3b\,\hat{\mathbf{y}} - z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S I
$\mathbf{B}_{12}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2}-y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_3a + z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S I
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4a + z_4c\cos\beta)\,\hat{\mathbf{x}} + y_4b\,\hat{\mathbf{y}} + z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S II
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2}+y_4\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_4a - z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S II
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4a - z_4c\cos\beta)\,\hat{\mathbf{x}} - y_4b\,\hat{\mathbf{y}} - z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S II
$\mathbf{B}_{16}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2}-y_4\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_4a + z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S II
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5a + z_5c\cos\beta)\,\hat{\mathbf{x}} + y_5b\,\hat{\mathbf{y}} + z_5c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S III
$\mathbf{B}_{18}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2}+y_5\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_5a - z_5c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S III
$\mathbf{B}_{19}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5a - z_5c\cos\beta)\,\hat{\mathbf{x}} - y_5b\,\hat{\mathbf{y}} - z_5c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S III
$\mathbf{B}_{20}$	=	$x_5\,\mathbf{a}_1 + \left(\frac{1}{2}-y_5\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_5a + z_5c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S III

References:
- N. Morimoto, *The Crystal Structure of Orpiment (As_2S_3) Refined*, Mineral. J. **1**, 160–169 (1954), doi:10.2465/minerj1953.1.160.

Geometry files:
- CIF: pp. 1538
- POSCAR: pp. 1539

Monoclinic Cu_2OSeO_3 Structure: A2B4C_mP28_14_abe_4e_e

http://aflow.org/prototype-encyclopedia/A2B4C_mP28_14_abe_4e_e

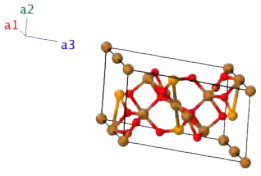


Prototype	:	Cu_2O_4Se
AFLOW prototype label	:	A2B4C_mP28_14_abe_4e_e
***Strukturbericht* designation**	:	None
Pearson symbol	:	mP28
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=A2B4C_mP28_14_abe_4e_e --params=`$a, b/a, c/a, \beta, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8$

- This is the monoclinic phase of Cu_2OSeO_3. There is also a cubic phase.

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	Cu I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Cu I
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	$(2b)$	Cu II
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(a + c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	$(2b)$	Cu II
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3a + z_3c\cos\beta)\,\hat{\mathbf{x}} + y_3b\,\hat{\mathbf{y}} + z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cu III

$$\mathbf{B}_{6} = -x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_3 a - z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_3\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_3\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cu III}$$

$$\mathbf{B}_{7} = -x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3 = (-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cu III}$$

$$\mathbf{B}_{8} = x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_3\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_3\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cu III}$$

$$\mathbf{B}_{9} = x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3 = (x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O I}$$

$$\mathbf{B}_{10} = -x_4\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_4 a - z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_4\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_4\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O I}$$

$$\mathbf{B}_{11} = -x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3 = (-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O I}$$

$$\mathbf{B}_{12} = x_4\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_4\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_4\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O I}$$

$$\mathbf{B}_{13} = x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3 = (x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O II}$$

$$\mathbf{B}_{14} = -x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_5 a - z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_5\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_5\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O II}$$

$$\mathbf{B}_{15} = -x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3 = (-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O II}$$

$$\mathbf{B}_{16} = x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_5\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_5\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O II}$$

$$\mathbf{B}_{17} = x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 = (x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O III}$$

$$\mathbf{B}_{18} = -x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_6 a - z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_6\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_6\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O III}$$

$$\mathbf{B}_{19} = -x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3 = (-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O III}$$

$$\mathbf{B}_{20} = x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_6\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O III}$$

$$\mathbf{B}_{21} = x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 = (x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O IV}$$

$$\mathbf{B}_{22} = -x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_7 a - z_7 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_7\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_7\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O IV}$$

$$\mathbf{B}_{23} = -x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3 = (-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O IV}$$

$$\mathbf{B}_{24} = x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_7 a + z_7 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_7\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_7\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O IV}$$

$$\mathbf{B}_{25} = x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 = (x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Se}$$

$$\mathbf{B}_{26} = -x_8\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_8 a - z_8 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_8\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_8\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Se}$$

$$\mathbf{B}_{27} = -x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3 = (-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Se}$$

$$\mathbf{B}_{28} = x_8\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_8 a + z_8 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_8\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_8\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Se}$$

References:

- H. Effenberger and F. Pertlik, *Die Kristallstrukturen der Kupfer(II)-oxo-selenite $Cu_2O(SeO_3)$ (kubisch und monoklin) und $Cu_4O(SeO_3)_3$ (monoklin und triklin)*, Monatshefte für Chemie - Chemical Monthly **117**, 887–896 (1986), doi:10.1007/BF00811258.

Found in:

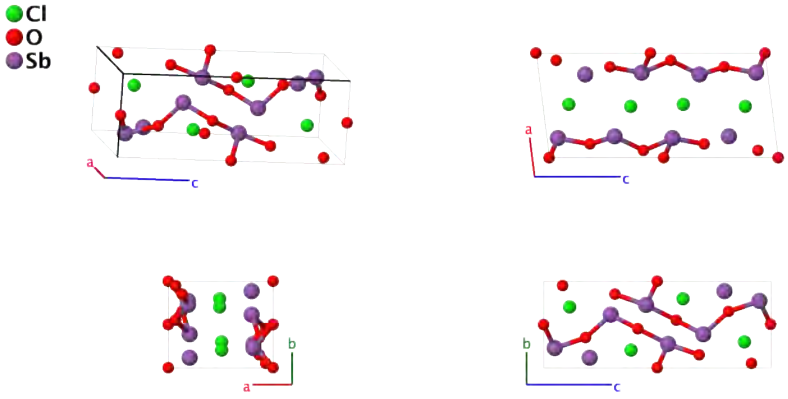
- P. Y. Portnichenko, J. Romhányi, Y. A. Onykiienko, A. Henschel, M. Schmidt, A. S. Cameron, M. A. Surmach, J. A. Lim, J. T. Park, A. Schneidewind, D. L. Abernathy, H. Rosner, J. van den Brink, and D. S. Inosov, *Magnon spectrum of the helimagnetic insulator Cu_2OSeO_3*, Nat. Commun. **7**, 10725 (2016), doi:10.1038/ncomms10725.

Geometry files:

- CIF: pp. 1539
- POSCAR: pp. 1539

$Sb_4O_5Cl_2$ Structure: A2B5C4_mP22_14_e_c2e_2e

http://aflow.org/prototype-encyclopedia/A2B5C4_mP22_14_e_c2e_2e



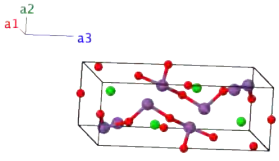
Prototype	:	$Cl_2O_5Sb_4$
AFLOW prototype label	:	A2B5C4_mP22_14_e_c2e_2e
***Strukturbericht* designation**	:	None
Pearson symbol	:	mP22
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=A2B5C4_mP22_14_e_c2e_2e --params=`$a, b/a, c/a, \beta, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6$

Other compounds with this structure

- $Sb_4O_5Br_2$

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	$(2c)$	O I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}b\,\hat{\mathbf{y}}$	$(2c)$	O I
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2a + z_2c\cos\beta)\,\hat{\mathbf{x}} + y_2b\,\hat{\mathbf{y}} + z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cl
$\mathbf{B}_4$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_2a - z_2c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cl
$\mathbf{B}_5$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2a - z_2c\cos\beta)\,\hat{\mathbf{x}} - y_2b\,\hat{\mathbf{y}} - z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cl

$\mathbf{B}_6$	=	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_2 a + z_2 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Cl
$\mathbf{B}_7$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	O II
$\mathbf{B}_8$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_3 a - z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	O II
$\mathbf{B}_9$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	O II
$\mathbf{B}_{10}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	O II
$\mathbf{B}_{11}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	O III
$\mathbf{B}_{12}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_4 a - z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	O III
$\mathbf{B}_{13}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	O III
$\mathbf{B}_{14}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	O III
$\mathbf{B}_{15}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Sb I
$\mathbf{B}_{16}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_5 a - z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Sb I
$\mathbf{B}_{17}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Sb I
$\mathbf{B}_{18}$	=	$x_5\,\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Sb I
$\mathbf{B}_{19}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Sb II
$\mathbf{B}_{20}$	=	$-x_6\,\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_6 a - z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Sb II
$\mathbf{B}_{21}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Sb II
$\mathbf{B}_{22}$	=	$x_6\,\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Sb II

References:

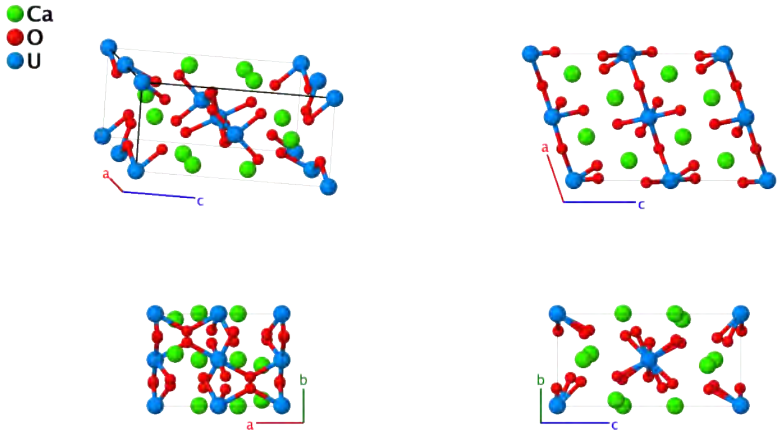
- C. Särnstrand, *The crystal structure of antimony(III) chloride oxide $Sb_4O_5Cl_2$*, Acta Crystallogr. Sect. B Struct. Sci. **34**, 2402–2407 (1978), doi:10.1107/S056774087800833X.

Geometry files:

- CIF: pp. 1539
- POSCAR: pp. 1540

Ca_2UO_5 Structure: A2B5C_mP32_14_2e_5e_ab

http://aflow.org/prototype-encyclopedia/A2B5C_mP32_14_2e_5e_ab



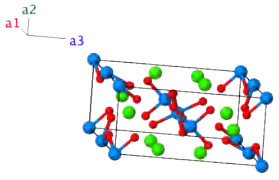
Prototype	:	Ca_2O_5U
AFLOW prototype label	:	A2B5C_mP32_14_2e_5e_ab
***Strukturbericht* designation**	:	None
Pearson symbol	:	mP32
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=A2B5C_mP32_14_2e_5e_ab --params=`$a, b/a, c/a, \beta, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9$

Other compounds with this structure

- Sr_2UO_5

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	U I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	U I
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	$(2b)$	U II
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(a + c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	$(2b)$	U II

$$\mathbf{B}_5 = x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3 = (x_3a + z_3c\cos\beta)\,\hat{\mathbf{x}} + y_3b\,\hat{\mathbf{y}} + z_3c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Ca I}$$

$$\mathbf{B}_6 = -x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_3a - z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_3\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_3\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Ca I}$$

$$\mathbf{B}_7 = -x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3 = (-x_3a - z_3c\cos\beta)\,\hat{\mathbf{x}} - y_3b\,\hat{\mathbf{y}} - z_3c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Ca I}$$

$$\mathbf{B}_8 = x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_3a + z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_3\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_3\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Ca I}$$

$$\mathbf{B}_9 = x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3 = (x_4a + z_4c\cos\beta)\,\hat{\mathbf{x}} + y_4b\,\hat{\mathbf{y}} + z_4c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Ca II}$$

$$\mathbf{B}_{10} = -x_4\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_4a - z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_4\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_4\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Ca II}$$

$$\mathbf{B}_{11} = -x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3 = (-x_4a - z_4c\cos\beta)\,\hat{\mathbf{x}} - y_4b\,\hat{\mathbf{y}} - z_4c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Ca II}$$

$$\mathbf{B}_{12} = x_4\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_4a + z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_4\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_4\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Ca II}$$

$$\mathbf{B}_{13} = x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3 = (x_5a + z_5c\cos\beta)\,\hat{\mathbf{x}} + y_5b\,\hat{\mathbf{y}} + z_5c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O I}$$

$$\mathbf{B}_{14} = -x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_5a - z_5c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_5\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_5\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O I}$$

$$\mathbf{B}_{15} = -x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3 = (-x_5a - z_5c\cos\beta)\,\hat{\mathbf{x}} - y_5b\,\hat{\mathbf{y}} - z_5c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O I}$$

$$\mathbf{B}_{16} = x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_5a + z_5c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_5\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_5\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O I}$$

$$\mathbf{B}_{17} = x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 = (x_6a + z_6c\cos\beta)\,\hat{\mathbf{x}} + y_6b\,\hat{\mathbf{y}} + z_6c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O II}$$

$$\mathbf{B}_{18} = -x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_6a - z_6c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_6\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_6\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O II}$$

$$\mathbf{B}_{19} = -x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3 = (-x_6a - z_6c\cos\beta)\,\hat{\mathbf{x}} - y_6b\,\hat{\mathbf{y}} - z_6c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O II}$$

$$\mathbf{B}_{20} = x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_6a + z_6c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_6\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O II}$$

$$\mathbf{B}_{21} = x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 = (x_7a + z_7c\cos\beta)\,\hat{\mathbf{x}} + y_7b\,\hat{\mathbf{y}} + z_7c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O III}$$

$$\mathbf{B}_{22} = -x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_7a - z_7c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_7\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_7\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O III}$$

$$\mathbf{B}_{23} = -x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3 = (-x_7a - z_7c\cos\beta)\,\hat{\mathbf{x}} - y_7b\,\hat{\mathbf{y}} - z_7c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O III}$$

$$\mathbf{B}_{24} = x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_7a + z_7c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_7\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_7\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O III}$$

$$\mathbf{B}_{25} = x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 = (x_8a + z_8c\cos\beta)\,\hat{\mathbf{x}} + y_8b\,\hat{\mathbf{y}} + z_8c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O IV}$$

$$\mathbf{B}_{26} = -x_8\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_8a - z_8c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_8\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_8\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O IV}$$

$\mathbf{B}_{27}$	=	$-x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8a - z_8c\cos\beta)\,\hat{\mathbf{x}} - y_8b\,\hat{\mathbf{y}} - z_8c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O IV
$\mathbf{B}_{28}$	=	$x_8\,\mathbf{a}_1 + \left(\frac{1}{2} - y_8\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_8a + z_8c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_8\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O IV
$\mathbf{B}_{29}$	=	$x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$(x_9a + z_9c\cos\beta)\,\hat{\mathbf{x}} + y_9b\,\hat{\mathbf{y}} + z_9c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O V
$\mathbf{B}_{30}$	=	$-x_9\,\mathbf{a}_1 + \left(\frac{1}{2} + y_9\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_9a - z_9c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_9\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_9\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O V
$\mathbf{B}_{31}$	=	$-x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$(-x_9a - z_9c\cos\beta)\,\hat{\mathbf{x}} - y_9b\,\hat{\mathbf{y}} - z_9c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O V
$\mathbf{B}_{32}$	=	$x_9\,\mathbf{a}_1 + \left(\frac{1}{2} - y_9\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_9a + z_9c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_9\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O V

References:

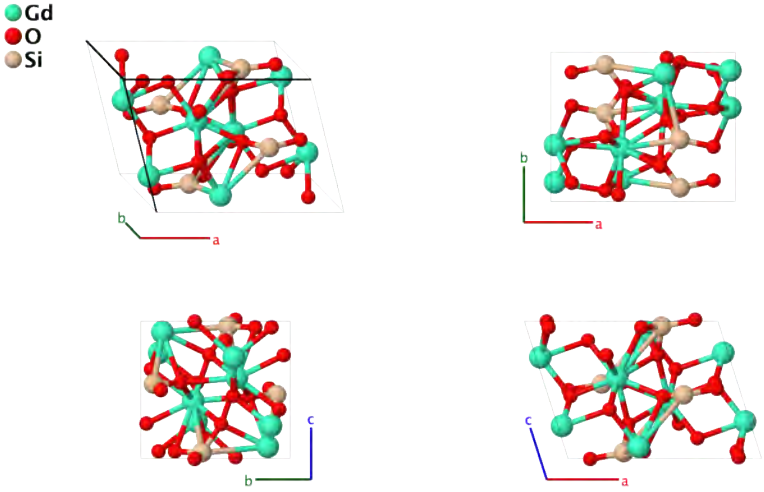
- B. O. Loopstra and H. M. Rietveld, *The structure of some alkaline-earth metal uranates*, Acta Crystallogr. Sect. B Struct. Sci. **25**, 787–791 (1969), doi:10.1107/S0567740869002974.

Geometry files:

- CIF: pp. 1540
- POSCAR: pp. 1540

Gd_2SiO_5 (RE_2SiO_5 X1) Structure: A2B5C_mP32_14_2e_5e_e

http://aflow.org/prototype-encyclopedia/A2B5C_mP32_14_2e_5e_e



Prototype	:	Gd_2O_5Si
AFLOW prototype label	:	A2B5C_mP32_14_2e_5e_e
***Strukturbericht* designation**	:	None
Pearson symbol	:	mP32
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=A2B5C_mP32_14_2e_5e_e --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8$

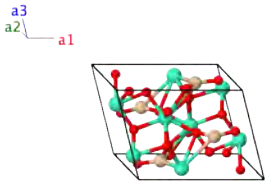
Other compounds with this structure

- Dy_2SiO_5, La_2SiO_5, Tb_2SiO_5, and Y_2SiO_5

- Compounds of the form $RESiO_5$ (RE = Rare Earth and related elements) crystallize in one of two forms (Wang, 2001 and Tian, 2016): X1, shown here, for rare earths between lanthanum and gadolinium, and X2, space group $C2/c$ #15 for the later rare earths, with the dysprosium, yttrium and ytterbium compounds existing in both structures.

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1a + z_1c\cos\beta)\,\hat{\mathbf{x}} + y_1b\,\hat{\mathbf{y}} + z_1c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Gd I
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2}+y_1\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_1a - z_1c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Gd I
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$(-x_1a - z_1c\cos\beta)\,\hat{\mathbf{x}} - y_1b\,\hat{\mathbf{y}} - z_1c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Gd I
$\mathbf{B}_4$	=	$x_1\,\mathbf{a}_1 + \left(\frac{1}{2}-y_1\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_1a + z_1c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Gd I
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2a + z_2c\cos\beta)\,\hat{\mathbf{x}} + y_2b\,\hat{\mathbf{y}} + z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Gd II
$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2}+y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_2a - z_2c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Gd II
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2a - z_2c\cos\beta)\,\hat{\mathbf{x}} - y_2b\,\hat{\mathbf{y}} - z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Gd II
$\mathbf{B}_8$	=	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2}-y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_2a + z_2c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Gd II
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3a + z_3c\cos\beta)\,\hat{\mathbf{x}} + y_3b\,\hat{\mathbf{y}} + z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2}+y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_3a - z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3a - z_3c\cos\beta)\,\hat{\mathbf{x}} - y_3b\,\hat{\mathbf{y}} - z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_{12}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2}-y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_3a + z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4a + z_4c\cos\beta)\,\hat{\mathbf{x}} + y_4b\,\hat{\mathbf{y}} + z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2}+y_4\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_4a - z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4a - z_4c\cos\beta)\,\hat{\mathbf{x}} - y_4b\,\hat{\mathbf{y}} - z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_{16}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2}-y_4\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_4a + z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5a + z_5c\cos\beta)\,\hat{\mathbf{x}} + y_5b\,\hat{\mathbf{y}} + z_5c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O III
$\mathbf{B}_{18}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2}+y_5\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_5a - z_5c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O III
$\mathbf{B}_{19}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5a - z_5c\cos\beta)\,\hat{\mathbf{x}} - y_5b\,\hat{\mathbf{y}} - z_5c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O III
$\mathbf{B}_{20}$	=	$x_5\,\mathbf{a}_1 + \left(\frac{1}{2}-y_5\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_5a + z_5c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O III
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6a + z_6c\cos\beta)\,\hat{\mathbf{x}} + y_6b\,\hat{\mathbf{y}} + z_6c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O IV

$\mathbf{B}_{22}$	=	$-x_6\,\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_6 a - z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	O IV
$\mathbf{B}_{23}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	O IV
$\mathbf{B}_{24}$	=	$x_6\,\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	O IV
$\mathbf{B}_{25}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	O V
$\mathbf{B}_{26}$	=	$-x_7\,\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_7 a - z_7 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	O V
$\mathbf{B}_{27}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	O V
$\mathbf{B}_{28}$	=	$x_7\,\mathbf{a}_1 + \left(\frac{1}{2} - y_7\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_7 a + z_7 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_7\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	O V
$\mathbf{B}_{29}$	=	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Si
$\mathbf{B}_{30}$	=	$-x_8\,\mathbf{a}_1 + \left(\frac{1}{2} + y_8\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_8 a - z_8 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_8\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Si
$\mathbf{B}_{31}$	=	$-x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Si
$\mathbf{B}_{32}$	=	$x_8\,\mathbf{a}_1 + \left(\frac{1}{2} - y_8\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_8 a + z_8 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_8\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Si

References:

- G. V. Anan'eva, A. M. Korovkin, T. I. Merkulyaeva, A. M. Morozova, M. V. Petrov, I. R. Savinova, V. R. Startsev, and P. P. Feofilov, *Growth of lanthanide oxyorthosilicate single crystals, and their structural and optical characteristics*, Inorg. Mat. **17**, 754–758 (1981). Translated from *Neorganicheskie Materialy*.
- J. Wang, S. Tian, G. Li, F. Liao, and X. Jing, *Preparation and X-ray characterization of low-temperature phases of R_2SiO_5 (R = rare earth elements)*, Mater. Res. Bull. **36**, 1855–1861 (2001), doi:10.1016/S0025-5408(01)00664-X.
- Z. Tian, L. Zheng, J. Wang, P. Wan, J. Li, and J. Wang, *Theoretical and experimental determination of the major thermo-mechanical properties of RE_2SiO_5 (RE = Tb, Dy, Ho, Er, Tm, Yb, Lu, and Y) for environmental and thermal barrier coating applications*, J. Am. Ceram. Soc. **36**, 189–202 (2016), doi:10.1016/j.jeurceramsoc.2015.09.013.

Found in:

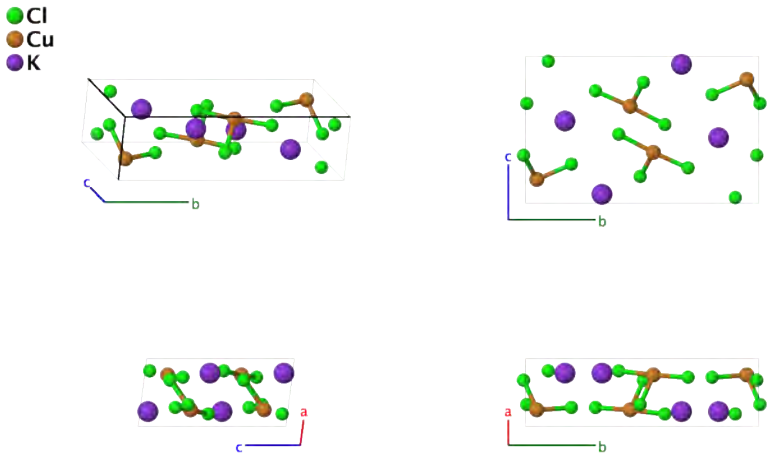
- P. Villars (Chief Editor), *Gd_2SiO_5 ($Gd_2[SiO_4]O$) Crystal Structure*, http://materials.springer.com/isp/crystallographic/docs/sd_0309353 (2016). PAULING FILE in: Inorganic Solid Phases, SpringerMaterials (online database), Springer, Heidelberg (ed.) SpringerMaterials.

Geometry files:

- CIF: pp. 1540
- POSCAR: pp. 1541

Sanguite ($KCuCl_3$) Structure: A3BC_mP20_14_3e_e_e

http://aflow.org/prototype-encyclopedia/A3BC_mP20_14_3e_e_e



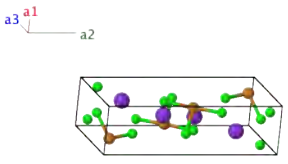
Prototype	:	Cl_3CuK
AFLOW prototype label	:	A3BC_mP20_14_3e_e_e
***Strukturbericht* designation**	:	None
Pearson symbol	:	mP20
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=A3BC_mP20_14_3e_e_e --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5$

Other compounds with this structure

- NH_4CuCl_3

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cl I
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2} + y_1\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_1 a - z_1 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_1\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cl I
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$(-x_1 a - z_1 c\cos\beta)\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} - z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cl I

$$
\begin{array}{lcccccc}
\mathbf{B}_4 & = & x_1\,\mathbf{a}_1 + \left(\frac{1}{2} - y_1\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3 & = & \left(\frac{1}{2}c\cos\beta + x_1 a + z_1 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{Cl I} \\
\mathbf{B}_5 & = & x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3 & = & (x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{Cl II} \\
\mathbf{B}_6 & = & -x_2\,\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3 & = & \left(\frac{1}{2}c\cos\beta - x_2 a - z_2 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{Cl II} \\
\mathbf{B}_7 & = & -x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3 & = & (-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} - z_2 c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{Cl II} \\
\mathbf{B}_8 & = & x_2\,\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3 & = & \left(\frac{1}{2}c\cos\beta + x_2 a + z_2 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{Cl II} \\
\mathbf{B}_9 & = & x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3 & = & (x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{Cl III} \\
\mathbf{B}_{10} & = & -x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3 & = & \left(\frac{1}{2}c\cos\beta - x_3 a - z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{Cl III} \\
\mathbf{B}_{11} & = & -x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3 & = & (-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{Cl III} \\
\mathbf{B}_{12} & = & x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3 & = & \left(\frac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{Cl III} \\
\mathbf{B}_{13} & = & x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3 & = & (x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{Cu} \\
\mathbf{B}_{14} & = & -x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3 & = & \left(\frac{1}{2}c\cos\beta - x_4 a - z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{Cu} \\
\mathbf{B}_{15} & = & -x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3 & = & (-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{Cu} \\
\mathbf{B}_{16} & = & x_4\,\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3 & = & \left(\frac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{Cu} \\
\mathbf{B}_{17} & = & x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3 & = & (x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{K} \\
\mathbf{B}_{18} & = & -x_5\,\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3 & = & \left(\frac{1}{2}c\cos\beta - x_5 a - z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{K} \\
\mathbf{B}_{19} & = & -x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3 & = & (-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{K} \\
\mathbf{B}_{20} & = & x_5\,\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3 & = & \left(\frac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{K}
\end{array}
$$

References:

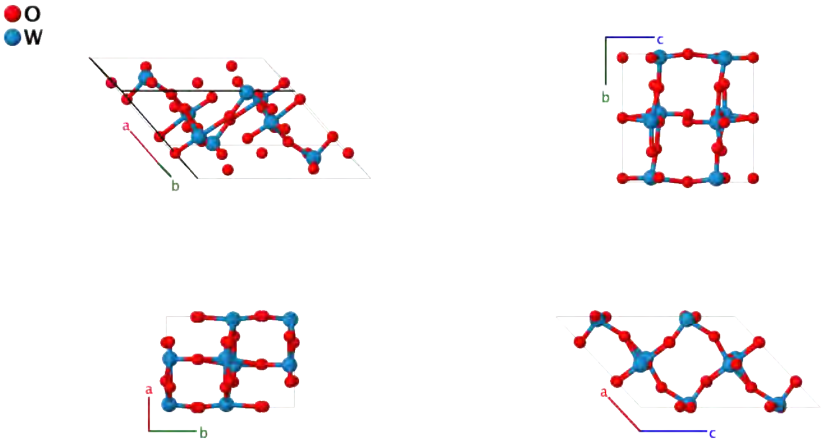
- R. D. Willett, C. Dwiggins, Jr., R. F. Kruh, and R. E. Rundle, *Crystal Structures of $KCuCl_3$ and NH_4CuCl_3*, J. Chem. Phys. **38**, 2429–2436 (1963), doi:10.1063/1.1733520.

Geometry files:

- CIF: pp. 1541
- POSCAR: pp. 1541

γ-WO_3 Structure: A3B_mP32_14_6e_2e

http://aflow.org/prototype-encyclopedia/A3B_mP32_14_6e_2e



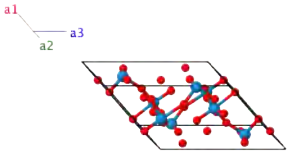
Prototype	:	O_3W
AFLOW prototype label	:	A3B_mP32_14_6e_2e
***Strukturbericht* designation**	:	None
Pearson symbol	:	mP32
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=A3B_mP32_14_6e_2e --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8$

- All stable phases of WO_3 are distortions of the cubic α-ReO_3 ($D0_9$) phase. Based on (Woodward, 1997 and Vogt, 1999), the known stable phases and their approximate temperature ranges are:
  - α-WO_3 (1010-1170 K) (Vogt, 1999)
  - β-WO_3 (600-1170 K) (Vogt, 1999)
  - γ-WO_3 (290-600 K) (Vogt, 1999), this structure
  - δ-WO_3 (230-290 K) (Diehl, 1978)
  - ϵ-WO_3 (below 23 K) (Woodward, 1997)

  Woodward's caption says "The transition temperatures display large hysteresis effects and universal agreement is not found in the literature."
- In addition, several other structures have been proposed and/or found:
  - The original $D0_{10}$ structure (Bräkken, 1931), (Hermann, 1937) superseded by δ-WO_3
  - Original β-WO_3 (Salje, 1977)
  - Hexagonal WO_3 (Gerand, 1979) (metastable)

Simple Monoclinic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= a\,\hat{\mathbf{x}}\\ \mathbf{a}_2 &= b\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	$(x_1a + z_1c\cos\beta)\,\hat{\mathbf{x}} + y_1b\,\hat{\mathbf{y}} + z_1c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_2$	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2}+y_1\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_1\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta - x_1a - z_1c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_3$	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	$(-x_1a - z_1c\cos\beta)\,\hat{\mathbf{x}} - y_1b\,\hat{\mathbf{y}} - z_1c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_4$	$x_1\,\mathbf{a}_1 + \left(\frac{1}{2}-y_1\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_1\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_1a + z_1c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_5$	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	$(x_2a + z_2c\cos\beta)\,\hat{\mathbf{x}} + y_2b\,\hat{\mathbf{y}} + z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_6$	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2}+y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_2\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta - x_2a - z_2c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_7$	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	$(-x_2a - z_2c\cos\beta)\,\hat{\mathbf{x}} - y_2b\,\hat{\mathbf{y}} - z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_8$	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2}-y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_2a + z_2c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_9$	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$(x_3a + z_3c\cos\beta)\,\hat{\mathbf{x}} + y_3b\,\hat{\mathbf{y}} + z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O III
$\mathbf{B}_{10}$	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2}+y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_3\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta - x_3a - z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O III
$\mathbf{B}_{11}$	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	$(-x_3a - z_3c\cos\beta)\,\hat{\mathbf{x}} - y_3b\,\hat{\mathbf{y}} - z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O III
$\mathbf{B}_{12}$	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2}-y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_3a + z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O III
$\mathbf{B}_{13}$	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$(x_4a + z_4c\cos\beta)\,\hat{\mathbf{x}} + y_4b\,\hat{\mathbf{y}} + z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O IV
$\mathbf{B}_{14}$	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2}+y_4\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_4\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta - x_4a - z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O IV
$\mathbf{B}_{15}$	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	$(-x_4a - z_4c\cos\beta)\,\hat{\mathbf{x}} - y_4b\,\hat{\mathbf{y}} - z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O IV
$\mathbf{B}_{16}$	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2}-y_4\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_4a + z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O IV
$\mathbf{B}_{17}$	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	$(x_5a + z_5c\cos\beta)\,\hat{\mathbf{x}} + y_5b\,\hat{\mathbf{y}} + z_5c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O V

$\mathbf{B}_{18}$	$=$	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}c\cos\beta - x_5 a - z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O V
$\mathbf{B}_{19}$	$=$	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	$=$	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O V
$\mathbf{B}_{20}$	$=$	$x_5\,\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O V
$\mathbf{B}_{21}$	$=$	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	$=$	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O VI
$\mathbf{B}_{22}$	$=$	$-x_6\,\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}c\cos\beta - x_6 a - z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O VI
$\mathbf{B}_{23}$	$=$	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	$=$	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O VI
$\mathbf{B}_{24}$	$=$	$x_6\,\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O VI
$\mathbf{B}_{25}$	$=$	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	$=$	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	W I
$\mathbf{B}_{26}$	$=$	$-x_7\,\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}c\cos\beta - x_7 a - z_7 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	W I
$\mathbf{B}_{27}$	$=$	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	$=$	$(-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	W I
$\mathbf{B}_{28}$	$=$	$x_7\,\mathbf{a}_1 + \left(\frac{1}{2} - y_7\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}c\cos\beta + x_7 a + z_7 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_7\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	W I
$\mathbf{B}_{29}$	$=$	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	$=$	$(x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	W II
$\mathbf{B}_{30}$	$=$	$-x_8\,\mathbf{a}_1 + \left(\frac{1}{2} + y_8\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}c\cos\beta - x_8 a - z_8 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_8\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	W II
$\mathbf{B}_{31}$	$=$	$-x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	$=$	$(-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	W II
$\mathbf{B}_{32}$	$=$	$x_8\,\mathbf{a}_1 + \left(\frac{1}{2} - y_8\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}c\cos\beta + x_8 a + z_8 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_8\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	W II

References:

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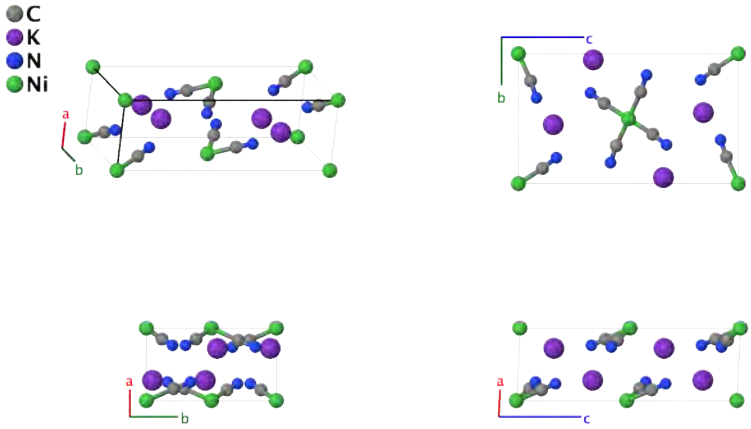
Geometry files:
- CIF: pp. 1542
- POSCAR: pp. 1542

$K_2Ni(CN)_4$ Structure: A4B2C4D_mP22_14_2e_e_2e_a

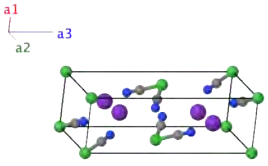
http://aflow.org/prototype-encyclopedia/A4B2C4D_mP22_14_2e_e_2e_a



Prototype	:	$C_4K_2N_4Ni$
AFLOW prototype label	:	A4B2C4D_mP22_14_2e_e_2e_a
***Strukturbericht* designation**	:	None
Pearson symbol	:	mP22
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=A4B2C4D_mP22_14_2e_e_2e_a --params=`$a, b/a, c/a, \beta, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6$

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	Ni
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Ni
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2a + z_2c\cos\beta)\,\hat{\mathbf{x}} + y_2b\,\hat{\mathbf{y}} + z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	C I
$\mathbf{B}_4$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_2a - z_2c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	C I
$\mathbf{B}_5$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2a - z_2c\cos\beta)\,\hat{\mathbf{x}} - y_2b\,\hat{\mathbf{y}} - z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	C I

$$\mathbf{B}_6 = x_2\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_2\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_2\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_2 a + z_2 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_2\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_2\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{C I}$$

$$\mathbf{B}_7 = x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3 = (x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{C II}$$

$$\mathbf{B}_8 = -x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_3 a - z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_3\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_3\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{C II}$$

$$\mathbf{B}_9 = -x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3 = (-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{C II}$$

$$\mathbf{B}_{10} = x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_3\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_3\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{C II}$$

$$\mathbf{B}_{11} = x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3 = (x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{K}$$

$$\mathbf{B}_{12} = -x_4\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_4 a - z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_4\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_4\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{K}$$

$$\mathbf{B}_{13} = -x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3 = (-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{K}$$

$$\mathbf{B}_{14} = x_4\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_4\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_4\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{K}$$

$$\mathbf{B}_{15} = x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3 = (x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{N I}$$

$$\mathbf{B}_{16} = -x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_5 a - z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_5\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_5\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{N I}$$

$$\mathbf{B}_{17} = -x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3 = (-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{N I}$$

$$\mathbf{B}_{18} = x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_5\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_5\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{N I}$$

$$\mathbf{B}_{19} = x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 = (x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{N II}$$

$$\mathbf{B}_{20} = -x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_6 a - z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_6\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_6\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{N II}$$

$$\mathbf{B}_{21} = -x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3 = (-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{N II}$$

$$\mathbf{B}_{22} = x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_6\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{N II}$$

References:

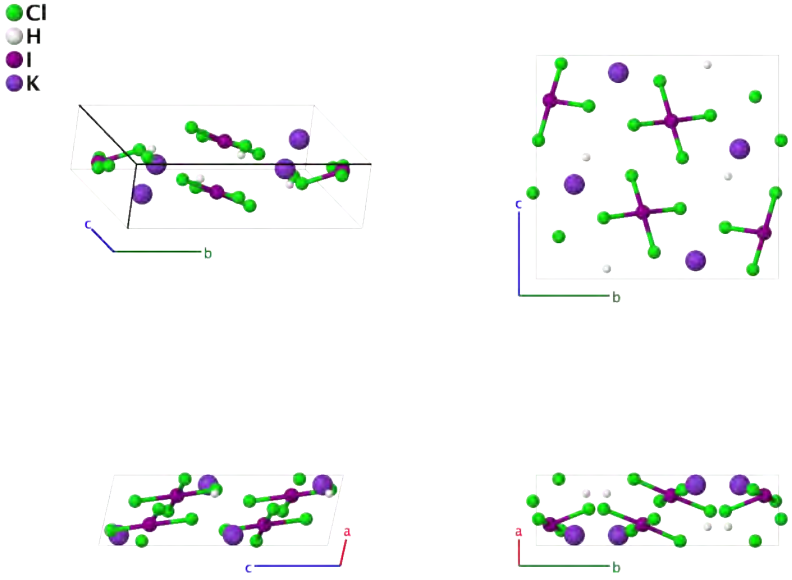
- N.-G. Vannerberg, *The Crystal Structure of $K_2Ni(CN)_4$*, Acta Chem. Scand. **18**, 2385–2391 (1964), doi:10.3891/acta.chem.scand.18-2385.

Geometry files:

- CIF: pp. 1542
- POSCAR: pp. 1543

$KICl_4{\cdot}H_2O$ ($H0_{10}$) Structure: A4BCD_mP28_14_4e_e_e_e

http://aflow.org/prototype-encyclopedia/A4BCD_mP28_14_4e_e_e_e

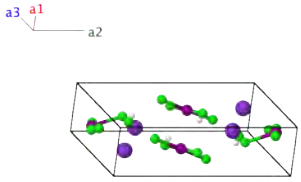


Prototype	:	$Cl_4IK \cdot H_2O$
AFLOW prototype label	:	A4BCD_mP28_14_4e_e_e_e
***Strukturbericht* designation**	:	$H0_{10}$
Pearson symbol	:	mP28
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=A4BCD_mP28_14_4e_e_e_e --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7$

- The structure of $KICl_4$ was originally determined by (Mooney, 1938) and assigned *Strukturbericht* designation $H0_{10}$ by (Herrmann, 1941). According to (Elema, 1963), "During Mooney's structure determination of $KIC1_4 \cdot H_2O$ it was not realized that the crystals contain water of crystallization." The correct structure is essentially the same as Mooney's with water molecules added into the vacancies, and we use $KICl_4 \cdot H_2O$ as our prototype for $H0_{10}$. To return to Mooney's structure, remove the water molecules, but of course this will not be a stable structure.
- The Wyckoff positions given by both (Mooney, 1938) and (Elema, 1963) are in the $P2_1/n$ setting of space group #14. We have used FINDSYM to translate this to the standard $P2_1/c$ setting.

Simple Monoclinic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= a\,\hat{\mathbf{x}}\\ \mathbf{a}_2 &= b\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	$(x_1a + z_1c\cos\beta)\,\hat{\mathbf{x}} + y_1b\,\hat{\mathbf{y}} + z_1c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cl I
$\mathbf{B}_2$	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2}+y_1\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_1\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta - x_1a - z_1c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cl I
$\mathbf{B}_3$	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	$(-x_1a - z_1c\cos\beta)\,\hat{\mathbf{x}} - y_1b\,\hat{\mathbf{y}} - z_1c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cl I
$\mathbf{B}_4$	$x_1\,\mathbf{a}_1 + \left(\frac{1}{2}-y_1\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_1\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_1a + z_1c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cl I
$\mathbf{B}_5$	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	$(x_2a + z_2c\cos\beta)\,\hat{\mathbf{x}} + y_2b\,\hat{\mathbf{y}} + z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cl II
$\mathbf{B}_6$	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2}+y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_2\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta - x_2a - z_2c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cl II
$\mathbf{B}_7$	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	$(-x_2a - z_2c\cos\beta)\,\hat{\mathbf{x}} - y_2b\,\hat{\mathbf{y}} - z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cl II
$\mathbf{B}_8$	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2}-y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_2a + z_2c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cl II
$\mathbf{B}_9$	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$(x_3a + z_3c\cos\beta)\,\hat{\mathbf{x}} + y_3b\,\hat{\mathbf{y}} + z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cl III
$\mathbf{B}_{10}$	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2}+y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_3\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta - x_3a - z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cl III
$\mathbf{B}_{11}$	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	$(-x_3a - z_3c\cos\beta)\,\hat{\mathbf{x}} - y_3b\,\hat{\mathbf{y}} - z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cl III
$\mathbf{B}_{12}$	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2}-y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_3a + z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cl III
$\mathbf{B}_{13}$	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$(x_4a + z_4c\cos\beta)\,\hat{\mathbf{x}} + y_4b\,\hat{\mathbf{y}} + z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cl IV
$\mathbf{B}_{14}$	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2}+y_4\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_4\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta - x_4a - z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cl IV
$\mathbf{B}_{15}$	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	$(-x_4a - z_4c\cos\beta)\,\hat{\mathbf{x}} - y_4b\,\hat{\mathbf{y}} - z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cl IV
$\mathbf{B}_{16}$	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2}-y_4\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_4a + z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cl IV

$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H_2O
$\mathbf{B}_{18}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2}+y_5\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_5 a - z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H_2O
$\mathbf{B}_{19}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H_2O
$\mathbf{B}_{20}$	=	$x_5\,\mathbf{a}_1 + \left(\frac{1}{2}-y_5\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H_2O
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	I
$\mathbf{B}_{22}$	=	$-x_6\,\mathbf{a}_1 + \left(\frac{1}{2}+y_6\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_6 a - z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_6\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	I
$\mathbf{B}_{23}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	I
$\mathbf{B}_{24}$	=	$x_6\,\mathbf{a}_1 + \left(\frac{1}{2}-y_6\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_6\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	I
$\mathbf{B}_{25}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	K
$\mathbf{B}_{26}$	=	$-x_7\,\mathbf{a}_1 + \left(\frac{1}{2}+y_7\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_7 a - z_7 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_7\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_7\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	K
$\mathbf{B}_{27}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	K
$\mathbf{B}_{28}$	=	$x_7\,\mathbf{a}_1 + \left(\frac{1}{2}-y_7\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_7 a + z_7 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_7\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_7\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	K

References:

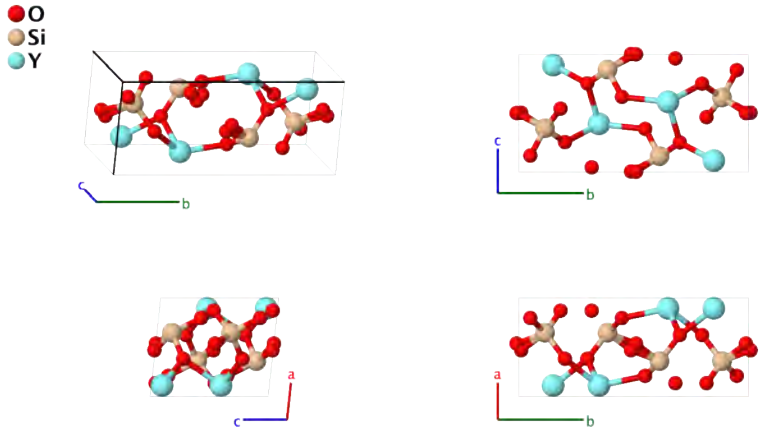
- R. J. Elema, J. L. de Boer, and A. Vos, *The refinement of the crystal structure of* $KICl_4 \cdot H_2O$, Acta Cryst. **16**, 243–247 (1963), doi:10.1107/S0365110X63000682.
- R. C. L. Mooney, *The Configuration of a Penthalogen Anion Group from the X-ray Structure Determination of Potassium Tetra-Chloriodide Crystals*, Zeitschrift für Kristallographie - Crystalline Materials **98**, 377–393 (1938), doi:10.1524/zkri.1938.98.1.377.
- K. Herrmann, ed., *Strukturbericht Band VI 1938* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1941).

Geometry files:

- CIF: pp. 1543
- POSCAR: pp. 1543

γ-$Y_2Si_2O_7$ Structure: A4BC_mP24_14_4e_e_e

http://aflow.org/prototype-encyclopedia/A4BC_mP24_14_4e_e_e



Prototype	:	$O_7Si_2Y_2$
AFLOW prototype label	:	A4BC_mP24_14_4e_e_e
***Strukturbericht* designation**	:	None
Pearson symbol	:	mP24
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=A4BC_mP24_14_4e_e_e --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6$

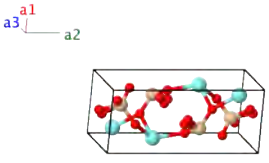
Other compounds with this structure

- γ-$Er_2Si_2O_7$ and γ-$Ho_2Si_2O_7$

- The O-III site is only 50% occupied. The Jmol image shows the two possible sites as a pair of very closely spaced oxygen atoms.
- (Christensen, 1997) refer to this as D-$Y_2Si_2O_7$, but we follow the classification of the $RE_2Si_2O_7$ structures in (Becerro, 2004) and refer to this as γ-$Y_2Si_2O_7$.

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$$
\begin{array}{lcccccc}
\mathbf{B}_1 & = & x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3 & = & (x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{O I} \\
\mathbf{B}_2 & = & -x_1\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_1\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_1\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}c\cos\beta - x_1 a - z_1 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_1\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_1\right) c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{O I} \\
\mathbf{B}_3 & = & -x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3 & = & (-x_1 a - z_1 c\cos\beta)\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} - z_1 c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{O I} \\
\mathbf{B}_4 & = & x_1\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_1\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_1\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}c\cos\beta + x_1 a + z_1 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_1\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_1\right) c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{O I} \\
\mathbf{B}_5 & = & x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3 & = & (x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{O II} \\
\mathbf{B}_6 & = & -x_2\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_2\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_2\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}c\cos\beta - x_2 a - z_2 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_2\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_2\right) c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{O II} \\
\mathbf{B}_7 & = & -x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3 & = & (-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} - z_2 c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{O II} \\
\mathbf{B}_8 & = & x_2\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_2\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_2\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}c\cos\beta + x_2 a + z_2 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_2\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_2\right) c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{O II} \\
\mathbf{B}_9 & = & x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3 & = & (x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{O III} \\
\mathbf{B}_{10} & = & -x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_3\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}c\cos\beta - x_3 a - z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_3\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_3\right) c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{O III} \\
\mathbf{B}_{11} & = & -x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3 & = & (-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{O III} \\
\mathbf{B}_{12} & = & x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_3\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_3\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_3\right) c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{O III} \\
\mathbf{B}_{13} & = & x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3 & = & (x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{O IV} \\
\mathbf{B}_{14} & = & -x_4\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_4\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}c\cos\beta - x_4 a - z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_4\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_4\right) c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{O IV} \\
\mathbf{B}_{15} & = & -x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3 & = & (-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{O IV} \\
\mathbf{B}_{16} & = & x_4\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_4\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_4\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_4\right) c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{O IV} \\
\mathbf{B}_{17} & = & x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3 & = & (x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{Si} \\
\mathbf{B}_{18} & = & -x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_5\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}c\cos\beta - x_5 a - z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_5\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_5\right) c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{Si} \\
\mathbf{B}_{19} & = & -x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3 & = & (-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{Si} \\
\mathbf{B}_{20} & = & x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_5\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_5\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_5\right) c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{Si} \\
\mathbf{B}_{21} & = & x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 & = & (x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{Y} \\
\mathbf{B}_{22} & = & -x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_6\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}c\cos\beta - x_6 a - z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_6\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_6\right) c\sin\beta\,\hat{\mathbf{z}} & (4e) & \text{Y}
\end{array}
$$

$\mathbf{B}_{23}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Y
$\mathbf{B}_{24}$	=	$x_6\,\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Y

References:

- A. N. Christensen, R. G. Hazell, and A. W. Hewat, *Synthesis, Crystal Growth and Structure Investigations of Rare-Earth Disilicates and Rare-Earth Oxyapatites*, Acta Chem. Scand. **51**, 37–43 (1997), doi:10.3891/acta.chem.scand.51-0037.

Found in:

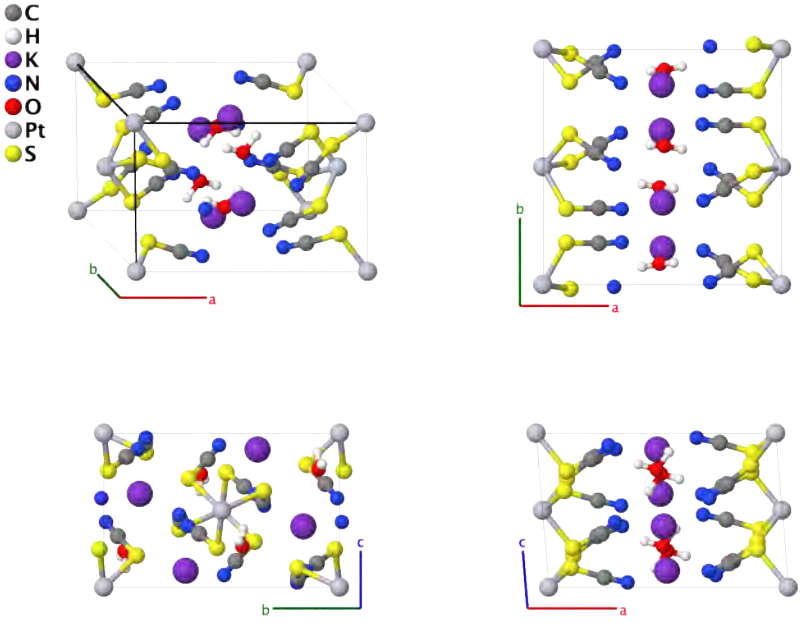
- A. I. Becerro and A. Escudero, *Revision of the crystallographic data of polymorphic $Y_2Si_2O_7$ and Y_2SiO_5 compounds*, Phase Transit. **77**, 1093–1102 (2004), doi:10.1080/01411590412331282814.

Geometry files:

- CIF: pp. 1543
- POSCAR: pp. 1544

$K_2Pt(SCN)_6 \cdot 2H_2O$ Structure: A6B4C2D6E2FG6_mP54_14_3e_2e_e_3e_e_a_3e

http://aflow.org/prototype-encyclopedia/A6B4C2D6E2FG6_mP54_14_3e_2e_e_3e_e_a_3e



Prototype	:	$C_6H_4K_2N_6O_2PtS_6$
AFLOW prototype label	:	A6B4C2D6E2FG6_mP54_14_3e_2e_e_3e_e_a_3e
***Strukturbericht* designation**	:	None
Pearson symbol	:	mP54
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=A6B4C2D6E2FG6_mP54_14_3e_2e_e_3e_e_a_3e --params=`$a, b/a, c/a, \beta, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}$

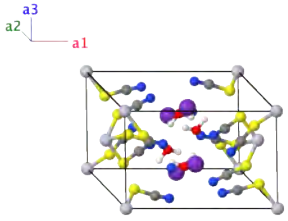
Other compounds with this structure

- $Rb_2Pt(SCN)_6 \cdot 2H_2O$ and $NH_{42}Pt(SCN)_6 \cdot 2H_2O$

- This is the hydrated form. For the anhydrous structure, see the $H6_3$ structure page.

Simple Monoclinic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= a\,\hat{\mathbf{x}}\\ \mathbf{a}_2 &= b\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	Pt
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Pt
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2a + z_2c\cos\beta)\,\hat{\mathbf{x}} + y_2b\,\hat{\mathbf{y}} + z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	C I
$\mathbf{B}_4$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2}+y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_2a - z_2c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	C I
$\mathbf{B}_5$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2a - z_2c\cos\beta)\,\hat{\mathbf{x}} - y_2b\,\hat{\mathbf{y}} - z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	C I
$\mathbf{B}_6$	=	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2}-y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_2a + z_2c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	C I
$\mathbf{B}_7$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3a + z_3c\cos\beta)\,\hat{\mathbf{x}} + y_3b\,\hat{\mathbf{y}} + z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	C II
$\mathbf{B}_8$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2}+y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_3a - z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	C II
$\mathbf{B}_9$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3a - z_3c\cos\beta)\,\hat{\mathbf{x}} - y_3b\,\hat{\mathbf{y}} - z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	C II
$\mathbf{B}_{10}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2}-y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_3a + z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	C II
$\mathbf{B}_{11}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4a + z_4c\cos\beta)\,\hat{\mathbf{x}} + y_4b\,\hat{\mathbf{y}} + z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	C III
$\mathbf{B}_{12}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2}+y_4\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_4a - z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	C III
$\mathbf{B}_{13}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4a - z_4c\cos\beta)\,\hat{\mathbf{x}} - y_4b\,\hat{\mathbf{y}} - z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	C III
$\mathbf{B}_{14}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2}-y_4\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_4a + z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	C III
$\mathbf{B}_{15}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5a + z_5c\cos\beta)\,\hat{\mathbf{x}} + y_5b\,\hat{\mathbf{y}} + z_5c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H I
$\mathbf{B}_{16}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2}+y_5\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_5a - z_5c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H I
$\mathbf{B}_{17}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5a - z_5c\cos\beta)\,\hat{\mathbf{x}} - y_5b\,\hat{\mathbf{y}} - z_5c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H I

$$\mathbf{B}_{18} = x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_5\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_5\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H I}$$

$$\mathbf{B}_{19} = x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 = (x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H II}$$

$$\mathbf{B}_{20} = -x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_6 a - z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_6\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_6\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H II}$$

$$\mathbf{B}_{21} = -x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3 = (-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H II}$$

$$\mathbf{B}_{22} = x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_6\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H II}$$

$$\mathbf{B}_{23} = x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 = (x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{K}$$

$$\mathbf{B}_{24} = -x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_7 a - z_7 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_7\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_7\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{K}$$

$$\mathbf{B}_{25} = -x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3 = (-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{K}$$

$$\mathbf{B}_{26} = x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_7 a + z_7 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_7\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_7\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{K}$$

$$\mathbf{B}_{27} = x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 = (x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{N I}$$

$$\mathbf{B}_{28} = -x_8\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_8 a - z_8 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_8\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_8\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{N I}$$

$$\mathbf{B}_{29} = -x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3 = (-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{N I}$$

$$\mathbf{B}_{30} = x_8\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_8 a + z_8 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_8\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_8\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{N I}$$

$$\mathbf{B}_{31} = x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 = (x_9 a + z_9 c\cos\beta)\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{N II}$$

$$\mathbf{B}_{32} = -x_9\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_9 a - z_9 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_9\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_9\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{N II}$$

$$\mathbf{B}_{33} = -x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3 = (-x_9 a - z_9 c\cos\beta)\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} - z_9 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{N II}$$

$$\mathbf{B}_{34} = x_9\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_9 a + z_9 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_9\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_9\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{N II}$$

$$\mathbf{B}_{35} = x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 = (x_{10} a + z_{10} c\cos\beta)\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{N III}$$

$$\mathbf{B}_{36} = -x_{10}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{10}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{10} a - z_{10} c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{10}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{10}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{N III}$$

$$\mathbf{B}_{37} = -x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 = (-x_{10} a - z_{10} c\cos\beta)\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} - z_{10} c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{N III}$$

$$\mathbf{B}_{38} = x_{10}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{10}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{10} a + z_{10} c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{10}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{10}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{N III}$$

$$\mathbf{B}_{39} = x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 = (x_{11} a + z_{11} c\cos\beta)\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + z_{11} c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O}$$

$\mathbf{B}_{40}$	=	$-x_{11}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{11}\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{11}\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_{11}a - z_{11}c\cos\beta\right)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{11}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{11}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O
$\mathbf{B}_{41}$	=	$-x_{11}\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3$	=	$(-x_{11}a - z_{11}c\cos\beta)\,\hat{\mathbf{x}} - y_{11}b\,\hat{\mathbf{y}} - z_{11}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O
$\mathbf{B}_{42}$	=	$x_{11}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{11}\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{11}\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{11}a + z_{11}c\cos\beta\right)\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{11}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{11}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O
$\mathbf{B}_{43}$	=	$x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3$	=	$(x_{12}a + z_{12}c\cos\beta)\,\hat{\mathbf{x}} + y_{12}b\,\hat{\mathbf{y}} + z_{12}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S I
$\mathbf{B}_{44}$	=	$-x_{12}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{12}\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{12}\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_{12}a - z_{12}c\cos\beta\right)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{12}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{12}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S I
$\mathbf{B}_{45}$	=	$-x_{12}\,\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3$	=	$(-x_{12}a - z_{12}c\cos\beta)\,\hat{\mathbf{x}} - y_{12}b\,\hat{\mathbf{y}} - z_{12}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S I
$\mathbf{B}_{46}$	=	$x_{12}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{12}\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{12}\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{12}a + z_{12}c\cos\beta\right)\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{12}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{12}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S I
$\mathbf{B}_{47}$	=	$x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3$	=	$(x_{13}a + z_{13}c\cos\beta)\,\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} + z_{13}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S II
$\mathbf{B}_{48}$	=	$-x_{13}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{13}\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{13}\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_{13}a - z_{13}c\cos\beta\right)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{13}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{13}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S II
$\mathbf{B}_{49}$	=	$-x_{13}\,\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 - z_{13}\,\mathbf{a}_3$	=	$(-x_{13}a - z_{13}c\cos\beta)\,\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} - z_{13}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S II
$\mathbf{B}_{50}$	=	$x_{13}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{13}\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{13}\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{13}a + z_{13}c\cos\beta\right)\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{13}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{13}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S II
$\mathbf{B}_{51}$	=	$x_{14}\,\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3$	=	$(x_{14}a + z_{14}c\cos\beta)\,\hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} + z_{14}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S III
$\mathbf{B}_{52}$	=	$-x_{14}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{14}\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{14}\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_{14}a - z_{14}c\cos\beta\right)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{14}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{14}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S III
$\mathbf{B}_{53}$	=	$-x_{14}\,\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 - z_{14}\,\mathbf{a}_3$	=	$(-x_{14}a - z_{14}c\cos\beta)\,\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} - z_{14}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S III
$\mathbf{B}_{54}$	=	$x_{14}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{14}\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{14}\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{14}a + z_{14}c\cos\beta\right)\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{14}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{14}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S III

References:

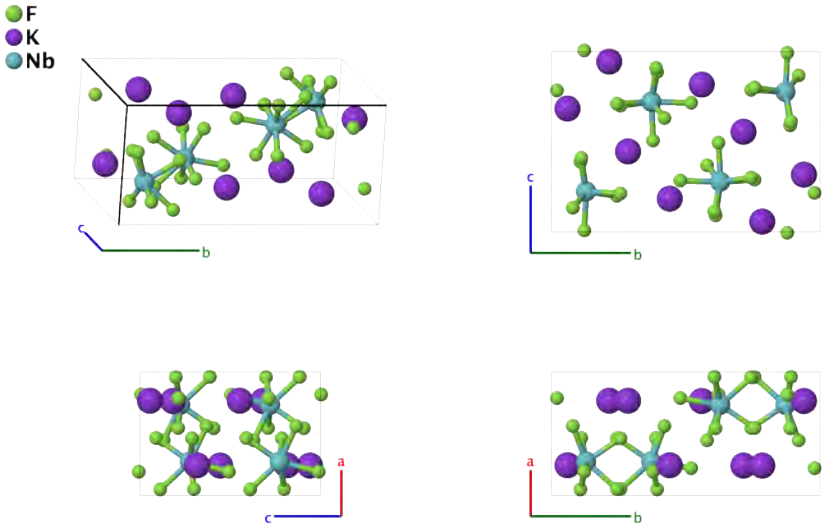
- J. Arpalahti, J. Hölsä, and R. Sillanpää, *Studies on Potassium Thiocyanatoplatinates. II. Crystal Structure of Potassium Hexathiocyanatoplatinate(IV) Dihydrate, $K_2Pt(SCN)_6 \cdot 2H_2O$*, Acta Chem. Scand. **47**, 1078–1082 (1993), doi:10.3891/acta.chem.scand.47-1078.

Geometry files:

- CIF: pp. 1544
- POSCAR: pp. 1544

K_2NbF_7 ($K6_2$) Structure: A7B2C_mP40_14_7e_2e_e

http://aflow.org/prototype-encyclopedia/A7B2C_mP40_14_7e_2e_e



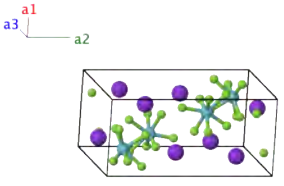
Prototype	:	F_7K_2Nb
AFLOW prototype label	:	A7B2C_mP40_14_7e_2e_e
***Strukturbericht* designation**	:	$K6_2$
Pearson symbol	:	mP40
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=A7B2C_mP40_14_7e_2e_e --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}$

Other compounds with this structure

- K_2TaF_7

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	F I

$$\mathbf{B}_2 = -x_1\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_1\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_1\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_1 a - z_1 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_1\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_1\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{F I}$$

$$\mathbf{B}_3 = -x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3 = (-x_1 a - z_1 c\cos\beta)\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} - z_1 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{F I}$$

$$\mathbf{B}_4 = x_1\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_1\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_1\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_1 a + z_1 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_1\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_1\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{F I}$$

$$\mathbf{B}_5 = x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3 = (x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{F II}$$

$$\mathbf{B}_6 = -x_2\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_2\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_2\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_2 a - z_2 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_2\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_2\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{F II}$$

$$\mathbf{B}_7 = -x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3 = (-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} - z_2 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{F II}$$

$$\mathbf{B}_8 = x_2\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_2\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_2\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_2 a + z_2 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_2\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_2\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{F II}$$

$$\mathbf{B}_9 = x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3 = (x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{F III}$$

$$\mathbf{B}_{10} = -x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_3 a - z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_3\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_3\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{F III}$$

$$\mathbf{B}_{11} = -x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3 = (-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{F III}$$

$$\mathbf{B}_{12} = x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_3\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_3\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{F III}$$

$$\mathbf{B}_{13} = x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3 = (x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{F IV}$$

$$\mathbf{B}_{14} = -x_4\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_4 a - z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_4\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_4\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{F IV}$$

$$\mathbf{B}_{15} = -x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3 = (-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{F IV}$$

$$\mathbf{B}_{16} = x_4\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_4\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_4\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{F IV}$$

$$\mathbf{B}_{17} = x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3 = (x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{F V}$$

$$\mathbf{B}_{18} = -x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_5 a - z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_5\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_5\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{F V}$$

$$\mathbf{B}_{19} = -x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3 = (-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{F V}$$

$$\mathbf{B}_{20} = x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_5\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_5\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{F V}$$

$$\mathbf{B}_{21} = x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 = (x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{F VI}$$

$$\mathbf{B}_{22} = -x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_6 a - z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_6\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_6\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{F VI}$$

$$\mathbf{B}_{23} = -x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3 = (-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{F VI}$$

$\mathbf{B}_{24}$	=	$x_6\,\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	F VI
$\mathbf{B}_{25}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	F VII
$\mathbf{B}_{26}$	=	$-x_7\,\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_7 a - z_7 c\cos\beta\right)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	F VII
$\mathbf{B}_{27}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	F VII
$\mathbf{B}_{28}$	=	$x_7\,\mathbf{a}_1 + \left(\frac{1}{2} - y_7\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_7 a + z_7 c\cos\beta\right)\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_7\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	F VII
$\mathbf{B}_{29}$	=	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	K I
$\mathbf{B}_{30}$	=	$-x_8\,\mathbf{a}_1 + \left(\frac{1}{2} + y_8\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_8 a - z_8 c\cos\beta\right)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_8\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	K I
$\mathbf{B}_{31}$	=	$-x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	K I
$\mathbf{B}_{32}$	=	$x_8\,\mathbf{a}_1 + \left(\frac{1}{2} - y_8\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_8 a + z_8 c\cos\beta\right)\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_8\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	K I
$\mathbf{B}_{33}$	=	$x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$(x_9 a + z_9 c\cos\beta)\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	K II
$\mathbf{B}_{34}$	=	$-x_9\,\mathbf{a}_1 + \left(\frac{1}{2} + y_9\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_9\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_9 a - z_9 c\cos\beta\right)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_9\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_9\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	K II
$\mathbf{B}_{35}$	=	$-x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$(-x_9 a - z_9 c\cos\beta)\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} - z_9 c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	K II
$\mathbf{B}_{36}$	=	$x_9\,\mathbf{a}_1 + \left(\frac{1}{2} - y_9\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_9 a + z_9 c\cos\beta\right)\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_9\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	K II
$\mathbf{B}_{37}$	=	$x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$(x_{10} a + z_{10} c\cos\beta)\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Nb
$\mathbf{B}_{38}$	=	$-x_{10}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{10}\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{10}\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_{10} a - z_{10} c\cos\beta\right)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{10}\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{10}\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Nb
$\mathbf{B}_{39}$	=	$-x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3$	=	$(-x_{10} a - z_{10} c\cos\beta)\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} - z_{10} c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Nb
$\mathbf{B}_{40}$	=	$x_{10}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{10}\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{10}\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{10} a + z_{10} c\cos\beta\right)\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{10}\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{10}\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Nb

References:

- G. M. Brown and L. A. Walker, *Refinement of the structure of potassium heptafluoroniobate, K_2NbF_7, from neutron-diffraction data*, Acta Cryst. **20**, 220–229 (1966), doi:10.1107/S0365110X66000458.

Found in:

- C. C. Torardi, L. H. Brixner, and G. Blasse, *Structure and luminescence of K_2TaF_7 and K_2NbF_7*, J. Solid State Chem. **67**, 21–25 (1987), doi:10.1016/0022-4596(87)90333-1.

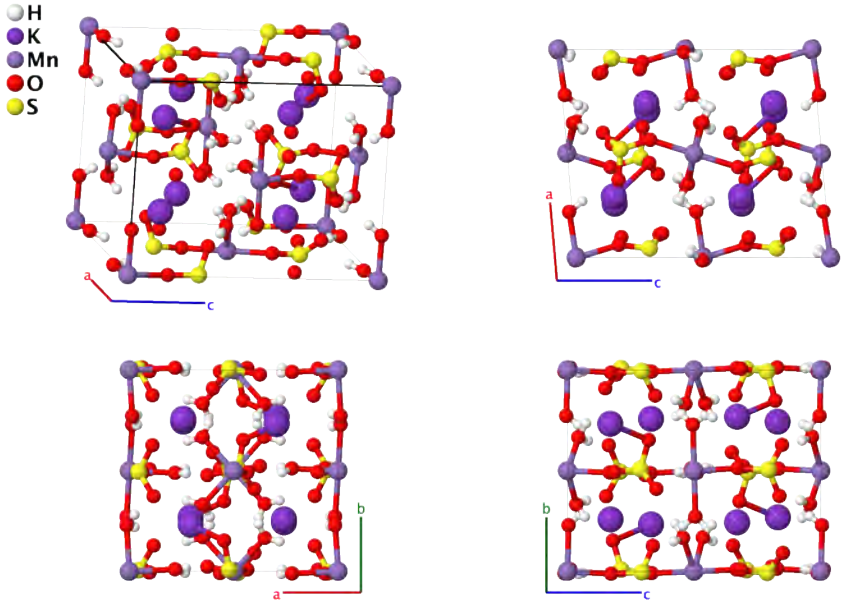
Geometry files:

- CIF: pp. 1545

- POSCAR: pp. 1545

Manganese-leonite 110 K [$K_2Mn(SO_4)_2{\cdot}4H_2O$] Structure: A8B2CD12E2_mP100_14_8e_2e_ad_12e_2e

http://aflow.org/prototype-encyclopedia/A8B2CD12E2_mP100_14_8e_2e_ad_12e_2e



Prototype	:	$H_8K_2MnO_{12}S_2$
AFLOW prototype label	:	A8B2CD12E2_mP100_14_8e_2e_ad_12e_2e
***Strukturbericht* designation**	:	None
Pearson symbol	:	mP100
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=A8B2CD12E2_mP100_14_8e_2e_ad_12e_2e --params=`$a, b/a, c/a, \beta, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}, x_{17}, y_{17}, z_{17}, x_{18}, y_{18}, z_{18}, x_{19}, y_{19}, z_{19}, x_{20}, y_{20}, z_{20}, x_{21}, y_{21}, z_{21}, x_{22}, y_{22}, z_{22}, x_{23}, y_{23}, z_{23}, x_{24}, y_{24}, z_{24}, x_{25}, y_{25}, z_{25}, x_{26}, y_{26}, z_{26}$

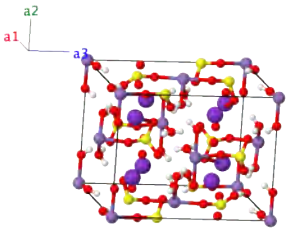
Other compounds with this structure

- $K_2Mg(SO_4)_2{\cdot}4H_2O$ (leonite)

- This is the intermediate-low temperature structure of leonite. For Mn-leonite, discussed here, it is stable below 168 K. The structure shown here was determined at 110 K.
- The room temperature structure is *Strukturbericht* $H4_{23}$, space group $C2/m$ #13. The current structure doubles the size of the unit cell and orders all of the SO_4 radicals.
- The intermediate temperature structure (168-205 K) has space group $P2_1/c$ #14.
- (Hertweck, 2001) give crystallographic information of this phase in the $P2_1/a$ setting of space group #14. We used FINDSYM to transform this to the $P2_1/c$ setting, which resulted in a switching of the a- and c-axes.

Simple Monoclinic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= a\,\hat{\mathbf{x}}\\ \mathbf{a}_2 &= b\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	Mn I
$\mathbf{B}_2$	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Mn I
$\mathbf{B}_3$	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	$\frac{1}{2}\,(a + c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	$(2d)$	Mn II
$\mathbf{B}_4$	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$	$(2d)$	Mn II
$\mathbf{B}_5$	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$(x_3a + z_3c\cos\beta)\,\hat{\mathbf{x}} + y_3b\,\hat{\mathbf{y}} + z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H I
$\mathbf{B}_6$	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta - x_3a - z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H I
$\mathbf{B}_7$	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	$(-x_3a - z_3c\cos\beta)\,\hat{\mathbf{x}} - y_3b\,\hat{\mathbf{y}} - z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H I
$\mathbf{B}_8$	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_3a + z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H I
$\mathbf{B}_9$	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$(x_4a + z_4c\cos\beta)\,\hat{\mathbf{x}} + y_4b\,\hat{\mathbf{y}} + z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H II
$\mathbf{B}_{10}$	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta - x_4a - z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H II
$\mathbf{B}_{11}$	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	$(-x_4a - z_4c\cos\beta)\,\hat{\mathbf{x}} - y_4b\,\hat{\mathbf{y}} - z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H II
$\mathbf{B}_{12}$	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_4a + z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H II
$\mathbf{B}_{13}$	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	$(x_5a + z_5c\cos\beta)\,\hat{\mathbf{x}} + y_5b\,\hat{\mathbf{y}} + z_5c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H III
$\mathbf{B}_{14}$	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta - x_5a - z_5c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H III
$\mathbf{B}_{15}$	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	$(-x_5a - z_5c\cos\beta)\,\hat{\mathbf{x}} - y_5b\,\hat{\mathbf{y}} - z_5c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H III
$\mathbf{B}_{16}$	$x_5\,\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_5a + z_5c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H III
$\mathbf{B}_{17}$	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	$(x_6a + z_6c\cos\beta)\,\hat{\mathbf{x}} + y_6b\,\hat{\mathbf{y}} + z_6c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H IV
$\mathbf{B}_{18}$	$-x_6\,\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta - x_6a - z_6c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H IV

$$
\begin{array}{lcccll}
\mathbf{B}_{19} & = & -x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3 & = & (-x_6 a - z_6 c\cos\beta)\;\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}} & (4e)\quad \text{H IV} \\
\mathbf{B}_{20} & = & x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_6\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right) c\sin\beta\,\hat{\mathbf{z}} & (4e)\quad \text{H IV} \\
\mathbf{B}_{21} & = & x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 & = & (x_7 a + z_7 c\cos\beta)\;\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}} & (4e)\quad \text{H V} \\
\mathbf{B}_{22} & = & -x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_7\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}c\cos\beta - x_7 a - z_7 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_7\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_7\right) c\sin\beta\,\hat{\mathbf{z}} & (4e)\quad \text{H V} \\
\mathbf{B}_{23} & = & -x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3 & = & (-x_7 a - z_7 c\cos\beta)\;\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}} & (4e)\quad \text{H V} \\
\mathbf{B}_{24} & = & x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_7\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}c\cos\beta + x_7 a + z_7 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_7\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_7\right) c\sin\beta\,\hat{\mathbf{z}} & (4e)\quad \text{H V} \\
\mathbf{B}_{25} & = & x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 & = & (x_8 a + z_8 c\cos\beta)\;\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}} & (4e)\quad \text{H VI} \\
\mathbf{B}_{26} & = & -x_8\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_8\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}c\cos\beta - x_8 a - z_8 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_8\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_8\right) c\sin\beta\,\hat{\mathbf{z}} & (4e)\quad \text{H VI} \\
\mathbf{B}_{27} & = & -x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3 & = & (-x_8 a - z_8 c\cos\beta)\;\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}} & (4e)\quad \text{H VI} \\
\mathbf{B}_{28} & = & x_8\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_8\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}c\cos\beta + x_8 a + z_8 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_8\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_8\right) c\sin\beta\,\hat{\mathbf{z}} & (4e)\quad \text{H VI} \\
\mathbf{B}_{29} & = & x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 & = & (x_9 a + z_9 c\cos\beta)\;\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\sin\beta\,\hat{\mathbf{z}} & (4e)\quad \text{H VII} \\
\mathbf{B}_{30} & = & -x_9\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_9\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}c\cos\beta - x_9 a - z_9 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_9\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_9\right) c\sin\beta\,\hat{\mathbf{z}} & (4e)\quad \text{H VII} \\
\mathbf{B}_{31} & = & -x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3 & = & (-x_9 a - z_9 c\cos\beta)\;\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} - z_9 c\sin\beta\,\hat{\mathbf{z}} & (4e)\quad \text{H VII} \\
\mathbf{B}_{32} & = & x_9\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}c\cos\beta + x_9 a + z_9 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_9\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_9\right) c\sin\beta\,\hat{\mathbf{z}} & (4e)\quad \text{H VII} \\
\mathbf{B}_{33} & = & x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 & = & (x_{10} a + z_{10} c\cos\beta)\;\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\sin\beta\,\hat{\mathbf{z}} & (4e)\quad \text{H VIII} \\
\mathbf{B}_{34} & = & -x_{10}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{10}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{10}\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}c\cos\beta - x_{10} a - z_{10} c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{10}\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{10}\right) c\sin\beta\,\hat{\mathbf{z}} & (4e)\quad \text{H VIII} \\
\mathbf{B}_{35} & = & -x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 & = & (-x_{10} a - z_{10} c\cos\beta)\;\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} - z_{10} c\sin\beta\,\hat{\mathbf{z}} & (4e)\quad \text{H VIII} \\
\mathbf{B}_{36} & = & x_{10}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{10}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{10}\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}c\cos\beta + x_{10} a + z_{10} c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{10}\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{10}\right) c\sin\beta\,\hat{\mathbf{z}} & (4e)\quad \text{H VIII} \\
\mathbf{B}_{37} & = & x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 & = & (x_{11} a + z_{11} c\cos\beta)\;\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + z_{11} c\sin\beta\,\hat{\mathbf{z}} & (4e)\quad \text{K I} \\
\mathbf{B}_{38} & = & -x_{11}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{11}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{11}\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}c\cos\beta - x_{11} a - z_{11} c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{11}\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{11}\right) c\sin\beta\,\hat{\mathbf{z}} & (4e)\quad \text{K I} \\
\mathbf{B}_{39} & = & -x_{11}\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3 & = & (-x_{11} a - z_{11} c\cos\beta)\;\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} - z_{11} c\sin\beta\,\hat{\mathbf{z}} & (4e)\quad \text{K I} \\
\mathbf{B}_{40} & = & x_{11}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{11}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{11}\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}c\cos\beta + x_{11} a + z_{11} c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{11}\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{11}\right) c\sin\beta\,\hat{\mathbf{z}} & (4e)\quad \text{K I}
\end{array}
$$

$$\mathbf{B}_{41} = x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 = (x_{12}a + z_{12}c\cos\beta)\,\hat{\mathbf{x}} + y_{12}b\,\hat{\mathbf{y}} + z_{12}c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{K II}$$

$$\mathbf{B}_{42} = -x_{12}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{12}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{12}a - z_{12}c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{12}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{12}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{K II}$$

$$\mathbf{B}_{43} = -x_{12}\,\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3 = (-x_{12}a - z_{12}c\cos\beta)\,\hat{\mathbf{x}} - y_{12}b\,\hat{\mathbf{y}} - z_{12}c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{K II}$$

$$\mathbf{B}_{44} = x_{12}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{12}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{12}a + z_{12}c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{12}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{12}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{K II}$$

$$\mathbf{B}_{45} = x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3 = (x_{13}a + z_{13}c\cos\beta)\,\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} + z_{13}c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O I}$$

$$\mathbf{B}_{46} = -x_{13}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{13}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{13}a - z_{13}c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{13}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{13}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O I}$$

$$\mathbf{B}_{47} = -x_{13}\,\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 - z_{13}\,\mathbf{a}_3 = (-x_{13}a - z_{13}c\cos\beta)\,\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} - z_{13}c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O I}$$

$$\mathbf{B}_{48} = x_{13}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{13}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{13}a + z_{13}c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{13}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{13}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O I}$$

$$\mathbf{B}_{49} = x_{14}\,\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3 = (x_{14}a + z_{14}c\cos\beta)\,\hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} + z_{14}c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O II}$$

$$\mathbf{B}_{50} = -x_{14}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{14}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{14}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{14}a - z_{14}c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{14}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{14}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O II}$$

$$\mathbf{B}_{51} = -x_{14}\,\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 - z_{14}\,\mathbf{a}_3 = (-x_{14}a - z_{14}c\cos\beta)\,\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} - z_{14}c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O II}$$

$$\mathbf{B}_{52} = x_{14}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{14}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{14}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{14}a + z_{14}c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{14}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{14}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O II}$$

$$\mathbf{B}_{53} = x_{15}\,\mathbf{a}_1 + y_{15}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 = (x_{15}a + z_{15}c\cos\beta)\,\hat{\mathbf{x}} + y_{15}b\,\hat{\mathbf{y}} + z_{15}c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O III}$$

$$\mathbf{B}_{54} = -x_{15}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{15}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{15}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{15}a - z_{15}c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{15}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{15}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O III}$$

$$\mathbf{B}_{55} = -x_{15}\,\mathbf{a}_1 - y_{15}\,\mathbf{a}_2 - z_{15}\,\mathbf{a}_3 = (-x_{15}a - z_{15}c\cos\beta)\,\hat{\mathbf{x}} - y_{15}b\,\hat{\mathbf{y}} - z_{15}c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O III}$$

$$\mathbf{B}_{56} = x_{15}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{15}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{15}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{15}a + z_{15}c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{15}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{15}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O III}$$

$$\mathbf{B}_{57} = x_{16}\,\mathbf{a}_1 + y_{16}\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 = (x_{16}a + z_{16}c\cos\beta)\,\hat{\mathbf{x}} + y_{16}b\,\hat{\mathbf{y}} + z_{16}c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O IV}$$

$$\mathbf{B}_{58} = -x_{16}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{16}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{16}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{16}a - z_{16}c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{16}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{16}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O IV}$$

$$\mathbf{B}_{59} = -x_{16}\,\mathbf{a}_1 - y_{16}\,\mathbf{a}_2 - z_{16}\,\mathbf{a}_3 = (-x_{16}a - z_{16}c\cos\beta)\,\hat{\mathbf{x}} - y_{16}b\,\hat{\mathbf{y}} - z_{16}c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O IV}$$

$$\mathbf{B}_{60} = x_{16}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{16}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{16}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{16}a + z_{16}c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{16}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{16}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O IV}$$

$$\mathbf{B}_{61} = x_{17}\,\mathbf{a}_1 + y_{17}\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 = (x_{17}a + z_{17}c\cos\beta)\,\hat{\mathbf{x}} + y_{17}b\,\hat{\mathbf{y}} + z_{17}c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O V}$$

$$\mathbf{B}_{62} = -x_{17}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{17}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{17}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{17}a - z_{17}c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{17}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{17}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O V}$$

$\mathbf{B}_{63}$	=	$-x_{17}\,\mathbf{a}_1 - y_{17}\,\mathbf{a}_2 - z_{17}\,\mathbf{a}_3$	=	$(-x_{17}a - z_{17}c\cos\beta)\;\hat{\mathbf{x}} - y_{17}b\,\hat{\mathbf{y}} - z_{17}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O V
$\mathbf{B}_{64}$	=	$x_{17}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{17}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{17}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{17}a + z_{17}c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{17}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{17}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O V
$\mathbf{B}_{65}$	=	$x_{18}\,\mathbf{a}_1 + y_{18}\,\mathbf{a}_2 + z_{18}\,\mathbf{a}_3$	=	$(x_{18}a + z_{18}c\cos\beta)\;\hat{\mathbf{x}} + y_{18}b\,\hat{\mathbf{y}} + z_{18}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O VI
$\mathbf{B}_{66}$	=	$-x_{18}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{18}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{18}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_{18}a - z_{18}c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{18}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{18}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O VI
$\mathbf{B}_{67}$	=	$-x_{18}\,\mathbf{a}_1 - y_{18}\,\mathbf{a}_2 - z_{18}\,\mathbf{a}_3$	=	$(-x_{18}a - z_{18}c\cos\beta)\;\hat{\mathbf{x}} - y_{18}b\,\hat{\mathbf{y}} - z_{18}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O VI
$\mathbf{B}_{68}$	=	$x_{18}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{18}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{18}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{18}a + z_{18}c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{18}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{18}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O VI
$\mathbf{B}_{69}$	=	$x_{19}\,\mathbf{a}_1 + y_{19}\,\mathbf{a}_2 + z_{19}\,\mathbf{a}_3$	=	$(x_{19}a + z_{19}c\cos\beta)\;\hat{\mathbf{x}} + y_{19}b\,\hat{\mathbf{y}} + z_{19}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O VII
$\mathbf{B}_{70}$	=	$-x_{19}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{19}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{19}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_{19}a - z_{19}c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{19}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{19}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O VII
$\mathbf{B}_{71}$	=	$-x_{19}\,\mathbf{a}_1 - y_{19}\,\mathbf{a}_2 - z_{19}\,\mathbf{a}_3$	=	$(-x_{19}a - z_{19}c\cos\beta)\;\hat{\mathbf{x}} - y_{19}b\,\hat{\mathbf{y}} - z_{19}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O VII
$\mathbf{B}_{72}$	=	$x_{19}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{19}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{19}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{19}a + z_{19}c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{19}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{19}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O VII
$\mathbf{B}_{73}$	=	$x_{20}\,\mathbf{a}_1 + y_{20}\,\mathbf{a}_2 + z_{20}\,\mathbf{a}_3$	=	$(x_{20}a + z_{20}c\cos\beta)\;\hat{\mathbf{x}} + y_{20}b\,\hat{\mathbf{y}} + z_{20}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O VIII
$\mathbf{B}_{74}$	=	$-x_{20}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{20}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{20}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_{20}a - z_{20}c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{20}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{20}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O VIII
$\mathbf{B}_{75}$	=	$-x_{20}\,\mathbf{a}_1 - y_{20}\,\mathbf{a}_2 - z_{20}\,\mathbf{a}_3$	=	$(-x_{20}a - z_{20}c\cos\beta)\;\hat{\mathbf{x}} - y_{20}b\,\hat{\mathbf{y}} - z_{20}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O VIII
$\mathbf{B}_{76}$	=	$x_{20}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{20}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{20}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{20}a + z_{20}c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{20}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{20}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O VIII
$\mathbf{B}_{77}$	=	$x_{21}\,\mathbf{a}_1 + y_{21}\,\mathbf{a}_2 + z_{21}\,\mathbf{a}_3$	=	$(x_{21}a + z_{21}c\cos\beta)\;\hat{\mathbf{x}} + y_{21}b\,\hat{\mathbf{y}} + z_{21}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O IX
$\mathbf{B}_{78}$	=	$-x_{21}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{21}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{21}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_{21}a - z_{21}c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{21}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{21}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O IX
$\mathbf{B}_{79}$	=	$-x_{21}\,\mathbf{a}_1 - y_{21}\,\mathbf{a}_2 - z_{21}\,\mathbf{a}_3$	=	$(-x_{21}a - z_{21}c\cos\beta)\;\hat{\mathbf{x}} - y_{21}b\,\hat{\mathbf{y}} - z_{21}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O IX
$\mathbf{B}_{80}$	=	$x_{21}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{21}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{21}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{21}a + z_{21}c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{21}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{21}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O IX
$\mathbf{B}_{81}$	=	$x_{22}\,\mathbf{a}_1 + y_{22}\,\mathbf{a}_2 + z_{22}\,\mathbf{a}_3$	=	$(x_{22}a + z_{22}c\cos\beta)\;\hat{\mathbf{x}} + y_{22}b\,\hat{\mathbf{y}} + z_{22}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O X
$\mathbf{B}_{82}$	=	$-x_{22}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{22}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{22}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_{22}a - z_{22}c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{22}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{22}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O X
$\mathbf{B}_{83}$	=	$-x_{22}\,\mathbf{a}_1 - y_{22}\,\mathbf{a}_2 - z_{22}\,\mathbf{a}_3$	=	$(-x_{22}a - z_{22}c\cos\beta)\;\hat{\mathbf{x}} - y_{22}b\,\hat{\mathbf{y}} - z_{22}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O X
$\mathbf{B}_{84}$	=	$x_{22}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{22}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{22}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{22}a + z_{22}c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{22}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{22}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O X

$\mathbf{B}_{85}$	=	$x_{23}\,\mathbf{a}_1 + y_{23}\,\mathbf{a}_2 + z_{23}\,\mathbf{a}_3$	=	$(x_{23}a + z_{23}c\cos\beta)\,\hat{\mathbf{x}} + y_{23}b\,\hat{\mathbf{y}} + z_{23}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O XI
$\mathbf{B}_{86}$	=	$-x_{23}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{23}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{23}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_{23}a - z_{23}c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{23}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{23}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O XI
$\mathbf{B}_{87}$	=	$-x_{23}\,\mathbf{a}_1 - y_{23}\,\mathbf{a}_2 - z_{23}\,\mathbf{a}_3$	=	$(-x_{23}a - z_{23}c\cos\beta)\,\hat{\mathbf{x}} - y_{23}b\,\hat{\mathbf{y}} - z_{23}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O XI
$\mathbf{B}_{88}$	=	$x_{23}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{23}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{23}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{23}a + z_{23}c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{23}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{23}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O XI
$\mathbf{B}_{89}$	=	$x_{24}\,\mathbf{a}_1 + y_{24}\,\mathbf{a}_2 + z_{24}\,\mathbf{a}_3$	=	$(x_{24}a + z_{24}c\cos\beta)\,\hat{\mathbf{x}} + y_{24}b\,\hat{\mathbf{y}} + z_{24}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O XII
$\mathbf{B}_{90}$	=	$-x_{24}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{24}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{24}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_{24}a - z_{24}c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{24}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{24}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O XII
$\mathbf{B}_{91}$	=	$-x_{24}\,\mathbf{a}_1 - y_{24}\,\mathbf{a}_2 - z_{24}\,\mathbf{a}_3$	=	$(-x_{24}a - z_{24}c\cos\beta)\,\hat{\mathbf{x}} - y_{24}b\,\hat{\mathbf{y}} - z_{24}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O XII
$\mathbf{B}_{92}$	=	$x_{24}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{24}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{24}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{24}a + z_{24}c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{24}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{24}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O XII
$\mathbf{B}_{93}$	=	$x_{25}\,\mathbf{a}_1 + y_{25}\,\mathbf{a}_2 + z_{25}\,\mathbf{a}_3$	=	$(x_{25}a + z_{25}c\cos\beta)\,\hat{\mathbf{x}} + y_{25}b\,\hat{\mathbf{y}} + z_{25}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S I
$\mathbf{B}_{94}$	=	$-x_{25}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{25}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{25}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_{25}a - z_{25}c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{25}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{25}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S I
$\mathbf{B}_{95}$	=	$-x_{25}\,\mathbf{a}_1 - y_{25}\,\mathbf{a}_2 - z_{25}\,\mathbf{a}_3$	=	$(-x_{25}a - z_{25}c\cos\beta)\,\hat{\mathbf{x}} - y_{25}b\,\hat{\mathbf{y}} - z_{25}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S I
$\mathbf{B}_{96}$	=	$x_{25}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{25}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{25}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{25}a + z_{25}c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{25}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{25}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S I
$\mathbf{B}_{97}$	=	$x_{26}\,\mathbf{a}_1 + y_{26}\,\mathbf{a}_2 + z_{26}\,\mathbf{a}_3$	=	$(x_{26}a + z_{26}c\cos\beta)\,\hat{\mathbf{x}} + y_{26}b\,\hat{\mathbf{y}} + z_{26}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S II
$\mathbf{B}_{98}$	=	$-x_{26}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{26}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{26}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_{26}a - z_{26}c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{26}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{26}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S II
$\mathbf{B}_{99}$	=	$-x_{26}\,\mathbf{a}_1 - y_{26}\,\mathbf{a}_2 - z_{26}\,\mathbf{a}_3$	=	$(-x_{26}a - z_{26}c\cos\beta)\,\hat{\mathbf{x}} - y_{26}b\,\hat{\mathbf{y}} - z_{26}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S II
$\mathbf{B}_{100}$	=	$x_{26}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{26}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{26}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{26}a + z_{26}c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{26}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{26}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S II

References:

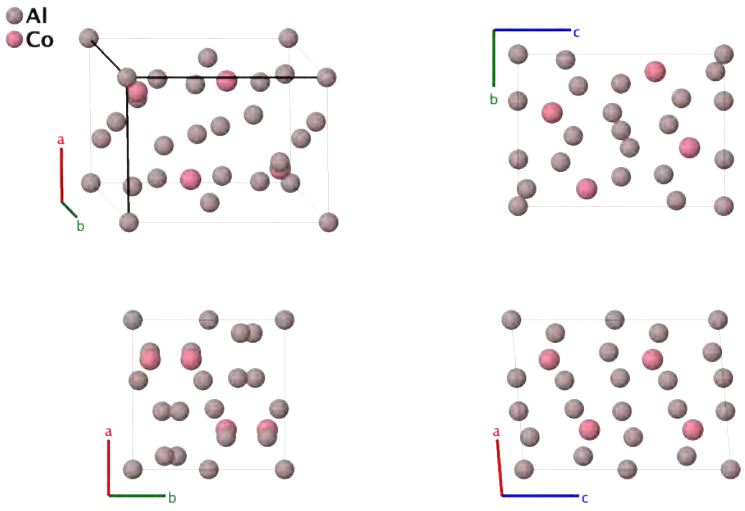
- B. Hertweck, G. Giester, and E. Libowitzky, *The crystal structures of the low-temperature phases of leonite-type compounds, $K_2Me(SO_4)_2{\cdot}4H_2O$ (Me^{2+} = Mg, Mn, Fe)*, Am. Mineral. **86**, 1282–1292 (2001), doi:10.2138/am-2001-1016.

Geometry files:

- CIF: pp. 1545
- POSCAR: pp. 1546

Co_2Al_9 ($D8_d$) Structure: A9B2_mP22_14_a4e_e

http://aflow.org/prototype-encyclopedia/A9B2_mP22_14_a4e_e



Prototype	:	Co_2Al_9
AFLOW prototype label	:	A9B2_mP22_14_a4e_e
***Strukturbericht* designation**	:	$D8_d$
Pearson symbol	:	mP22
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=A9B2_mP22_14_a4e_e --params=`$a, b/a, c/a, \beta, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6$

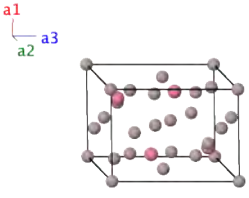
Other compounds with this structure

- Ir_2Al_9 and Rh_2Al_9

- (Douglas, 1950) gives the Wyckoff positions in terms of the $P2_1/a$ setting of space group #14. We have rotated the axes to change this to our standard $P2_1/c$ orientation.

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($2a$)	Al I

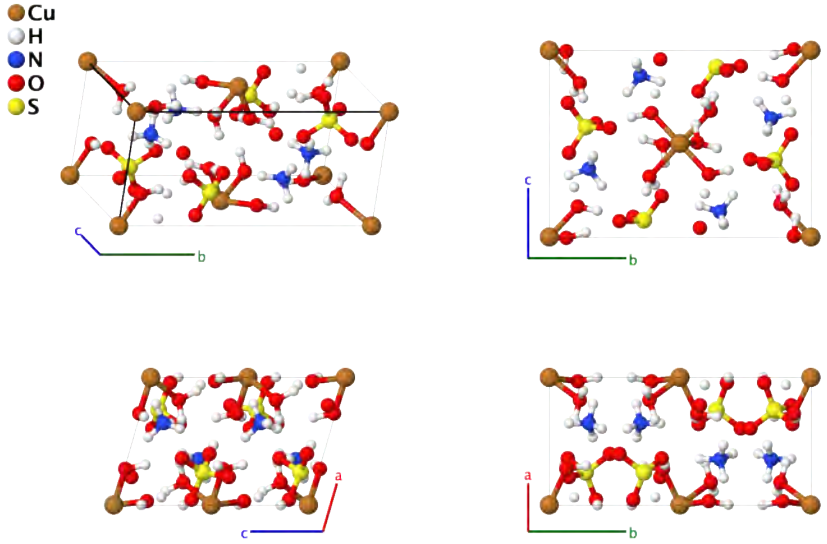
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Al I
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2a + z_2c\cos\beta)\,\hat{\mathbf{x}} + y_2b\,\hat{\mathbf{y}} + z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Al II
$\mathbf{B}_4$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2}+y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_2a - z_2c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Al II
$\mathbf{B}_5$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2a - z_2c\cos\beta)\,\hat{\mathbf{x}} - y_2b\,\hat{\mathbf{y}} - z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Al II
$\mathbf{B}_6$	=	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2}-y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_2a + z_2c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Al II
$\mathbf{B}_7$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3a + z_3c\cos\beta)\,\hat{\mathbf{x}} + y_3b\,\hat{\mathbf{y}} + z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Al III
$\mathbf{B}_8$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2}+y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_3a - z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Al III
$\mathbf{B}_9$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3a - z_3c\cos\beta)\,\hat{\mathbf{x}} - y_3b\,\hat{\mathbf{y}} - z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Al III
$\mathbf{B}_{10}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2}-y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_3a + z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Al III
$\mathbf{B}_{11}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4a + z_4c\cos\beta)\,\hat{\mathbf{x}} + y_4b\,\hat{\mathbf{y}} + z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Al IV
$\mathbf{B}_{12}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2}+y_4\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_4a - z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Al IV
$\mathbf{B}_{13}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4a - z_4c\cos\beta)\,\hat{\mathbf{x}} - y_4b\,\hat{\mathbf{y}} - z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Al IV
$\mathbf{B}_{14}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2}-y_4\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_4a + z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Al IV
$\mathbf{B}_{15}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5a + z_5c\cos\beta)\,\hat{\mathbf{x}} + y_5b\,\hat{\mathbf{y}} + z_5c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Al V
$\mathbf{B}_{16}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2}+y_5\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_5a - z_5c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Al V
$\mathbf{B}_{17}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5a - z_5c\cos\beta)\,\hat{\mathbf{x}} - y_5b\,\hat{\mathbf{y}} - z_5c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Al V
$\mathbf{B}_{18}$	=	$x_5\,\mathbf{a}_1 + \left(\frac{1}{2}-y_5\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_5a + z_5c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Al V
$\mathbf{B}_{19}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6a + z_6c\cos\beta)\,\hat{\mathbf{x}} + y_6b\,\hat{\mathbf{y}} + z_6c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Co
$\mathbf{B}_{20}$	=	$-x_6\,\mathbf{a}_1 + \left(\frac{1}{2}+y_6\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_6a - z_6c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_6\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Co
$\mathbf{B}_{21}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6a - z_6c\cos\beta)\,\hat{\mathbf{x}} - y_6b\,\hat{\mathbf{y}} - z_6c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Co
$\mathbf{B}_{22}$	=	$x_6\,\mathbf{a}_1 + \left(\frac{1}{2}-y_6\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_6a + z_6c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_6\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Co

References:
- A. M. B. Douglas, *The Structure of Co_2Al_9*, Acta Crystallogr. Sect. B Struct. Sci. **3**, 19–24 (1950), doi:10.1107/S0365110X50000057.

Geometry files:
- CIF: pp. 1546
- POSCAR: pp. 1547

Tutton salt [$Cu(NH_4)_2(SO_4)_2{\cdot}6H_2O$, $H4_4$] Structure: AB20C2D14E2_mP78_14_a_10e_e_7e_e

http://aflow.org/prototype-encyclopedia/AB20C2D14E2_mP78_14_a_10e_e_7e_e



Prototype	:	$CuH_{20}N_2O_{14}S_2$
AFLOW prototype label	:	AB20C2D14E2_mP78_14_a_10e_e_7e_e
***Strukturbericht* designation**	:	$H4_4$
Pearson symbol	:	mP78
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=AB20C2D14E2_mP78_14_a_10e_e_7e_e --params=`$a, b/a, c/a, \beta, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}, x_{17}, y_{17}, z_{17}, x_{18}, y_{18}, z_{18}, x_{19}, y_{19}, z_{19}, x_{20}, y_{20}, z_{20}$

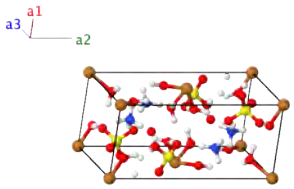
Other compounds with this structure

- $Cd(NH_4)_2(SO_4)_2{\cdot}6H_2O$, $Co(NH_4)_2(SO_4)_2{\cdot}6H_2O$, $Fe(NH_4)_2(SO_4)_2{\cdot}6H_2O$, $Mg(NH_4)_2(SO_4)_2{\cdot}6H_2O$, $Mn(NH_4)_2(SO_4)_2{\cdot}6H_2O$, $Ni(NH_4)_2(SO_4)_2{\cdot}6H_2O$, $V(NH_4)_2(SO_4)_2{\cdot}6H_2O$, $Zn(NH_4)_2(SO_4)_2{\cdot}6H_2O$, $Mg(NH_4)_2(SeO_4)_2{\cdot}6H_2O$, $Co(KSO_4)_2{\cdot}6H_2O$, $Cu(KSeO_4)_2{\cdot}6H_2O$, $Cu(KSO_4)_2{\cdot}6H_2O$, $Mg(KSO_4)_2{\cdot}6H_2O$, $Mg(TlSO_4)_2{\cdot}6H_2O$, and $Ni(KSO_4)_2{\cdot}6H_2O$

- Tutton Salts have the form $AB_2(SO_4)_2{\cdot}6H_2O$, were A is a divalent ion and B is a monovalent ion. (Hermann, 1937) formally lists $Mg(NH_4)_2(SO_4)_2{\cdot}6H_2O$ as the prototype, but notes that one may have "other monovalent metals in place of NH_4 and divalent [metals] in place of Mg." We choose $Cu(NH_4)_2(SO_4)_2{\cdot}6H_2O$ as the prototype because (Brown, 1969) were able to locate all of the hydrogen ions in agreement with expectation, *i.e.*, H_2O molecules have properly bent H-O bonds, and the hydrogen atoms in the ammonium ion form a tetrahedron around the nitrogen atom.
- The atomic positions were originally given in the $P2_1/a$ setting of space group #14. We used FINDSYM to convert this to the standard $P2_1/c$ setting. This resulted in the a and c axes being swapped.

Simple Monoclinic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= a\,\hat{\mathbf{x}}\\ \mathbf{a}_2 &= b\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	Cu
$\mathbf{B}_2$	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Cu
$\mathbf{B}_3$	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	$(x_2a + z_2c\cos\beta)\,\hat{\mathbf{x}} + y_2b\,\hat{\mathbf{y}} + z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H I
$\mathbf{B}_4$	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2}+y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_2\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta - x_2a - z_2c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H I
$\mathbf{B}_5$	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	$(-x_2a - z_2c\cos\beta)\,\hat{\mathbf{x}} - y_2b\,\hat{\mathbf{y}} - z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H I
$\mathbf{B}_6$	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2}-y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_2a + z_2c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H I
$\mathbf{B}_7$	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$(x_3a + z_3c\cos\beta)\,\hat{\mathbf{x}} + y_3b\,\hat{\mathbf{y}} + z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H II
$\mathbf{B}_8$	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2}+y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_3\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta - x_3a - z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H II
$\mathbf{B}_9$	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	$(-x_3a - z_3c\cos\beta)\,\hat{\mathbf{x}} - y_3b\,\hat{\mathbf{y}} - z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H II
$\mathbf{B}_{10}$	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2}-y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_3a + z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H II
$\mathbf{B}_{11}$	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$(x_4a + z_4c\cos\beta)\,\hat{\mathbf{x}} + y_4b\,\hat{\mathbf{y}} + z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H III
$\mathbf{B}_{12}$	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2}+y_4\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_4\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta - x_4a - z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H III
$\mathbf{B}_{13}$	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	$(-x_4a - z_4c\cos\beta)\,\hat{\mathbf{x}} - y_4b\,\hat{\mathbf{y}} - z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H III
$\mathbf{B}_{14}$	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2}-y_4\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_4a + z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H III
$\mathbf{B}_{15}$	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	$(x_5a + z_5c\cos\beta)\,\hat{\mathbf{x}} + y_5b\,\hat{\mathbf{y}} + z_5c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H IV
$\mathbf{B}_{16}$	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2}+y_5\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_5\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta - x_5a - z_5c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H IV
$\mathbf{B}_{17}$	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	$(-x_5a - z_5c\cos\beta)\,\hat{\mathbf{x}} - y_5b\,\hat{\mathbf{y}} - z_5c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H IV
$\mathbf{B}_{18}$	$x_5\,\mathbf{a}_1 + \left(\frac{1}{2}-y_5\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_5a + z_5c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H IV

$$\mathbf{B}_{19} = x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 = (x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H V}$$

$$\mathbf{B}_{20} = -x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_6 a - z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_6\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_6\right) c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H V}$$

$$\mathbf{B}_{21} = -x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3 = (-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H V}$$

$$\mathbf{B}_{22} = x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_6\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_6\right) c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H V}$$

$$\mathbf{B}_{23} = x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 = (x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H VI}$$

$$\mathbf{B}_{24} = -x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_7 a - z_7 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_7\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_7\right) c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H VI}$$

$$\mathbf{B}_{25} = -x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3 = (-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H VI}$$

$$\mathbf{B}_{26} = x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_7 a + z_7 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_7\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_7\right) c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H VI}$$

$$\mathbf{B}_{27} = x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 = (x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H VII}$$

$$\mathbf{B}_{28} = -x_8\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_8 a - z_8 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_8\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_8\right) c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H VII}$$

$$\mathbf{B}_{29} = -x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3 = (-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H VII}$$

$$\mathbf{B}_{30} = x_8\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_8 a + z_8 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_8\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_8\right) c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H VII}$$

$$\mathbf{B}_{31} = x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 = (x_9 a + z_9 c\cos\beta)\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H VIII}$$

$$\mathbf{B}_{32} = -x_9\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_9 a - z_9 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_9\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_9\right) c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H VIII}$$

$$\mathbf{B}_{33} = -x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3 = (-x_9 a - z_9 c\cos\beta)\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} - z_9 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H VIII}$$

$$\mathbf{B}_{34} = x_9\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_9 a + z_9 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_9\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_9\right) c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H VIII}$$

$$\mathbf{B}_{35} = x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 = (x_{10} a + z_{10} c\cos\beta)\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H IX}$$

$$\mathbf{B}_{36} = -x_{10}\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_{10}\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{10} a - z_{10} c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_{10}\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_{10}\right) c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H IX}$$

$$\mathbf{B}_{37} = -x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 = (-x_{10} a - z_{10} c\cos\beta)\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} - z_{10} c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H IX}$$

$$\mathbf{B}_{38} = x_{10}\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_{10}\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{10} a + z_{10} c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_{10}\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{10}\right) c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H IX}$$

$$\mathbf{B}_{39} = x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 = (x_{11} a + z_{11} c\cos\beta)\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + z_{11} c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H X}$$

$$\mathbf{B}_{40} = -x_{11}\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_{11}\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{11} a - z_{11} c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_{11}\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_{11}\right) c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H X}$$

$$\mathbf{B}_{41} = -x_{11}\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3 = (-x_{11}a - z_{11}c\cos\beta)\,\hat{\mathbf{x}} - y_{11}b\,\hat{\mathbf{y}} - z_{11}c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H X}$$

$$\mathbf{B}_{42} = x_{11}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{11}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{11}a + z_{11}c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{11}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{11}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H X}$$

$$\mathbf{B}_{43} = x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 = (x_{12}a + z_{12}c\cos\beta)\,\hat{\mathbf{x}} + y_{12}b\,\hat{\mathbf{y}} + z_{12}c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{N}$$

$$\mathbf{B}_{44} = -x_{12}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{12}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{12}a - z_{12}c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{12}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{12}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{N}$$

$$\mathbf{B}_{45} = -x_{12}\,\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3 = (-x_{12}a - z_{12}c\cos\beta)\,\hat{\mathbf{x}} - y_{12}b\,\hat{\mathbf{y}} - z_{12}c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{N}$$

$$\mathbf{B}_{46} = x_{12}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{12}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{12}a + z_{12}c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{12}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{12}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{N}$$

$$\mathbf{B}_{47} = x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3 = (x_{13}a + z_{13}c\cos\beta)\,\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} + z_{13}c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O I}$$

$$\mathbf{B}_{48} = -x_{13}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{13}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{13}a - z_{13}c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{13}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{13}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O I}$$

$$\mathbf{B}_{49} = -x_{13}\,\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 - z_{13}\,\mathbf{a}_3 = (-x_{13}a - z_{13}c\cos\beta)\,\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} - z_{13}c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O I}$$

$$\mathbf{B}_{50} = x_{13}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{13}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{13}a + z_{13}c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{13}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{13}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O I}$$

$$\mathbf{B}_{51} = x_{14}\,\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3 = (x_{14}a + z_{14}c\cos\beta)\,\hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} + z_{14}c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O II}$$

$$\mathbf{B}_{52} = -x_{14}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{14}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{14}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{14}a - z_{14}c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{14}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{14}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O II}$$

$$\mathbf{B}_{53} = -x_{14}\,\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 - z_{14}\,\mathbf{a}_3 = (-x_{14}a - z_{14}c\cos\beta)\,\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} - z_{14}c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O II}$$

$$\mathbf{B}_{54} = x_{14}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{14}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{14}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{14}a + z_{14}c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{14}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{14}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O II}$$

$$\mathbf{B}_{55} = x_{15}\,\mathbf{a}_1 + y_{15}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 = (x_{15}a + z_{15}c\cos\beta)\,\hat{\mathbf{x}} + y_{15}b\,\hat{\mathbf{y}} + z_{15}c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O III}$$

$$\mathbf{B}_{56} = -x_{15}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{15}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{15}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{15}a - z_{15}c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{15}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{15}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O III}$$

$$\mathbf{B}_{57} = -x_{15}\,\mathbf{a}_1 - y_{15}\,\mathbf{a}_2 - z_{15}\,\mathbf{a}_3 = (-x_{15}a - z_{15}c\cos\beta)\,\hat{\mathbf{x}} - y_{15}b\,\hat{\mathbf{y}} - z_{15}c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O III}$$

$$\mathbf{B}_{58} = x_{15}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{15}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{15}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{15}a + z_{15}c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{15}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{15}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O III}$$

$$\mathbf{B}_{59} = x_{16}\,\mathbf{a}_1 + y_{16}\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 = (x_{16}a + z_{16}c\cos\beta)\,\hat{\mathbf{x}} + y_{16}b\,\hat{\mathbf{y}} + z_{16}c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O IV}$$

$$\mathbf{B}_{60} = -x_{16}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{16}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{16}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{16}a - z_{16}c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{16}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{16}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O IV}$$

$$\mathbf{B}_{61} = -x_{16}\,\mathbf{a}_1 - y_{16}\,\mathbf{a}_2 - z_{16}\,\mathbf{a}_3 = (-x_{16}a - z_{16}c\cos\beta)\,\hat{\mathbf{x}} - y_{16}b\,\hat{\mathbf{y}} - z_{16}c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O IV}$$

$$\mathbf{B}_{62} = x_{16}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{16}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{16}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{16}a + z_{16}c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{16}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{16}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O IV}$$

$\mathbf{B}_{63}$	=	$x_{17}\,\mathbf{a}_1 + y_{17}\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3$	=	$(x_{17}a + z_{17}c\cos\beta)\,\hat{\mathbf{x}} + y_{17}b\,\hat{\mathbf{y}} + z_{17}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O V
$\mathbf{B}_{64}$	=	$-x_{17}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{17}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{17}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_{17}a - z_{17}c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{17}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{17}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O V
$\mathbf{B}_{65}$	=	$-x_{17}\,\mathbf{a}_1 - y_{17}\,\mathbf{a}_2 - z_{17}\,\mathbf{a}_3$	=	$(-x_{17}a - z_{17}c\cos\beta)\,\hat{\mathbf{x}} - y_{17}b\,\hat{\mathbf{y}} - z_{17}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O V
$\mathbf{B}_{66}$	=	$x_{17}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{17}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{17}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{17}a + z_{17}c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{17}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{17}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O V
$\mathbf{B}_{67}$	=	$x_{18}\,\mathbf{a}_1 + y_{18}\,\mathbf{a}_2 + z_{18}\,\mathbf{a}_3$	=	$(x_{18}a + z_{18}c\cos\beta)\,\hat{\mathbf{x}} + y_{18}b\,\hat{\mathbf{y}} + z_{18}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O VI
$\mathbf{B}_{68}$	=	$-x_{18}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{18}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{18}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_{18}a - z_{18}c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{18}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{18}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O VI
$\mathbf{B}_{69}$	=	$-x_{18}\,\mathbf{a}_1 - y_{18}\,\mathbf{a}_2 - z_{18}\,\mathbf{a}_3$	=	$(-x_{18}a - z_{18}c\cos\beta)\,\hat{\mathbf{x}} - y_{18}b\,\hat{\mathbf{y}} - z_{18}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O VI
$\mathbf{B}_{70}$	=	$x_{18}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{18}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{18}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{18}a + z_{18}c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{18}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{18}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O VI
$\mathbf{B}_{71}$	=	$x_{19}\,\mathbf{a}_1 + y_{19}\,\mathbf{a}_2 + z_{19}\,\mathbf{a}_3$	=	$(x_{19}a + z_{19}c\cos\beta)\,\hat{\mathbf{x}} + y_{19}b\,\hat{\mathbf{y}} + z_{19}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O VII
$\mathbf{B}_{72}$	=	$-x_{19}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{19}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{19}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_{19}a - z_{19}c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{19}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{19}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O VII
$\mathbf{B}_{73}$	=	$-x_{19}\,\mathbf{a}_1 - y_{19}\,\mathbf{a}_2 - z_{19}\,\mathbf{a}_3$	=	$(-x_{19}a - z_{19}c\cos\beta)\,\hat{\mathbf{x}} - y_{19}b\,\hat{\mathbf{y}} - z_{19}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O VII
$\mathbf{B}_{74}$	=	$x_{19}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{19}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{19}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{19}a + z_{19}c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{19}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{19}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O VII
$\mathbf{B}_{75}$	=	$x_{20}\,\mathbf{a}_1 + y_{20}\,\mathbf{a}_2 + z_{20}\,\mathbf{a}_3$	=	$(x_{20}a + z_{20}c\cos\beta)\,\hat{\mathbf{x}} + y_{20}b\,\hat{\mathbf{y}} + z_{20}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S
$\mathbf{B}_{76}$	=	$-x_{20}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{20}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{20}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_{20}a - z_{20}c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{20}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{20}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S
$\mathbf{B}_{77}$	=	$-x_{20}\,\mathbf{a}_1 - y_{20}\,\mathbf{a}_2 - z_{20}\,\mathbf{a}_3$	=	$(-x_{20}a - z_{20}c\cos\beta)\,\hat{\mathbf{x}} - y_{20}b\,\hat{\mathbf{y}} - z_{20}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S
$\mathbf{B}_{78}$	=	$x_{20}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{20}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{20}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{20}a + z_{20}c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{20}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{20}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S

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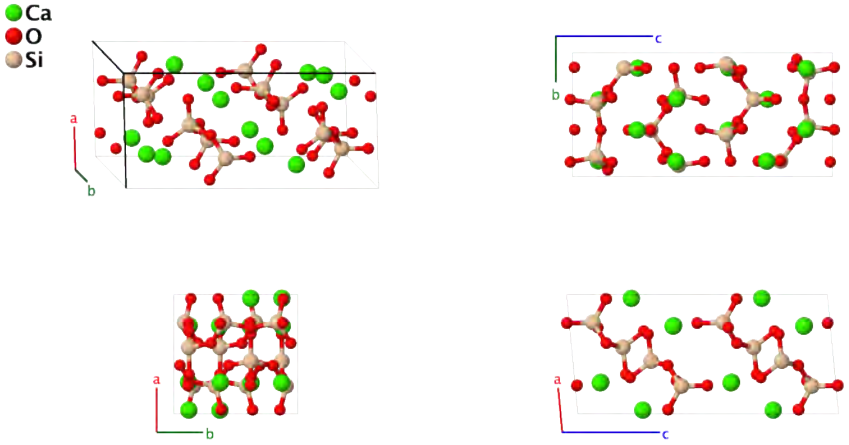
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1547
- POSCAR: pp. 1547

Parawollastonite ($CaSiO_3$, $S3_3$ (II)) Structure: AB3C_mP60_14_3e_9e_3e

http://aflow.org/prototype-encyclopedia/AB3C_mP60_14_3e_9e_3e

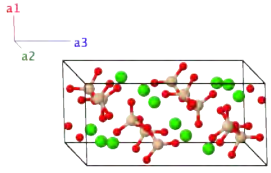


Prototype	:	CaO_3Si
AFLOW prototype label	:	AB3C_mP60_14_3e_9e_3e
***Strukturbericht* designation**	:	$S3_3$ (II)
Pearson symbol	:	mP60
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=AB3C_mP60_14_3e_9e_3e --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}$

- (Trojer, 1968) refined the original structure of (Barnick, 1936). While Barnick referred to this structure as "wollastonite," in modern terminology it is called "parawollastonite," with the original name used for the triclinic $CaSiO_3$ form.
- (Gottfried, 1938) gave this structure the $S3_3$ designation, but (Gottfried, 1937) had already used this label for crancrinite, $Na_6Ca_2Al_6Si_6O_{24}(CO_3)_2$. We will refer to parawollastonite as $S3_3$ (II), and crancrinite as $S3_3$ (I).

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Ca I

$$\mathbf{B}_2 = -x_1\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_1\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_1\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_1 a - z_1 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_1\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_1\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Ca I}$$

$$\mathbf{B}_3 = -x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3 = (-x_1 a - z_1 c\cos\beta)\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} - z_1 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Ca I}$$

$$\mathbf{B}_4 = x_1\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_1\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_1\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_1 a + z_1 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_1\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_1\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Ca I}$$

$$\mathbf{B}_5 = x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3 = (x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Ca II}$$

$$\mathbf{B}_6 = -x_2\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_2\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_2\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_2 a - z_2 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_2\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_2\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Ca II}$$

$$\mathbf{B}_7 = -x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3 = (-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} - z_2 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Ca II}$$

$$\mathbf{B}_8 = x_2\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_2\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_2\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_2 a + z_2 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_2\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_2\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Ca II}$$

$$\mathbf{B}_9 = x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3 = (x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Ca III}$$

$$\mathbf{B}_{10} = -x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_3 a - z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_3\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_3\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Ca III}$$

$$\mathbf{B}_{11} = -x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3 = (-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Ca III}$$

$$\mathbf{B}_{12} = x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_3\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_3\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Ca III}$$

$$\mathbf{B}_{13} = x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3 = (x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O I}$$

$$\mathbf{B}_{14} = -x_4\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_4 a - z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_4\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_4\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O I}$$

$$\mathbf{B}_{15} = -x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3 = (-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O I}$$

$$\mathbf{B}_{16} = x_4\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_4\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_4\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O I}$$

$$\mathbf{B}_{17} = x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3 = (x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O II}$$

$$\mathbf{B}_{18} = -x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_5 a - z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_5\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_5\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O II}$$

$$\mathbf{B}_{19} = -x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3 = (-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O II}$$

$$\mathbf{B}_{20} = x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_5\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_5\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O II}$$

$$\mathbf{B}_{21} = x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 = (x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O III}$$

$$\mathbf{B}_{22} = -x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_6 a - z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_6\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_6\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O III}$$

$$\mathbf{B}_{23} = -x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3 = (-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O III}$$

$$\mathbf{B}_{24} = x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_6\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O III}$$

$$\mathbf{B}_{25} = x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 = (x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O IV}$$

$$\mathbf{B}_{26} = -x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_7 a - z_7 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_7\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_7\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O IV}$$

$$\mathbf{B}_{27} = -x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3 = (-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O IV}$$

$$\mathbf{B}_{28} = x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_7 a + z_7 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_7\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_7\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O IV}$$

$$\mathbf{B}_{29} = x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 = (x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O V}$$

$$\mathbf{B}_{30} = -x_8\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_8 a - z_8 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_8\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_8\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O V}$$

$$\mathbf{B}_{31} = -x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3 = (-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O V}$$

$$\mathbf{B}_{32} = x_8\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_8 a + z_8 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_8\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_8\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O V}$$

$$\mathbf{B}_{33} = x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 = (x_9 a + z_9 c\cos\beta)\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O VI}$$

$$\mathbf{B}_{34} = -x_9\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_9 a - z_9 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_9\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_9\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O VI}$$

$$\mathbf{B}_{35} = -x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3 = (-x_9 a - z_9 c\cos\beta)\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} - z_9 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O VI}$$

$$\mathbf{B}_{36} = x_9\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_9 a + z_9 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_9\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_9\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O VI}$$

$$\mathbf{B}_{37} = x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 = (x_{10} a + z_{10} c\cos\beta)\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O VII}$$

$$\mathbf{B}_{38} = -x_{10}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{10}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{10} a - z_{10} c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{10}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{10}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O VII}$$

$$\mathbf{B}_{39} = -x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 = (-x_{10} a - z_{10} c\cos\beta)\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} - z_{10} c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O VII}$$

$$\mathbf{B}_{40} = x_{10}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{10}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{10} a + z_{10} c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{10}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{10}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O VII}$$

$$\mathbf{B}_{41} = x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 = (x_{11} a + z_{11} c\cos\beta)\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + z_{11} c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O VIII}$$

$$\mathbf{B}_{42} = -x_{11}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{11}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{11} a - z_{11} c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{11}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{11}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O VIII}$$

$$\mathbf{B}_{43} = -x_{11}\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3 = (-x_{11} a - z_{11} c\cos\beta)\,\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} - z_{11} c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O VIII}$$

$$\mathbf{B}_{44} = x_{11}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{11}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{11} a + z_{11} c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{11}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{11}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O VIII}$$

$$\mathbf{B}_{45} = x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 = (x_{12} a + z_{12} c\cos\beta)\,\hat{\mathbf{x}} + y_{12} b\,\hat{\mathbf{y}} + z_{12} c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O IX}$$

$$\mathbf{B}_{46} = -x_{12}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{12}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{12}a - z_{12}c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{12}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{12}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O IX}$$

$$\mathbf{B}_{47} = -x_{12}\,\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3 = (-x_{12}a - z_{12}c\cos\beta)\,\hat{\mathbf{x}} - y_{12}b\,\hat{\mathbf{y}} - z_{12}c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O IX}$$

$$\mathbf{B}_{48} = x_{12}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{12}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{12}a + z_{12}c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{12}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{12}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O IX}$$

$$\mathbf{B}_{49} = x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3 = (x_{13}a + z_{13}c\cos\beta)\,\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} + z_{13}c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Si I}$$

$$\mathbf{B}_{50} = -x_{13}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{13}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{13}a - z_{13}c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{13}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{13}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Si I}$$

$$\mathbf{B}_{51} = -x_{13}\,\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 - z_{13}\,\mathbf{a}_3 = (-x_{13}a - z_{13}c\cos\beta)\,\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} - z_{13}c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Si I}$$

$$\mathbf{B}_{52} = x_{13}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{13}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{13}a + z_{13}c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{13}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{13}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Si I}$$

$$\mathbf{B}_{53} = x_{14}\,\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3 = (x_{14}a + z_{14}c\cos\beta)\,\hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} + z_{14}c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Si II}$$

$$\mathbf{B}_{54} = -x_{14}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{14}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{14}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{14}a - z_{14}c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{14}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{14}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Si II}$$

$$\mathbf{B}_{55} = -x_{14}\,\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 - z_{14}\,\mathbf{a}_3 = (-x_{14}a - z_{14}c\cos\beta)\,\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} - z_{14}c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Si II}$$

$$\mathbf{B}_{56} = x_{14}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{14}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{14}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{14}a + z_{14}c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{14}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{14}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Si II}$$

$$\mathbf{B}_{57} = x_{15}\,\mathbf{a}_1 + y_{15}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 = (x_{15}a + z_{15}c\cos\beta)\,\hat{\mathbf{x}} + y_{15}b\,\hat{\mathbf{y}} + z_{15}c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Si III}$$

$$\mathbf{B}_{58} = -x_{15}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{15}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{15}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{15}a - z_{15}c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{15}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{15}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Si III}$$

$$\mathbf{B}_{59} = -x_{15}\,\mathbf{a}_1 - y_{15}\,\mathbf{a}_2 - z_{15}\,\mathbf{a}_3 = (-x_{15}a - z_{15}c\cos\beta)\,\hat{\mathbf{x}} - y_{15}b\,\hat{\mathbf{y}} - z_{15}c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Si III}$$

$$\mathbf{B}_{60} = x_{15}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{15}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{15}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{15}a + z_{15}c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{15}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{15}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Si III}$$

References:

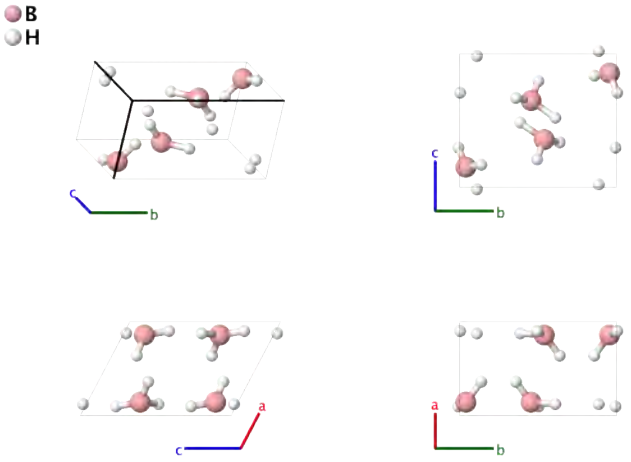
- F. J. Trojer, *The crystal structure of parawollastonite*, Zeitschrift für Kristallographie - Crystalline Materials **127**, 291–308 (1968), doi:10.1524/zkri.1968.127.16.291.
- M. Barnick, *Strukturuntersuchung des natürlichen Wollastonits* (1936). Dissertation.
- C. Gottfried, ed., *Strukturbericht Band IV 1936* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1938).
- C. Gottfried and F. Schossberger, eds., *Strukturbericht Band III 1933-1935* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1548
- POSCAR: pp. 1548

β-B_2H_6 Structure: AB3_mP16_14_e_3e

http://aflow.org/prototype-encyclopedia/AB3_mP16_14_e_3e.beta

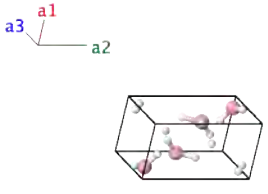


Prototype	:	β-B_2H_6
AFLOW prototype label	:	AB3_mP16_14_e_3e
***Strukturbericht* designation**	:	None
Pearson symbol	:	mP16
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=AB3_mP16_14_e_3e --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4$

- (Mark, 1925) studied a variety of B_2H_6 structures, including the one given *Strukturbericht* designation $D4_1$ by (Ewald, 1931). (Smith, 1965) refined this structure, including the hydrogen positions.
- This structure shares the same AFLOW designation, AB3_mP16_14_e_3e, as the B_2H_6 structure defined by (Yao, 2011). If we remove the hydrogen atoms from either of these structures we get the A10 (α-Hg) structure, but that is more closely associated with the Yao structure than it is with this one.
- (Smith, 1965) presented the crystallographic information for this structure in the $P2_1/n$, unique axis c, setting of space group #14. We used FINDSYM to put the information into the standard $P2_1/c$, unique axis b, setting. This involved modifications of the primitive vectors beyond a simple rotation.

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

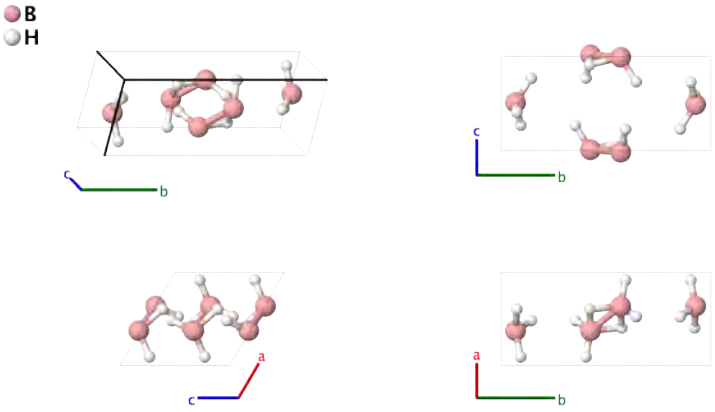
		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1a + z_1c\cos\beta)\,\hat{\mathbf{x}} + y_1b\,\hat{\mathbf{y}} + z_1c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	B
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2} + y_1\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_1a - z_1c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	B
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$(-x_1a - z_1c\cos\beta)\,\hat{\mathbf{x}} - y_1b\,\hat{\mathbf{y}} - z_1c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	B
$\mathbf{B}_4$	=	$x_1\,\mathbf{a}_1 + \left(\frac{1}{2} - y_1\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_1a + z_1c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	B
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2a + z_2c\cos\beta)\,\hat{\mathbf{x}} + y_2b\,\hat{\mathbf{y}} + z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H I
$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_2a - z_2c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H I
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2a - z_2c\cos\beta)\,\hat{\mathbf{x}} - y_2b\,\hat{\mathbf{y}} - z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H I
$\mathbf{B}_8$	=	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_2a + z_2c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H I
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3a + z_3c\cos\beta)\,\hat{\mathbf{x}} + y_3b\,\hat{\mathbf{y}} + z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H II
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_3a - z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H II
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3a - z_3c\cos\beta)\,\hat{\mathbf{x}} - y_3b\,\hat{\mathbf{y}} - z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H II
$\mathbf{B}_{12}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_3a + z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H II
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4a + z_4c\cos\beta)\,\hat{\mathbf{x}} + y_4b\,\hat{\mathbf{y}} + z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H III
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_4a - z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H III
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4a - z_4c\cos\beta)\,\hat{\mathbf{x}} - y_4b\,\hat{\mathbf{y}} - z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H III
$\mathbf{B}_{16}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_4a + z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H III

References:
- H. W. Smith and W. N. Lipscomb, *Single-Crystal X-Ray Diffraction Study of β-Diborane*, J. Chem. Phys. **43**, 1060–1064 (1965), doi:10.1063/1.1696820.
- H. Mark and E. Pohland, *IV. Über die Gitterstruktur des Äthans und des Diborans*, Zeitschrift für Kristallographie - Crystalline Materials **62**, 103–112 (1925), doi:10.1524/zkri.1925.62.1.103.
- P. P. Ewald and C. Hermann, eds., *Strukturbericht 1913-1928* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1931).
- Y. Yao and R. Hoffmann, *BH_3 under Pressure: Leaving the Molecular Diborane Motif*, J. Am. Chem. Soc. **133**, 21002–21009 (2011), doi:10.1021/ja2092568.

Geometry files:
- CIF: pp. 1549
- POSCAR: pp. 1549

B_2H_6 ($P2_1/c$) Structure: AB3_mP16_14_e_3e

http://aflow.org/prototype-encyclopedia/AB3_mP16_14_e_3e

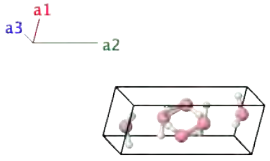


Prototype	:	B_2H_6
AFLOW prototype label	:	AB3_mP16_14_e_3e
***Strukturbericht* designation**	:	None
Pearson symbol	:	mP16
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=AB3_mP16_14_e_3e --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4$

- (Mark, 1925) studied a variety of B_2H_6 structures, including the one given *Strukturbericht* designation $D4_1$ by (Ewald, 1931). Very little information was given about this structure, other than Ewalds's statement "The position of the boron atoms is exactly the same as that of the Hg atoms in the A10 (α-Hg) structure." (Yao, 2011) found the current structure computationally, and note that it is likely the structure found by (Mark, 1925).
- This structure shares the same AFLOW designation, AB3_mP16_14_e_3e, as the β-B_2H_6 structure described by (Smith, 1965).

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$ =	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$ =	$(x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\sin\beta\,\hat{\mathbf{z}}$	($4e$)	B
$\mathbf{B}_2$ =	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2}+y_1\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_1\right)\mathbf{a}_3$ =	$\left(\frac{1}{2}c\cos\beta - x_1 a - z_1 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	($4e$)	B

$$\mathbf{B}_3 = -x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3 = (-x_1 a - z_1 c\cos\beta)\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} - z_1 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{B}$$

$$\mathbf{B}_4 = x_1\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_1\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_1\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_1 a + z_1 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_1\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_1\right) c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{B}$$

$$\mathbf{B}_5 = x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3 = (x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H I}$$

$$\mathbf{B}_6 = -x_2\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_2\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_2\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_2 a - z_2 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_2\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_2\right) c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H I}$$

$$\mathbf{B}_7 = -x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3 = (-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} - z_2 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H I}$$

$$\mathbf{B}_8 = x_2\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_2\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_2\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_2 a + z_2 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_2\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_2\right) c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H I}$$

$$\mathbf{B}_9 = x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3 = (x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H II}$$

$$\mathbf{B}_{10} = -x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_3 a - z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_3\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_3\right) c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H II}$$

$$\mathbf{B}_{11} = -x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3 = (-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H II}$$

$$\mathbf{B}_{12} = x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_3\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_3\right) c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H II}$$

$$\mathbf{B}_{13} = x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3 = (x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H III}$$

$$\mathbf{B}_{14} = -x_4\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_4 a - z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_4\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_4\right) c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H III}$$

$$\mathbf{B}_{15} = -x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3 = (-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H III}$$

$$\mathbf{B}_{16} = x_4\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_4\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_4\right) c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{H III}$$

References:

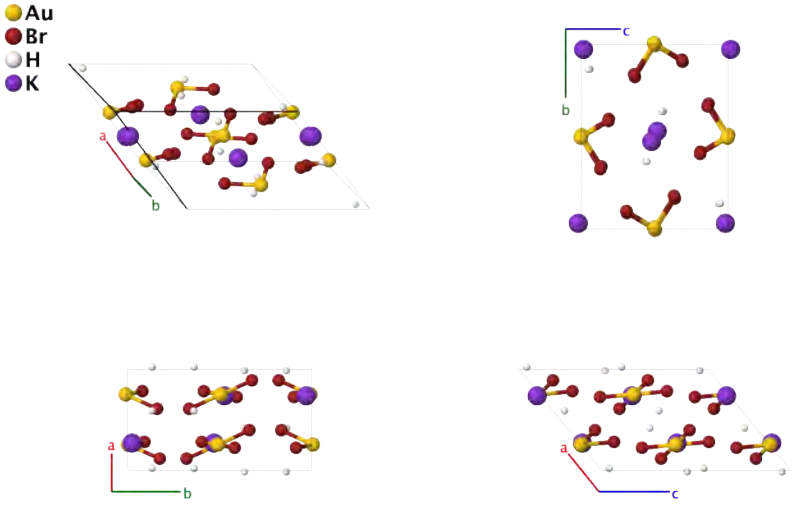
- Y. Yao and R. Hoffmann, *BH_3 under Pressure: Leaving the Molecular Diborane Motif*, J. Am. Chem. Soc. **133**, 21002–21009 (2011), doi:10.1021/ja2092568.
- H. Mark and E. Pohland, *IV. Über die Gitterstruktur des Äthans und des Diborans*, Zeitschrift für Kristallographie - Crystalline Materials **62**, 103–112 (1925), doi:10.1524/zkri.1925.62.1.103.
- P. P. Ewald and C. Hermann, eds., *Strukturbericht 1913-1928* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1931).
- H. W. Smith and W. N. Lipscomb, *Single-Crystal X-Ray Diffraction Study of β-Diborane*, J. Chem. Phys. **43**, 1060–1064 (1965), doi:10.1063/1.1696820.

Geometry files:

- CIF: pp. 1549
- POSCAR: pp. 1549

$KAuBr_4{\cdot}2H_2O$ ($H4_{19}$) Structure: AB4C2D_mP32_14_e_4e_2e_e

http://aflow.org/prototype-encyclopedia/AB4C2D_mP32_14_e_4e_2e_e

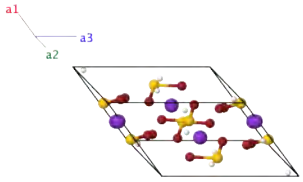


Prototype	:	$AuBr_4(H_2O)_2K$
AFLOW prototype label	:	AB4C2D_mP32_14_e_4e_2e_e
***Strukturbericht* designation**	:	$H4_{19}$
Pearson symbol	:	mP32
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=AB4C2D_mP32_14_e_4e_2e_e --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8$

- (Cox, 1936) determined lattice constants for $KAuBr_4{\cdot}H_2O$, which were very close to those later found by (Omrani, 1986). However, Cox *et al.* found that the (100) planes (in our orientation) with odd indices showed very weak diffraction spots. Neglecting these spots gives a unit cell which has a lattice constant a (in our orientation) which is half the size of that found by Omrani *et al*, and so the Pearson symbol mP16. Fitting this cell into space group $P2_1/c$ #14 then required that the gold atom be at the ($2a$) Wyckoff position and the potassium atom at ($2b$), while the later paper found that these atoms are slightly displaced from these points.
- The results of (Omrani, 1986) are very close to those of (Cox, 1936) and can indeed reduce to the Cox *et al.* structure by allowing some uncertainty in the atomic positions. Given this, and the fact that the former reference actually found the correct unit cell, we will use the more modern work as our prototype for *Strukturbericht* symbol $H4_{19}$.
- (Omrani, 1986) gave the unit cell and Wyckoff positions in terms of setting $P2_1/n$ of space group #14. Changing this to the standard $P2_1/c$ setting required a rotation of the lattice and a significant rewriting of the primitive vectors.
- The anhydrous form of this compound can be seen at $KAuBr_4$ (AB4C_mP24_14_ab_4e_e).

Simple Monoclinic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= a\,\hat{\mathbf{x}}\\ \mathbf{a}_2 &= b\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	$(x_1a + z_1c\cos\beta)\,\hat{\mathbf{x}} + y_1b\,\hat{\mathbf{y}} + z_1c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Au
$\mathbf{B}_2$	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2}+y_1\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_1\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta - x_1a - z_1c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Au
$\mathbf{B}_3$	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	$(-x_1a - z_1c\cos\beta)\,\hat{\mathbf{x}} - y_1b\,\hat{\mathbf{y}} - z_1c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Au
$\mathbf{B}_4$	$x_1\,\mathbf{a}_1 + \left(\frac{1}{2}-y_1\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_1\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_1a + z_1c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Au
$\mathbf{B}_5$	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	$(x_2a + z_2c\cos\beta)\,\hat{\mathbf{x}} + y_2b\,\hat{\mathbf{y}} + z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Br I
$\mathbf{B}_6$	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2}+y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_2\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta - x_2a - z_2c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Br I
$\mathbf{B}_7$	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	$(-x_2a - z_2c\cos\beta)\,\hat{\mathbf{x}} - y_2b\,\hat{\mathbf{y}} - z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Br I
$\mathbf{B}_8$	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2}-y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_2a + z_2c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Br I
$\mathbf{B}_9$	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$(x_3a + z_3c\cos\beta)\,\hat{\mathbf{x}} + y_3b\,\hat{\mathbf{y}} + z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Br II
$\mathbf{B}_{10}$	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2}+y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_3\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta - x_3a - z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Br II
$\mathbf{B}_{11}$	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	$(-x_3a - z_3c\cos\beta)\,\hat{\mathbf{x}} - y_3b\,\hat{\mathbf{y}} - z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Br II
$\mathbf{B}_{12}$	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2}-y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_3a + z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Br II
$\mathbf{B}_{13}$	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$(x_4a + z_4c\cos\beta)\,\hat{\mathbf{x}} + y_4b\,\hat{\mathbf{y}} + z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Br III
$\mathbf{B}_{14}$	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2}+y_4\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_4\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta - x_4a - z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Br III
$\mathbf{B}_{15}$	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	$(-x_4a - z_4c\cos\beta)\,\hat{\mathbf{x}} - y_4b\,\hat{\mathbf{y}} - z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Br III
$\mathbf{B}_{16}$	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2}-y_4\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_4a + z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Br III
$\mathbf{B}_{17}$	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	$(x_5a + z_5c\cos\beta)\,\hat{\mathbf{x}} + y_5b\,\hat{\mathbf{y}} + z_5c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Br IV

$\mathbf{B}_{18}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_5 a - z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Br IV
$\mathbf{B}_{19}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Br IV
$\mathbf{B}_{20}$	=	$x_5\,\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Br IV
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H_2O I
$\mathbf{B}_{22}$	=	$-x_6\,\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_6 a - z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H_2O I
$\mathbf{B}_{23}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H_2O I
$\mathbf{B}_{24}$	=	$x_6\,\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H_2O I
$\mathbf{B}_{25}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H_2O II
$\mathbf{B}_{26}$	=	$-x_7\,\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_7 a - z_7 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H_2O II
$\mathbf{B}_{27}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H_2O II
$\mathbf{B}_{28}$	=	$x_7\,\mathbf{a}_1 + \left(\frac{1}{2} - y_7\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_7 a + z_7 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_7\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H_2O II
$\mathbf{B}_{29}$	=	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	K
$\mathbf{B}_{30}$	=	$-x_8\,\mathbf{a}_1 + \left(\frac{1}{2} + y_8\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_8 a - z_8 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_8\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	K
$\mathbf{B}_{31}$	=	$-x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	K
$\mathbf{B}_{32}$	=	$x_8\,\mathbf{a}_1 + \left(\frac{1}{2} - y_8\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_8 a + z_8 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_8\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	K

References:

- H. Omrani, F. Théobald, and H. Vivier, *Structure of potassium tetrabromoaurate(III) dihydrate*, Acta Crystallogr. C **42**, 1091–1092 (1986), doi:10.1107/S0108270186093344.
- E. G. Cox and K. C. Webster, *The stereochemistry of quadricovalent atoms: tervalent gold*, J. Chem. Soc. pp. 1635–1637 (1936), doi:10.1039/JR9360001635.

Found in:

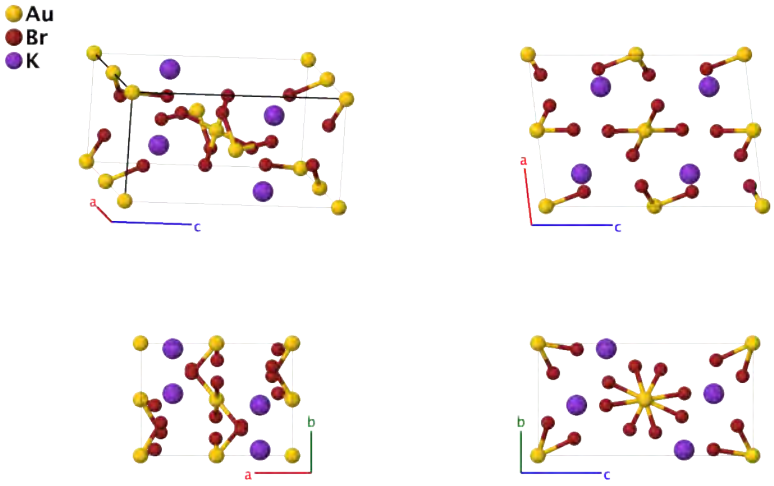
- R. Welter, H. Omrani, and R. Vangelisti, *Sodium tetrabromoaurate(III) dihydrate*, Acta Crystallogr. E **57**, i8–i9 (2001), doi:10.1107/S1600536800021140.

Geometry files:

- CIF: pp. 1550
- POSCAR: pp. 1550

Anhydrous $KAuBr_4$ Structure: AB4C_mP24_14_ab_4e_e

http://aflow.org/prototype-encyclopedia/AB4C_mP24_14_ab_4e_e

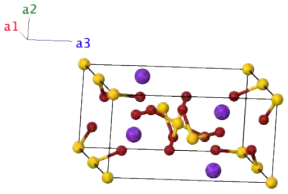


Prototype	:	$AuBr_4K$
AFLOW prototype label	:	AB4C_mP24_14_ab_4e_e
***Strukturbericht* designation**	:	None
Pearson symbol	:	mP24
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=AB4C_mP24_14_ab_4e_e --params=`$a, b/a, c/a, \beta, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7$

- This hydrated form of this compound is $KAuBr_4{\cdot}2H_2O$ ($H4_{19}$).

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($2a$)	Au I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	($2a$)	Au I
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	($2b$)	Au II
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(a + c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	($2b$)	Au II

$$\mathbf{B}_5 = x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3 = (x_3a + z_3c\cos\beta)\,\hat{\mathbf{x}} + y_3b\,\hat{\mathbf{y}} + z_3c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Br I}$$

$$\mathbf{B}_6 = -x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_3a - z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_3\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_3\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Br I}$$

$$\mathbf{B}_7 = -x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3 = (-x_3a - z_3c\cos\beta)\,\hat{\mathbf{x}} - y_3b\,\hat{\mathbf{y}} - z_3c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Br I}$$

$$\mathbf{B}_8 = x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_3a + z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_3\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_3\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Br I}$$

$$\mathbf{B}_9 = x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3 = (x_4a + z_4c\cos\beta)\,\hat{\mathbf{x}} + y_4b\,\hat{\mathbf{y}} + z_4c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Br II}$$

$$\mathbf{B}_{10} = -x_4\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_4a - z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_4\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_4\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Br II}$$

$$\mathbf{B}_{11} = -x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3 = (-x_4a - z_4c\cos\beta)\,\hat{\mathbf{x}} - y_4b\,\hat{\mathbf{y}} - z_4c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Br II}$$

$$\mathbf{B}_{12} = x_4\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_4a + z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_4\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_4\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Br II}$$

$$\mathbf{B}_{13} = x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3 = (x_5a + z_5c\cos\beta)\,\hat{\mathbf{x}} + y_5b\,\hat{\mathbf{y}} + z_5c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Br III}$$

$$\mathbf{B}_{14} = -x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_5a - z_5c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_5\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_5\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Br III}$$

$$\mathbf{B}_{15} = -x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3 = (-x_5a - z_5c\cos\beta)\,\hat{\mathbf{x}} - y_5b\,\hat{\mathbf{y}} - z_5c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Br III}$$

$$\mathbf{B}_{16} = x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_5a + z_5c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_5\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_5\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Br III}$$

$$\mathbf{B}_{17} = x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 = (x_6a + z_6c\cos\beta)\,\hat{\mathbf{x}} + y_6b\,\hat{\mathbf{y}} + z_6c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Br IV}$$

$$\mathbf{B}_{18} = -x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_6a - z_6c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_6\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_6\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Br IV}$$

$$\mathbf{B}_{19} = -x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3 = (-x_6a - z_6c\cos\beta)\,\hat{\mathbf{x}} - y_6b\,\hat{\mathbf{y}} - z_6c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Br IV}$$

$$\mathbf{B}_{20} = x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_6a + z_6c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_6\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Br IV}$$

$$\mathbf{B}_{21} = x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 = (x_7a + z_7c\cos\beta)\,\hat{\mathbf{x}} + y_7b\,\hat{\mathbf{y}} + z_7c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{K}$$

$$\mathbf{B}_{22} = -x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_7a - z_7c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_7\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_7\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{K}$$

$$\mathbf{B}_{23} = -x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3 = (-x_7a - z_7c\cos\beta)\,\hat{\mathbf{x}} - y_7b\,\hat{\mathbf{y}} - z_7c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{K}$$

$$\mathbf{B}_{24} = x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_7a + z_7c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_7\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_7\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{K}$$

References:

- H. Omrani, R. Welter, and R. Vangelisti, *Potassium tetrabromoaurate(III)*, Acta Crystallogr. C **55**, 13–14 (1999), doi:10.1107/S010827019801110X.

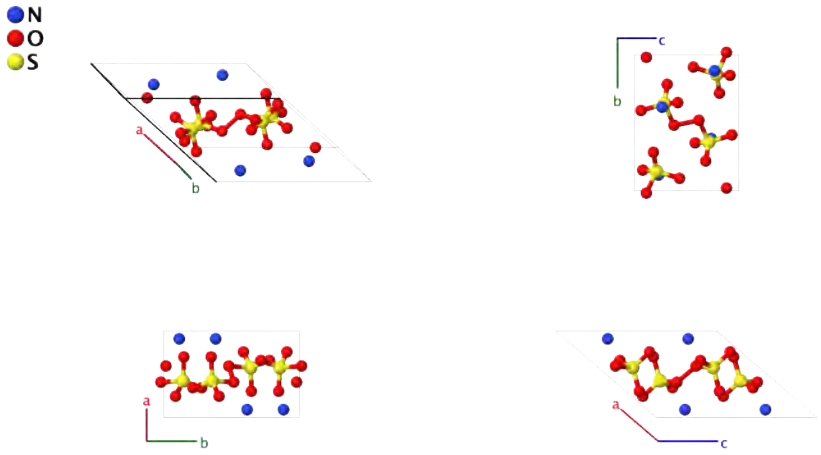
Found in:

- R. Welter, H. Omrani, and R. Vangelisti, *Sodium tetrabromoaurate(III) dihydrate*, Acta Crystallogr. E **57**, i8–i9 (2001), doi:10.1107/S1600536800021140.

Geometry files:
- CIF: pp. 1550
- POSCAR: pp. 1551

Ammonium Persulfate [$(NH_4)_2S_2O_8$, $K4_1$] Structure: AB4C_mP24_14_e_4e_e

http://aflow.org/prototype-encyclopedia/AB4C_mP24_14_e_4e_e.K41



Prototype	:	$(NH_4)O_4S$
AFLOW prototype label	:	AB4C_mP24_14_e_4e_e
***Strukturbericht* designation**	:	$K4_1$
Pearson symbol	:	mP24
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=AB4C_mP24_14_e_4e_e --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6$

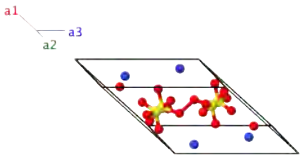
Other compounds with this structure

- $CaSO_4$

- This structure has the same AFLOW label as monazite, $LaPO_4$. The structures are generated by the same symmetry operations with different sets of parameters (`--params`) specified in their corresponding CIF files.

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$$
\begin{array}{lcccccc}
\mathbf{B}_1 & = & x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3 & = & (x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\sin\beta\,\hat{\mathbf{z}} & (4e) & \mathrm{NH_4} \\
\mathbf{B}_2 & = & -x_1\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_1\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_1\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}c\cos\beta - x_1 a - z_1 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_1\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_1\right) c\sin\beta\,\hat{\mathbf{z}} & (4e) & \mathrm{NH_4} \\
\mathbf{B}_3 & = & -x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3 & = & (-x_1 a - z_1 c\cos\beta)\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} - z_1 c\sin\beta\,\hat{\mathbf{z}} & (4e) & \mathrm{NH_4} \\
\mathbf{B}_4 & = & x_1\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_1\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_1\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}c\cos\beta + x_1 a + z_1 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_1\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_1\right) c\sin\beta\,\hat{\mathbf{z}} & (4e) & \mathrm{NH_4} \\
\mathbf{B}_5 & = & x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3 & = & (x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}} & (4e) & \mathrm{O\ I} \\
\mathbf{B}_6 & = & -x_2\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_2\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_2\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}c\cos\beta - x_2 a - z_2 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_2\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_2\right) c\sin\beta\,\hat{\mathbf{z}} & (4e) & \mathrm{O\ I} \\
\mathbf{B}_7 & = & -x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3 & = & (-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} - z_2 c\sin\beta\,\hat{\mathbf{z}} & (4e) & \mathrm{O\ I} \\
\mathbf{B}_8 & = & x_2\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_2\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_2\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}c\cos\beta + x_2 a + z_2 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_2\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_2\right) c\sin\beta\,\hat{\mathbf{z}} & (4e) & \mathrm{O\ I} \\
\mathbf{B}_9 & = & x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3 & = & (x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}} & (4e) & \mathrm{O\ II} \\
\mathbf{B}_{10} & = & -x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_3\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}c\cos\beta - x_3 a - z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_3\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_3\right) c\sin\beta\,\hat{\mathbf{z}} & (4e) & \mathrm{O\ II} \\
\mathbf{B}_{11} & = & -x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3 & = & (-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}} & (4e) & \mathrm{O\ II} \\
\mathbf{B}_{12} & = & x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_3\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_3\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_3\right) c\sin\beta\,\hat{\mathbf{z}} & (4e) & \mathrm{O\ II} \\
\mathbf{B}_{13} & = & x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3 & = & (x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}} & (4e) & \mathrm{O\ III} \\
\mathbf{B}_{14} & = & -x_4\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_4\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}c\cos\beta - x_4 a - z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_4\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_4\right) c\sin\beta\,\hat{\mathbf{z}} & (4e) & \mathrm{O\ III} \\
\mathbf{B}_{15} & = & -x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3 & = & (-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}} & (4e) & \mathrm{O\ III} \\
\mathbf{B}_{16} & = & x_4\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_4\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_4\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_4\right) c\sin\beta\,\hat{\mathbf{z}} & (4e) & \mathrm{O\ III} \\
\mathbf{B}_{17} & = & x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3 & = & (x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}} & (4e) & \mathrm{O\ IV} \\
\mathbf{B}_{18} & = & -x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_5\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}c\cos\beta - x_5 a - z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_5\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_5\right) c\sin\beta\,\hat{\mathbf{z}} & (4e) & \mathrm{O\ IV} \\
\mathbf{B}_{19} & = & -x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3 & = & (-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}} & (4e) & \mathrm{O\ IV} \\
\mathbf{B}_{20} & = & x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_5\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_5\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_5\right) c\sin\beta\,\hat{\mathbf{z}} & (4e) & \mathrm{O\ IV} \\
\mathbf{B}_{21} & = & x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 & = & (x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}} & (4e) & \mathrm{S} \\
\mathbf{B}_{22} & = & -x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_6\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}c\cos\beta - x_6 a - z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_6\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_6\right) c\sin\beta\,\hat{\mathbf{z}} & (4e) & \mathrm{S}
\end{array}
$$

$\mathbf{B}_{23}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	S
$\mathbf{B}_{24}$	=	$x_6\,\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	S

References:

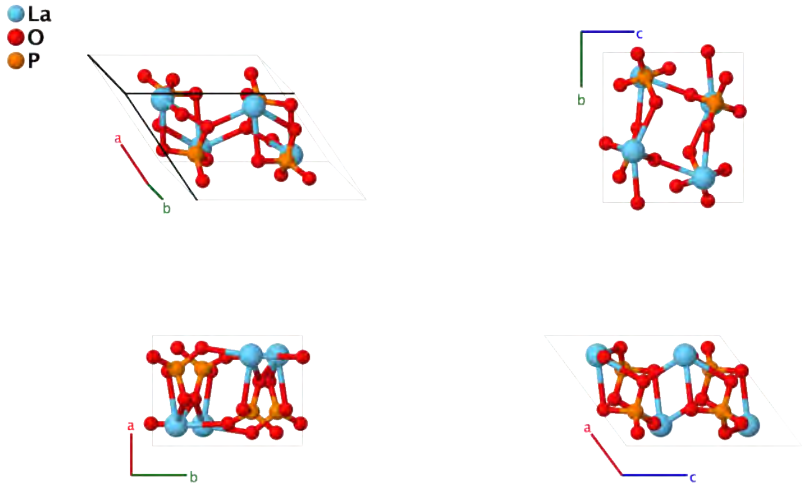
- B. K. Sivertsen and H. Sorum, *A reinvestigation of the crystal structure of ammonium persulfate, $(NH_4)_2S_2O_8$*, Zeitschrift für Kristallographie - Crystalline Materials **130**, 449–460 (1969), doi:10.1524/zkri.1969.130.16.449.

Geometry files:

- CIF: pp. 1551
- POSCAR: pp. 1551

Monazite ($LaPO_4$) Structure: AB4C_mP24_14_e_4e_e

http://aflow.org/prototype-encyclopedia/AB4C_mP24_14_e_4e_e



Prototype	:	LaO_4P
AFLOW prototype label	:	AB4C_mP24_14_e_4e_e
***Strukturbericht* designation**	:	None
Pearson symbol	:	mP24
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=AB4C_mP24_14_e_4e_e --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6$

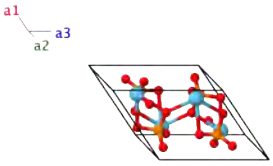
Other compounds with this structure

- $CePO_4$ (monazite-Ce), $PrPO_4$ (monazite-Pr), $NdPO_4$ (monazite-Nd), $SmPO_4$ (monazite-Sm), $EuPO_4$ (monazite-Eu), $LaAsO_4$ (gasparite-La), and $CeAsO_4$ (gasparite-Ce)

- Monazites (RE-PO_4) and gasparites (RE-AsO_4) can have a mixture of rare earth elements in the RE slot of the formula. Technically these minerals are called monazite-(X) and gasparite-(X), where "X" is the predominant rare earth element in the sample. All of the structures are similar, but we must pick one as the prototype, so we take the first entry in (Ni, 1995) to define the class.
- (Ni, 1995) gives the structural data in the $P2_1/n$ setting of space group #14. We used FINDSYM to change this to our standard $P2_1/c$ setting. This involves a change of primitive vectors as well as a rotation of the crystal.
- This structure has the same AFLOW label as NH_4SO_4, $K4_1$. The structures are generated by the same symmetry operations with different sets of parameters (`--params`) specified in their corresponding CIF files.

Simple Monoclinic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= a\,\hat{\mathbf{x}}\\ \mathbf{a}_2 &= b\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	$(x_1a + z_1c\cos\beta)\,\hat{\mathbf{x}} + y_1b\,\hat{\mathbf{y}} + z_1c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	La
$\mathbf{B}_2$	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2}+y_1\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_1\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta - x_1a - z_1c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	La
$\mathbf{B}_3$	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	$(-x_1a - z_1c\cos\beta)\,\hat{\mathbf{x}} - y_1b\,\hat{\mathbf{y}} - z_1c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	La
$\mathbf{B}_4$	$x_1\,\mathbf{a}_1 + \left(\frac{1}{2}-y_1\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_1\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_1a + z_1c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	La
$\mathbf{B}_5$	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	$(x_2a + z_2c\cos\beta)\,\hat{\mathbf{x}} + y_2b\,\hat{\mathbf{y}} + z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_6$	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2}+y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_2\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta - x_2a - z_2c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_7$	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	$(-x_2a - z_2c\cos\beta)\,\hat{\mathbf{x}} - y_2b\,\hat{\mathbf{y}} - z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_8$	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2}-y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_2a + z_2c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_9$	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$(x_3a + z_3c\cos\beta)\,\hat{\mathbf{x}} + y_3b\,\hat{\mathbf{y}} + z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_{10}$	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2}+y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_3\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta - x_3a - z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_{11}$	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	$(-x_3a - z_3c\cos\beta)\,\hat{\mathbf{x}} - y_3b\,\hat{\mathbf{y}} - z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_{12}$	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2}-y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_3a + z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_{13}$	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$(x_4a + z_4c\cos\beta)\,\hat{\mathbf{x}} + y_4b\,\hat{\mathbf{y}} + z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O III
$\mathbf{B}_{14}$	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2}+y_4\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_4\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta - x_4a - z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O III
$\mathbf{B}_{15}$	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	$(-x_4a - z_4c\cos\beta)\,\hat{\mathbf{x}} - y_4b\,\hat{\mathbf{y}} - z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O III
$\mathbf{B}_{16}$	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2}-y_4\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_4a + z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O III
$\mathbf{B}_{17}$	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	$(x_5a + z_5c\cos\beta)\,\hat{\mathbf{x}} + y_5b\,\hat{\mathbf{y}} + z_5c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O IV

$\mathbf{B}_{18}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_5 a - z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O IV
$\mathbf{B}_{19}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O IV
$\mathbf{B}_{20}$	=	$x_5\,\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O IV
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	P
$\mathbf{B}_{22}$	=	$-x_6\,\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_6 a - z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	P
$\mathbf{B}_{23}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	P
$\mathbf{B}_{24}$	=	$x_6\,\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	P

References:

- Y. Ni, J. M. Hughes, and A. N. Mariano, *Crystal chemistry of the monazite and xenotime structures*, Am. Mineral. **80**, 21–26 (1995).

Found in:

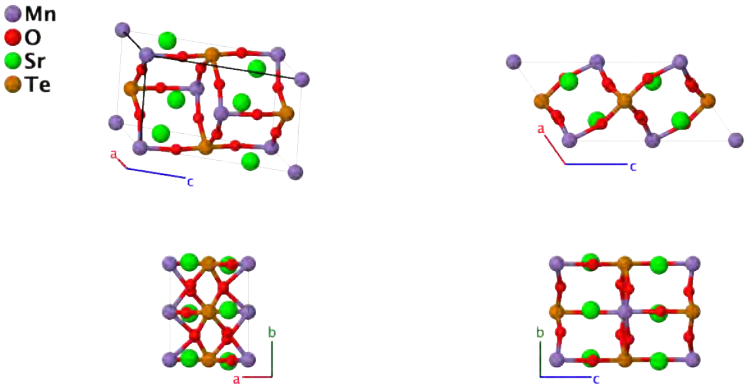
- O. S. Vereshchagin, S. N. Britvin, E. N. Perova, A. I. Brusnitsyn, Y. S. Polekhovsky, V. V. Shilovskikh, V. N. Bocharov, A. van der Burgt, S. Cuchet, and N. Meisser, *Gasparite-(La), La(AsO$_4$), a new mineral from Mn ores of the Ushkatyn-III deposit, Central Kazakhstan, and metamorphic rocks of the Wanni glacier, Switzerland*, Am. Mineral. **104**, 1469–1480 (2019), doi:10.2138/am-2019-7028.

Geometry files:

- CIF: pp. 1551
- POSCAR: pp. 1552

Sr_2MnTeO_6 Structure: AB6C2D_mP20_14_a_3e_e_d

http://aflow.org/prototype-encyclopedia/AB6C2D_mP20_14_a_3e_e_d



Prototype	:	MnO_6Sr_2Te
AFLOW prototype label	:	AB6C2D_mP20_14_a_3e_e_d
***Strukturbericht* designation**	:	None
Pearson symbol	:	mP20
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=AB6C2D_mP20_14_a_3e_e_d --params=`$a, b/a, c/a, \beta, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6$

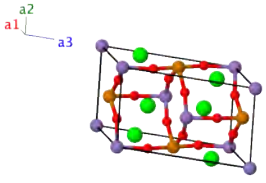
Other compounds with this structure

- Cs_2RbDyF_6

- This is the ground state structure of the double perovskite Sr_2MnTeO_6. Above 250 °C, it transforms into the Sr_2NiTeO_6 structure. Above 550 °C it transforms into the Sr_2NiWO_6 structure, and above 675 °C it transforms into the cubic perovskite $E2_1$ structure. (Ortega-San Martin, 2004).
- (Ortega-San Martin, 2004) give the structure in the $P2_1/n$ setting of space group #14. We used FINDSYM to transform this to the standard $P2_1/c$ setting, which uses a different set of basis vectors.

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	Mn
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	$(2a)$	Mn

$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(a + c\cos\beta)\;\hat{\mathbf{x}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	$(2d)$	Te
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$	$(2d)$	Te
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3a + z_3c\cos\beta)\;\hat{\mathbf{x}} + y_3b\,\hat{\mathbf{y}} + z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_3a - z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_7$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3a - z_3c\cos\beta)\;\hat{\mathbf{x}} - y_3b\,\hat{\mathbf{y}} - z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_8$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_3a + z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_9$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4a + z_4c\cos\beta)\;\hat{\mathbf{x}} + y_4b\,\hat{\mathbf{y}} + z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_{10}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_4a - z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_{11}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4a - z_4c\cos\beta)\;\hat{\mathbf{x}} - y_4b\,\hat{\mathbf{y}} - z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_{12}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_4a + z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_{13}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5a + z_5c\cos\beta)\;\hat{\mathbf{x}} + y_5b\,\hat{\mathbf{y}} + z_5c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O III
$\mathbf{B}_{14}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_5a - z_5c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O III
$\mathbf{B}_{15}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5a - z_5c\cos\beta)\;\hat{\mathbf{x}} - y_5b\,\hat{\mathbf{y}} - z_5c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O III
$\mathbf{B}_{16}$	=	$x_5\,\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_5a + z_5c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O III
$\mathbf{B}_{17}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6a + z_6c\cos\beta)\;\hat{\mathbf{x}} + y_6b\,\hat{\mathbf{y}} + z_6c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Sr
$\mathbf{B}_{18}$	=	$-x_6\,\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_6a - z_6c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Sr
$\mathbf{B}_{19}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6a - z_6c\cos\beta)\;\hat{\mathbf{x}} - y_6b\,\hat{\mathbf{y}} - z_6c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Sr
$\mathbf{B}_{20}$	=	$x_6\,\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_6a + z_6c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Sr

References:

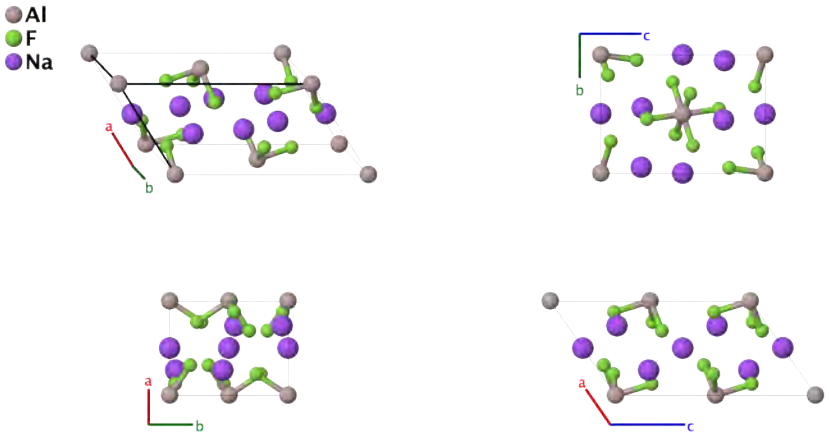
- L. Ortega-San Martin, J. P. Chapman, E. Hernández-Bocanegra, M. Insausti, M. I. Arriortua, and T. Rojo, *Structural phase transitions in the ordered double perovskite Sr_2MnTeO_6*, J. Phys.: Condens. Matter **16**, 3879–3888 (2004), doi:10.1088/0953-8984/16/23/008.

Geometry files:

- CIF: pp. 1552
- POSCAR: pp. 1552

Cryolite (Na_3AlF_6, $J2_6$) Structure: AB6C3_mP20_14_a_3e_de

http://aflow.org/prototype-encyclopedia/AB6C3_mP20_14_a_3e_de

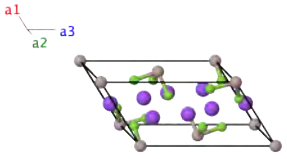


Prototype	:	AlF_6Na_3
AFLOW prototype label	:	AB6C3_mP20_14_a_3e_de
***Strukturbericht* designation**	:	$J2_6$
Pearson symbol	:	mP20
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=AB6C3_mP20_14_a_3e_de --params=`$a, b/a, c/a, \beta, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6$

- Cryolite undergoes a phase transition to an orthorhombic *Immm* phase at 890 K. Here we use structural data taken at 295 K.
- (Yang, 1993) gives the Wyckoff positions for the orientation $P2_1/n$ of space group #14. We used FINDSYM to change this to the standard $P2_1/c$ orientation. This included a change in the primitive vectors and the shift of the sodium atom on the ($2b$) site to the ($2d$) site.

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($2a$)	Al
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	($2a$)	Al

$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(a + c\cos\beta)\;\hat{\mathbf{x}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	(2*d*)	Na I
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$	(2*d*)	Na I
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3a + z_3c\cos\beta)\;\hat{\mathbf{x}} + y_3b\,\hat{\mathbf{y}} + z_3c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	F I
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_3a - z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	F I
$\mathbf{B}_7$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3a - z_3c\cos\beta)\;\hat{\mathbf{x}} - y_3b\,\hat{\mathbf{y}} - z_3c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	F I
$\mathbf{B}_8$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_3a + z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	F I
$\mathbf{B}_9$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4a + z_4c\cos\beta)\;\hat{\mathbf{x}} + y_4b\,\hat{\mathbf{y}} + z_4c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	F II
$\mathbf{B}_{10}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_4a - z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	F II
$\mathbf{B}_{11}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4a - z_4c\cos\beta)\;\hat{\mathbf{x}} - y_4b\,\hat{\mathbf{y}} - z_4c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	F II
$\mathbf{B}_{12}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_4a + z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	F II
$\mathbf{B}_{13}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5a + z_5c\cos\beta)\;\hat{\mathbf{x}} + y_5b\,\hat{\mathbf{y}} + z_5c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	F III
$\mathbf{B}_{14}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_5a - z_5c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	F III
$\mathbf{B}_{15}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5a - z_5c\cos\beta)\;\hat{\mathbf{x}} - y_5b\,\hat{\mathbf{y}} - z_5c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	F III
$\mathbf{B}_{16}$	=	$x_5\,\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_5a + z_5c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	F III
$\mathbf{B}_{17}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6a + z_6c\cos\beta)\;\hat{\mathbf{x}} + y_6b\,\hat{\mathbf{y}} + z_6c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Na II
$\mathbf{B}_{18}$	=	$-x_6\,\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_6a - z_6c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Na II
$\mathbf{B}_{19}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6a - z_6c\cos\beta)\;\hat{\mathbf{x}} - y_6b\,\hat{\mathbf{y}} - z_6c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Na II
$\mathbf{B}_{20}$	=	$x_6\,\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_6a + z_6c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Na II

References:

- H. Yang, S. Ghose, and D. M. Hatch, *Ferroelastic phase transition in cryolite, Na_3AlF_6, a mixed fluoride perovskite: High temperature single crystal X-ray diffraction study and symmetry analysis of the transition mechanism*, Phys. Chem. Miner. **19**, 528–544 (1993), doi:10.1007/BF00203053.

Found in:

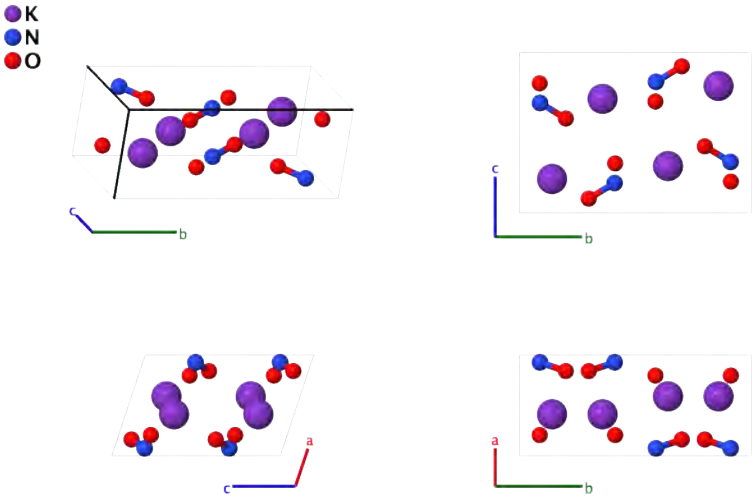
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1552
- POSCAR: pp. 1553

KNO_2 III Structure: ABC2_mP16_14_e_e_2e

http://aflow.org/prototype-encyclopedia/ABC2_mP16_14_e_e_2e.KNO2

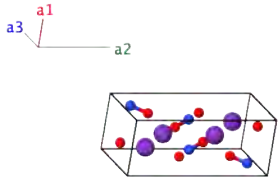


Prototype	:	KNO_2
AFLOW prototype label	:	ABC2_mP16_14_e_e_2e
***Strukturbericht* designation**	:	None
Pearson symbol	:	mP16
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=ABC2_mP16_14_e_e_2e --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4$

- This is *not* the KNO_2 structure found by (Ziegler, 1936) and given the *Strukturbericht* designation $F5_{11}$. That structure is now considered obsolete.

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	K
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2}+y_1\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_1 a - z_1 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	K

$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$(-x_1 a - z_1 c\cos\beta)\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} - z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	K
$\mathbf{B}_4$	=	$x_1\,\mathbf{a}_1 + \left(\frac{1}{2} - y_1\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_1 a + z_1 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_1\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	K
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	N
$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_2 a - z_2 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	N
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} - z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	N
$\mathbf{B}_8$	=	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_2 a + z_2 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	N
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_3 a - z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_{12}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_4 a - z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_{16}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O II

References:

- N. Onoda-Yamamuro, H. Honda, R. Ikeda, O. Yamamuro, T. Matsuo, K. Oikawa, T. Kamiyama, and F. Izumi, *Neutron powder diffraction study of the low-temperature phases of KNO_2*, J. Phys.: Condens. Matter **10**, 3341–3351 (1998), doi:10.1088/0953-8984/10/15/011.
- G. E. Ziegler, *The Crystal Structure of Potassium Nitrite, KNO_2*, Zeitschrift für Kristallographie - Crystalline Materials **34**, 491–499 (1936), doi:10.1524/zkri.1936.94.1.491.

Found in:

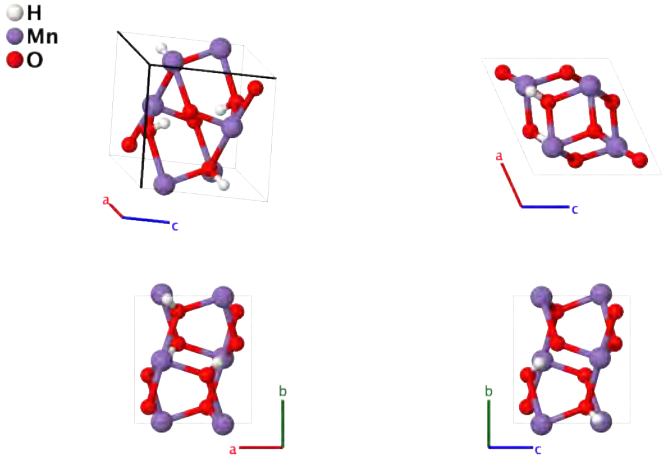
- C. Duan, W. N. Mei, R. W. Smith, J. Liu, M. M. Ossowski, and J. R. Hardy, *Order-disorder phase transitions in KNO_2, $CsNO_2$, and $TlNO_2$ crystals: A molecular dynamics study*, Phys. Rev. B **63**, 144105 (2001), doi:10.1103/PhysRevB.63.144105.

Geometry files:

- CIF: pp. 1553
- POSCAR: pp. 1553

Manganite (γ-MnO(OH), $E0_6$) Structure: ABC2_mP16_14_e_e_2e

http://aflow.org/prototype-encyclopedia/ABC2_mP16_14_e_e_2e

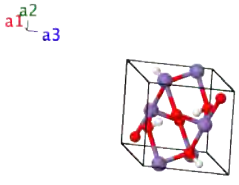


Prototype	:	$MnHO_2$
AFLOW prototype label	:	ABC2_mP16_14_e_e_2e
***Strukturbericht* designation**	:	$E0_6$
Pearson symbol	:	mP16
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=ABC2_mP16_14_e_e_2e --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4$

- (Buerger, 1936) originally determined the positions of the manganese and oxygen atoms using the $B2_1/d$ setting of space group #14. This structure was given the *Strukturbericht* designation $E0_6$ in (Gottfried, 1938). Much later, (Kohler, 1997) found the positions of the hydrogen atoms, and presented the data in the standard $P2_1/c$ setting of space group #14.

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$ =	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H

$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2}+y_1\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_1 a - z_1 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$(-x_1 a - z_1 c\cos\beta)\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} - z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H
$\mathbf{B}_4$	=	$x_1\,\mathbf{a}_1 + \left(\frac{1}{2}-y_1\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_1 a + z_1 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Mn
$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2}+y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_2 a - z_2 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Mn
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} - z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Mn
$\mathbf{B}_8$	=	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2}-y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_2 a + z_2 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Mn
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2}+y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_3 a - z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_{12}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2}-y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2}+y_4\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_4 a - z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_{16}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2}-y_4\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O II

References:

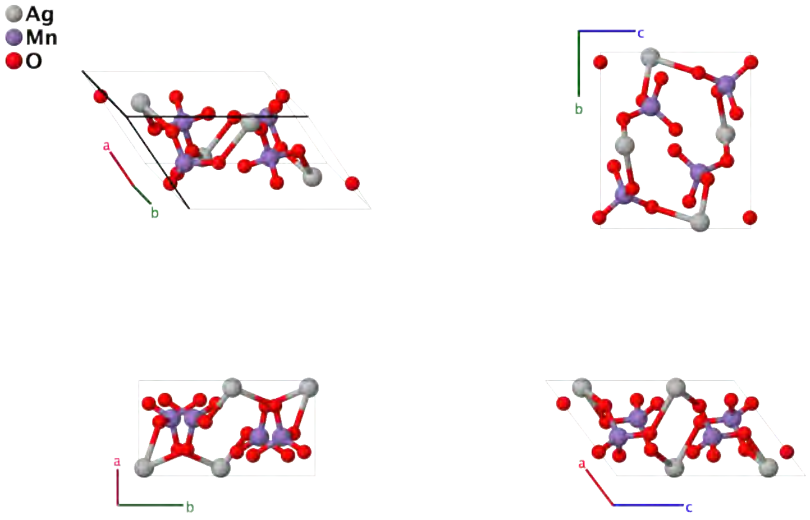
- T. Kohler, T. Armbruster, and E. Libowitzky, *Hydrogen Bonding and Jahn-Teller Distortion in Groutite, α-MnOOH, and Manganite, γ-MnOOH, and Their Relations to the Manganese Dioxides Ramsdellite and Pyrolusite*, J. Solid State Chem. **133**, 486–500 (1997), doi:10.1006/jssc.1997.7516.
- C. Gottfried, ed., *Strukturbericht Band IV 1936* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1938).
- M. J. Buerger, *The Symmetry and Crystal Structure of Manganite, Mn(OH)O*, Zeitschrift für Kristallographie - Crystalline Materials **95**, 163–174 (1936), doi:10.1524/zkri.1936.95.1.163.

Geometry files:

- CIF: pp. 1553
- POSCAR: pp. 1554

$AgMnO_4$ ($H0_9$) Structure: ABC4_mP24_14_e_e_4e

http://aflow.org/prototype-encyclopedia/ABC4_mP24_14_e_e_4e



Prototype	:	$AgMnO_4$
AFLOW prototype label	:	ABC4_mP24_14_e_e_4e
***Strukturbericht* designation**	:	$H0_9$
Pearson symbol	:	mP24
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=ABC4_mP24_14_e_e_4e --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6$

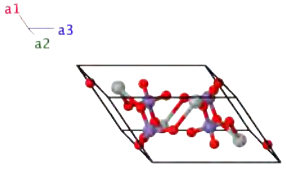
Other compounds with this structure

- $CaSO_4$, $SrCrO_4$, $InVO_4$, $CsSO_4$, and $(NH_4)SO_4$

- The structure of $AgMnO_4$ was originally determined by (Sasvári, 1938) and refined by (Boonstra, 1968). In particular, Boonstra found that the actual positions of the Ag and Mn atoms were switched compared to the structure found by Sasvári. As the lattice constants remain approximately the same and there was no change in the space group or the occupied Wyckoff positions, we use the Boonstra structure to define *Strukturbericht* designation $H0_9$.
- Boonstra gave the structure in setting $P2_1/n$ of space group #14. We used FINDSYM to translate the structure to the standard $P2_1/c$ setting. This included a redefinition of the primitive vectors of the lattice.

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Ag
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2}+y_1\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_1 a - z_1 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Ag
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$(-x_1 a - z_1 c\cos\beta)\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} - z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Ag
$\mathbf{B}_4$	=	$x_1\,\mathbf{a}_1 + \left(\frac{1}{2}-y_1\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_1 a + z_1 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Ag
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Mn
$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2}+y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_2 a - z_2 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Mn
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} - z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Mn
$\mathbf{B}_8$	=	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2}-y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_2 a + z_2 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Mn
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2}+y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_3 a - z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_{12}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2}-y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2}+y_4\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_4 a - z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_{16}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2}-y_4\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O III
$\mathbf{B}_{18}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2}+y_5\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_5 a - z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O III
$\mathbf{B}_{19}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O III
$\mathbf{B}_{20}$	=	$x_5\,\mathbf{a}_1 + \left(\frac{1}{2}-y_5\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O III
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O IV

$$\mathbf{B}_{22} = -x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_6 a - z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_6\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_6\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O IV}$$

$$\mathbf{B}_{23} = -x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3 = (-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O IV}$$

$$\mathbf{B}_{24} = x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_6\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O IV}$$

References:

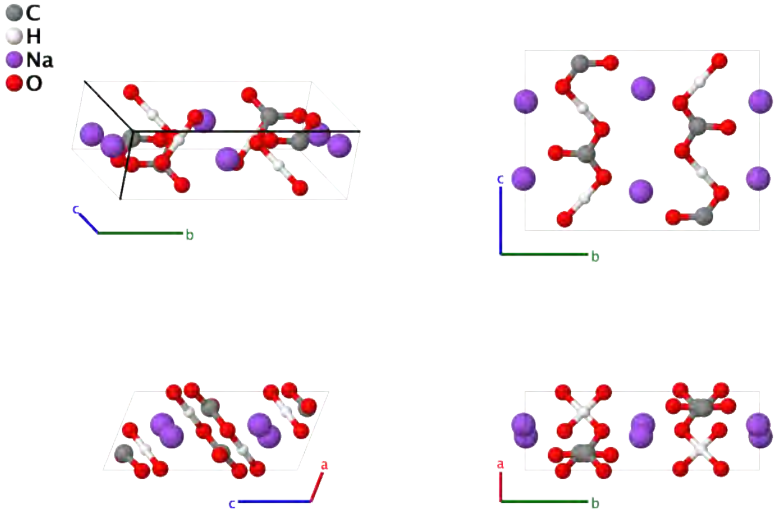
- E. G. Boonstra, *The Crystal Structure of Silver Permanganate*, Acta Crystallogr. Sect. B Struct. Sci. **24**, 1053–1062 (1968), doi:10.1107/S0567740868003699.
- K. Sasvári, *Die Struktur des Silberpermanganats $AgMnO_4$*, Zeitschrift für Kristallographie - Crystalline Materials **99**, 9–15 (1938), doi:10.1524/zkri.1938.99.1.9.

Geometry files:

- CIF: pp. 1554
- POSCAR: pp. 1554

Nahcolite ($NaHCO_3$, $G0_{12}$) Structure: ABCD3_mP24_14_e_e_e_3e

http://aflow.org/prototype-encyclopedia/ABCD3_mP24_14_e_e_e_3e



Prototype	:	$CHNaO_3$
AFLOW prototype label	:	ABCD3_mP24_14_e_e_e_3e
***Strukturbericht* designation**	:	$G0_{12}$
Pearson symbol	:	mP24
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=ABCD3_mP24_14_e_e_e_3e --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6$

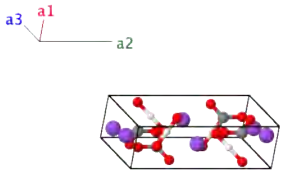
Other compounds with this structure

- $KHCO_3$

- This is more commonly known as sodium bicarbonate or baking soda.
- (Sass, 1962) do not locate the hydrogen atoms.
- (Zachariasen, 1933) applied chemical and crystallographic reasoning and placed the hydrogen atoms halfway between each O II atom and the nearest O III atom. We use this structure, while realizing that the hydrogen positions are only approximate.

Simple Monoclinic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= a\,\hat{\mathbf{x}}\\ \mathbf{a}_2 &= b\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	$(x_1a + z_1c\cos\beta)\,\hat{\mathbf{x}} + y_1b\,\hat{\mathbf{y}} + z_1c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	C
$\mathbf{B}_2$	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2}+y_1\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_1\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta - x_1a - z_1c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	C
$\mathbf{B}_3$	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	$(-x_1a - z_1c\cos\beta)\,\hat{\mathbf{x}} - y_1b\,\hat{\mathbf{y}} - z_1c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	C
$\mathbf{B}_4$	$x_1\,\mathbf{a}_1 + \left(\frac{1}{2}-y_1\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_1\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_1a + z_1c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	C
$\mathbf{B}_5$	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	$(x_2a + z_2c\cos\beta)\,\hat{\mathbf{x}} + y_2b\,\hat{\mathbf{y}} + z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H
$\mathbf{B}_6$	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2}+y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_2\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta - x_2a - z_2c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H
$\mathbf{B}_7$	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	$(-x_2a - z_2c\cos\beta)\,\hat{\mathbf{x}} - y_2b\,\hat{\mathbf{y}} - z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H
$\mathbf{B}_8$	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2}-y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_2a + z_2c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H
$\mathbf{B}_9$	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$(x_3a + z_3c\cos\beta)\,\hat{\mathbf{x}} + y_3b\,\hat{\mathbf{y}} + z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Na
$\mathbf{B}_{10}$	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2}+y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_3\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta - x_3a - z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Na
$\mathbf{B}_{11}$	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	$(-x_3a - z_3c\cos\beta)\,\hat{\mathbf{x}} - y_3b\,\hat{\mathbf{y}} - z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Na
$\mathbf{B}_{12}$	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2}-y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_3a + z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Na
$\mathbf{B}_{13}$	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$(x_4a + z_4c\cos\beta)\,\hat{\mathbf{x}} + y_4b\,\hat{\mathbf{y}} + z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_{14}$	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2}+y_4\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_4\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta - x_4a - z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_{15}$	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	$(-x_4a - z_4c\cos\beta)\,\hat{\mathbf{x}} - y_4b\,\hat{\mathbf{y}} - z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_{16}$	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2}-y_4\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta + x_4a + z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_{17}$	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	$(x_5a + z_5c\cos\beta)\,\hat{\mathbf{x}} + y_5b\,\hat{\mathbf{y}} + z_5c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O II

$$\mathbf{B}_{18} = -x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_5 a - z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_5\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_5\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O II}$$

$$\mathbf{B}_{19} = -x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3 = (-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O II}$$

$$\mathbf{B}_{20} = x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_5\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_5\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O II}$$

$$\mathbf{B}_{21} = x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 = (x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O III}$$

$$\mathbf{B}_{22} = -x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_6 a - z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_6\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_6\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O III}$$

$$\mathbf{B}_{23} = -x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3 = (-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O III}$$

$$\mathbf{B}_{24} = x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_6\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{O III}$$

References:

- R. L. Sass and R. F. Scheuerman, *The Crystal Structure of Sodium Bicarbonate*, Acta Cryst. **15**, 77–81 (1962), doi:10.1107/S0365110X62000158.
- W. H. Zachariasen, *The Crystal Lattice of Sodium Bicarbonate, $NaHCO_3$*, J. Chem. Phys. **1**, 634–639 (1933), doi:10.1063/1.1749342.

Found in:

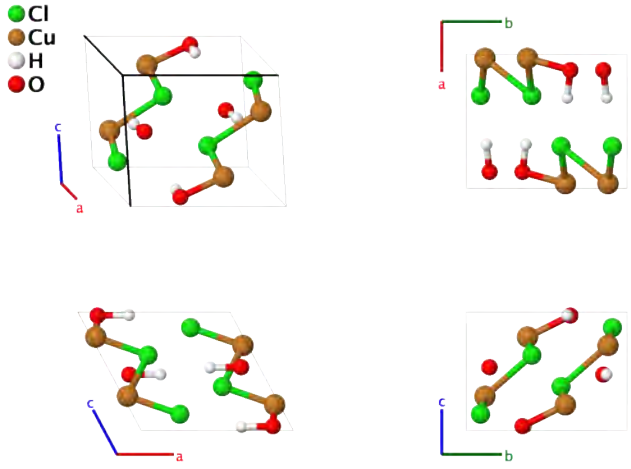
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1554
- POSCAR: pp. 1555

Cu(OH)Cl Structure: ABCD_mP16_14_e_e_e_e

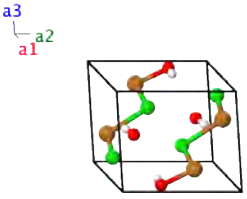
http://aflow.org/prototype-encyclopedia/ABCD_mP16_14_e_e_e_e



Prototype	:	ClCuHO
AFLOW prototype label	:	ABCD_mP16_14_e_e_e_e
***Strukturbericht* designation**	:	None
Pearson symbol	:	mP16
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=ABCD_mP16_14_e_e_e_e --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4$

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$ =	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$ =	$(x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cl
$\mathbf{B}_2$ =	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2} + y_1\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$ =	$\left(\frac{1}{2}c\cos\beta - x_1 a - z_1 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_1\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cl
$\mathbf{B}_3$ =	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$ =	$(-x_1 a - z_1 c\cos\beta)\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} - z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cl
$\mathbf{B}_4$ =	$x_1\,\mathbf{a}_1 + \left(\frac{1}{2} - y_1\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$ =	$\left(\frac{1}{2}c\cos\beta + x_1 a + z_1 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_1\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cl

$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2a + z_2c\cos\beta)\,\hat{\mathbf{x}} + y_2b\,\hat{\mathbf{y}} + z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cu
$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_2a - z_2c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cu
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2a - z_2c\cos\beta)\,\hat{\mathbf{x}} - y_2b\,\hat{\mathbf{y}} - z_2c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cu
$\mathbf{B}_8$	=	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_2a + z_2c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cu
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3a + z_3c\cos\beta)\,\hat{\mathbf{x}} + y_3b\,\hat{\mathbf{y}} + z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_3a - z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3a - z_3c\cos\beta)\,\hat{\mathbf{x}} - y_3b\,\hat{\mathbf{y}} - z_3c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H
$\mathbf{B}_{12}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_3a + z_3c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4a + z_4c\cos\beta)\,\hat{\mathbf{x}} + y_4b\,\hat{\mathbf{y}} + z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_4a - z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4a - z_4c\cos\beta)\,\hat{\mathbf{x}} - y_4b\,\hat{\mathbf{y}} - z_4c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O
$\mathbf{B}_{16}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_4a + z_4c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O

References:

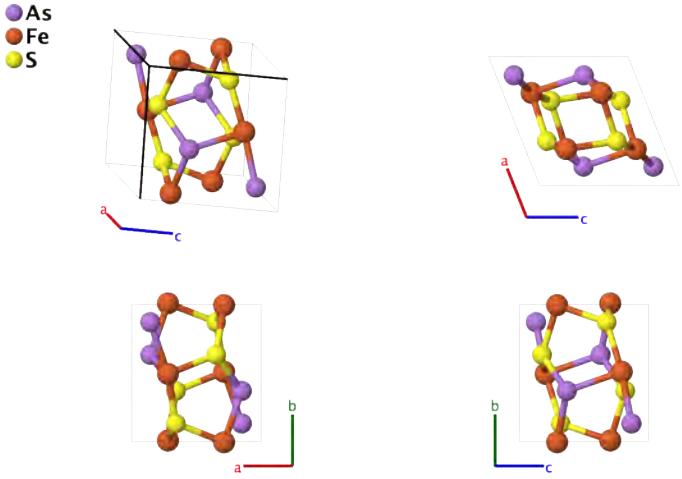
- Y. Cudennec, A. Riou, Y. Gérault, and A. Lecerf, *Synthesis and Crystal Structures of Cd(OH)Cl and Cu(OH)Cl and Relationship to Brucite Type*, J. Solid State Chem. **151**, 308–312 (2000), doi:10.1006/jssc.2000.8659.

Geometry files:

- CIF: pp. 1555
- POSCAR: pp. 1555

Arsenopyrite (FeAsS, $E0_7$) Structure: ABC_mP12_14_e_e_e

http://aflow.org/prototype-encyclopedia/ABC_mP12_14_e_e_e



Prototype	:	AsFeS
AFLOW prototype label	:	ABC_mP12_14_e_e_e
***Strukturbericht* designation**	:	$E0_7$
Pearson symbol	:	mP12
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=ABC_mP12_14_e_e_e --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3$

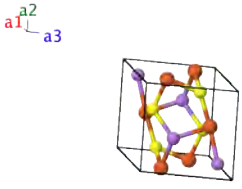
Other compounds with this structure

- FeSbS (gudmundite), $CoSb_2$, IrN_2, MoO_2 (tugarinovite), VO_2, $IrAs_2$ (iridarsenite), $CoAs_2$ (clinosafflorite), and HfO_2

- This is very similar to *C*43 baddeleyite, ZrO_2.

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\sin\beta\,\hat{\mathbf{z}}$	($4e$)	As

$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2}+y_1\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_1 a - z_1 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$(-x_1 a - z_1 c\cos\beta)\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} - z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As
$\mathbf{B}_4$	=	$x_1\,\mathbf{a}_1 + \left(\frac{1}{2}-y_1\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_1 a + z_1 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Fe
$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2}+y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_2 a - z_2 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Fe
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} - z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Fe
$\mathbf{B}_8$	=	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2}-y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_2 a + z_2 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Fe
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2}+y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_3 a - z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S
$\mathbf{B}_{12}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2}-y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S

References:

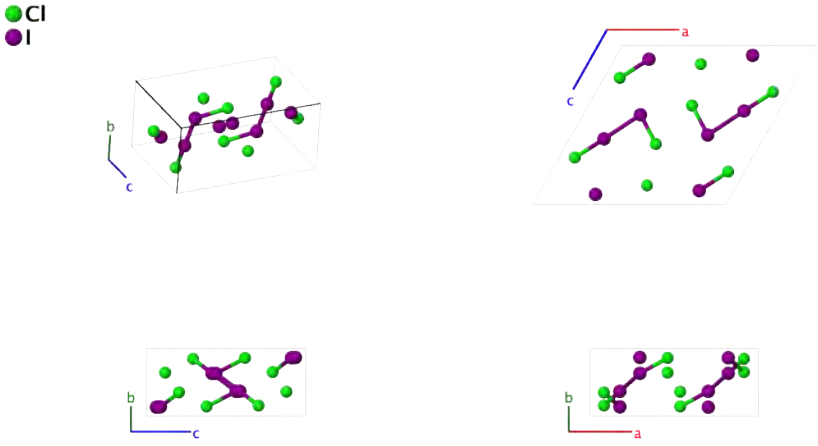
- L. Bindi, Y. Moëlo, P. Léone, and M. Suchaud, *Stoichiometric Arsenopyrite, FeAsS, from La Roche-Balue Quarry, Loire-Atlantique, France: Crystal Structure And Mössbauer Study*, Can. Mineral. **50**, 471–479 (2012), doi:10.3749/canmin.50.2.471.

Geometry files:

- CIF: pp. 1556
- POSCAR: pp. 1556

α-ICl Structure: AB_mP16_14_2e_2e

http://aflow.org/prototype-encyclopedia/AB_mP16_14_2e_2e.ICl

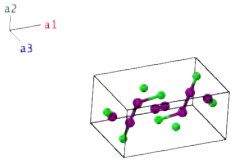


Prototype	:	ClI
AFLOW prototype label	:	AB_mP16_14_2e_2e
***Strukturbericht* designation**	:	None
Pearson symbol	:	mP16
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=AB_mP16_14_2e_2e --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4$

- LiAs (AB_mP16_14_2e_2e) and α-ICl (AB_mP16_14_2e_2e) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters (`--params`) specified in their corresponding CIF files.

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$ =	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cl I
$\mathbf{B}_2$ =	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2}+y_1\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_1 a - z_1 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cl I
$\mathbf{B}_3$ =	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$(-x_1 a - z_1 c\cos\beta)\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} - z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cl I

$\mathbf{B}_4$	=	$x_1\,\mathbf{a}_1 + \left(\frac{1}{2} - y_1\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_1 a + z_1 c\cos\beta\right)\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_1\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Cl I
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Cl II
$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_2 a - z_2 c\cos\beta\right)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Cl II
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} - z_2 c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Cl II
$\mathbf{B}_8$	=	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_2 a + z_2 c\cos\beta\right)\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Cl II
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	I I
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_3 a - z_3 c\cos\beta\right)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	I I
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	I I
$\mathbf{B}_{12}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	I I
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	I II
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_4 a - z_4 c\cos\beta\right)\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	I II
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	I II
$\mathbf{B}_{16}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	I II

References:

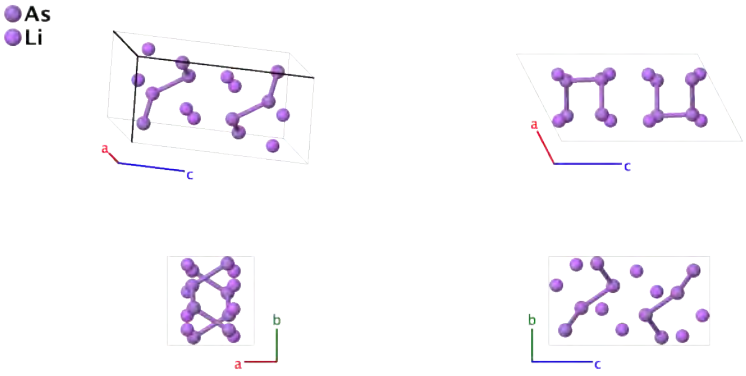
- K. H. Boswijk, J. van der Heide, A. Vos, and E. H. Wiebenga, *The crystal structure of α-ICl*, Acta Cryst. **9**, 274–277 (1956), doi:10.1107/S0365110X56000760.

Geometry files:

- CIF: pp. 1556
- POSCAR: pp. 1556

LiAs Structure: AB_mP16_14_2e_2e

http://aflow.org/prototype-encyclopedia/AB_mP16_14_2e_2e



Prototype	:	AsLi
AFLOW prototype label	:	AB_mP16_14_2e_2e
***Strukturbericht* designation**	:	None
Pearson symbol	:	mP16
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=AB_mP16_14_2e_2e --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4$

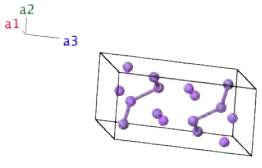
Other compounds with this structure

- NaSb and KSb

- LiAs (AB_mP16_14_2e_2e) and α-ICl (AB_mP16_14_2e_2e) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters (`--params`) specified in their corresponding CIF files.

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	$(x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As I
$\mathbf{B}_2$	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2} + y_1\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	$\left(\frac{1}{2}c\cos\beta - x_1 a - z_1 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_1\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As I
$\mathbf{B}_3$	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	$(-x_1 a - z_1 c\cos\beta)\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} - z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As I

$\mathbf{B}_4$	=	$x_1\,\mathbf{a}_1 + \left(\frac{1}{2} - y_1\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_1 a + z_1 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_1\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As I
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As II
$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_2 a - z_2 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As II
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} - z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As II
$\mathbf{B}_8$	=	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_2 a + z_2 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As II
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Li I
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_3 a - z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Li I
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Li I
$\mathbf{B}_{12}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Li I
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Li II
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_4 a - z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Li II
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Li II
$\mathbf{B}_{16}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Li II

References:

- D. T. Cromer, *The Crystal Structure of LiAs*, Acta Cryst. **12**, 36–41 (1959), doi:10.1107/S0365110X59000111.

Found in:

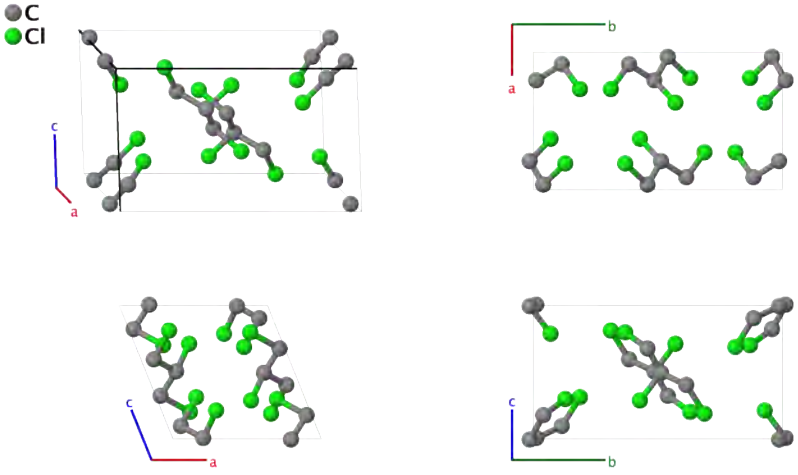
- P. Villars and L. Calvert, *Pearson's Handbook of Crystallographic Data for Intermetallic Phases* (ASM International, Materials Park, OH, 1991), 2nd edn.

Geometry files:

- CIF: pp. 1556
- POSCAR: pp. 1557

ϵ-1,2,3,4,5,6-Hexachlorocyclohexane (C_6Cl_6) Structure: AB_mP24_14_3e_3e

http://aflow.org/prototype-encyclopedia/AB_mP24_14_3e_3e

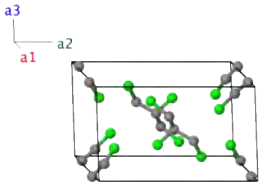


Prototype	:	C_6Cl_6
AFLOW prototype label	:	AB_mP24_14_3e_3e
***Strukturbericht* designation**	:	None
Pearson symbol	:	mP24
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=AB_mP24_14_3e_3e --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6$

- This unit cell contains two benzene molecules, with every hydrogen replaced by a chlorine atom, causing considerable distortion. To see this it will be necessary to use a supercell view of at least 212.

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$ =	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$ =	$(x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	C I
$\mathbf{B}_2$ =	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2}+y_1\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_1\right)\mathbf{a}_3$ =	$\left(\frac{1}{2}c\cos\beta - x_1 a - z_1 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	C I

$$\mathbf{B}_3 = -x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3 = (-x_1 a - z_1 c\cos\beta)\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} - z_1 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{C I}$$

$$\mathbf{B}_4 = x_1\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_1\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_1\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_1 a + z_1 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_1\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_1\right) c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{C I}$$

$$\mathbf{B}_5 = x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3 = (x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{C II}$$

$$\mathbf{B}_6 = -x_2\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_2\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_2\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_2 a - z_2 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_2\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_2\right) c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{C II}$$

$$\mathbf{B}_7 = -x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3 = (-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} - z_2 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{C II}$$

$$\mathbf{B}_8 = x_2\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_2\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_2\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_2 a + z_2 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_2\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_2\right) c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{C II}$$

$$\mathbf{B}_9 = x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3 = (x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{C III}$$

$$\mathbf{B}_{10} = -x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_3 a - z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_3\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_3\right) c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{C III}$$

$$\mathbf{B}_{11} = -x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3 = (-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{C III}$$

$$\mathbf{B}_{12} = x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_3\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_3\right) c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{C III}$$

$$\mathbf{B}_{13} = x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3 = (x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cl I}$$

$$\mathbf{B}_{14} = -x_4\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_4 a - z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_4\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_4\right) c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cl I}$$

$$\mathbf{B}_{15} = -x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3 = (-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cl I}$$

$$\mathbf{B}_{16} = x_4\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_4\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_4\right) c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cl I}$$

$$\mathbf{B}_{17} = x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3 = (x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cl II}$$

$$\mathbf{B}_{18} = -x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_5 a - z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_5\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_5\right) c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cl II}$$

$$\mathbf{B}_{19} = -x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3 = (-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cl II}$$

$$\mathbf{B}_{20} = x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_5\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_5\right) c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cl II}$$

$$\mathbf{B}_{21} = x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 = (x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cl III}$$

$$\mathbf{B}_{22} = -x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_6 a - z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_6\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_6\right) c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cl III}$$

$$\mathbf{B}_{23} = -x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3 = (-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cl III}$$

$$\mathbf{B}_{24} = x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_6\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right) c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{Cl III}$$

References:

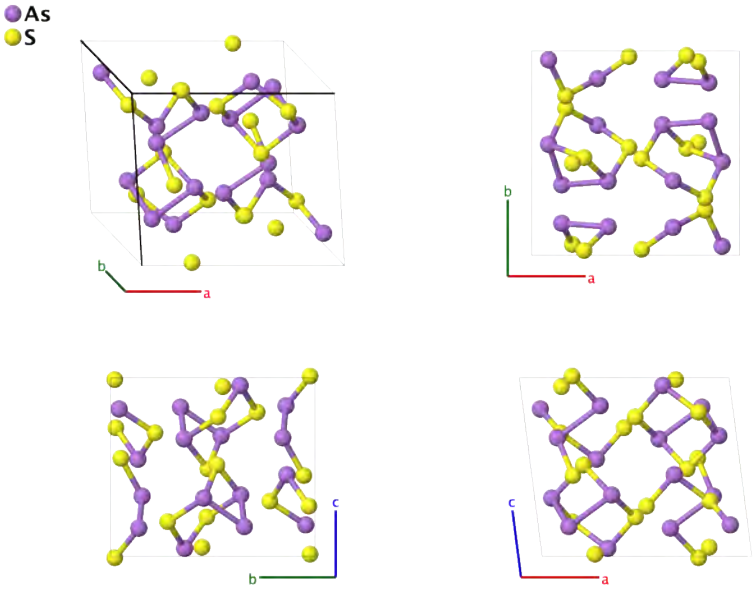
- N. Norman, *The Crystal Structure of the Epsilon Isomer of 1,2,3,4,5,6-Hexachloro*cyclohexane, Acta Chem. Scand. **4**, 251–259 (1950), doi:10.3891/acta.chem.scand.04-0251.

Geometry files:

- CIF: pp. 1557
- POSCAR: pp. 1557

Pararealgar (AsS) Structure: AB_mP32_14_4e_4e

http://aflow.org/prototype-encyclopedia/AB_mP32_14_4e_4e.pararealgar



Prototype	:	AsS
AFLOW prototype label	:	AB_mP32_14_4e_4e
***Strukturbericht* designation**	:	None
Pearson symbol	:	mP32
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=AB_mP32_14_4e_4e --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8$

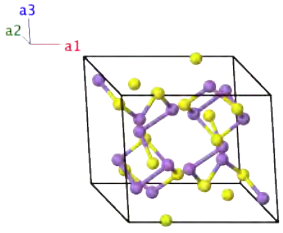
Other compounds with this structure

- AsSe

- Realgar (B_l) transforms to pararealgar on exposure to visible light with wavelengths in the range 500-670 nm (Douglass, 1992).

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As I
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2}+y_1\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_1 a - z_1 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As I
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$(-x_1 a - z_1 c\cos\beta)\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} - z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As I
$\mathbf{B}_4$	=	$x_1\,\mathbf{a}_1 + \left(\frac{1}{2}-y_1\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_1 a + z_1 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As I
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As II
$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2}+y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_2 a - z_2 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As II
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} - z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As II
$\mathbf{B}_8$	=	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2}-y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_2 a + z_2 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As II
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As III
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2}+y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_3 a - z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As III
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As III
$\mathbf{B}_{12}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2}-y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As III
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As IV
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2}+y_4\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_4 a - z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As IV
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As IV
$\mathbf{B}_{16}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2}-y_4\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As IV
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S I
$\mathbf{B}_{18}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2}+y_5\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_5 a - z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S I
$\mathbf{B}_{19}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S I
$\mathbf{B}_{20}$	=	$x_5\,\mathbf{a}_1 + \left(\frac{1}{2}-y_5\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S I
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S II

$\mathbf{B}_{22}$	$=$	$-x_6\,\mathbf{a}_1+\left(\frac{1}{2}+y_6\right)\mathbf{a}_2+\left(\frac{1}{2}-z_6\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}c\cos\beta-x_6a-z_6c\cos\beta\right)\hat{\mathbf{x}}+\left(\frac{1}{2}+y_6\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S II
$\mathbf{B}_{23}$	$=$	$-x_6\,\mathbf{a}_1-y_6\,\mathbf{a}_2-z_6\,\mathbf{a}_3$	$=$	$(-x_6a-z_6c\cos\beta)\,\hat{\mathbf{x}}-y_6b\,\hat{\mathbf{y}}-z_6c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S II
$\mathbf{B}_{24}$	$=$	$x_6\,\mathbf{a}_1+\left(\frac{1}{2}-y_6\right)\mathbf{a}_2+\left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}c\cos\beta+x_6a+z_6c\cos\beta\right)\hat{\mathbf{x}}+\left(\frac{1}{2}-y_6\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S II
$\mathbf{B}_{25}$	$=$	$x_7\,\mathbf{a}_1+y_7\,\mathbf{a}_2+z_7\,\mathbf{a}_3$	$=$	$(x_7a+z_7c\cos\beta)\,\hat{\mathbf{x}}+y_7b\,\hat{\mathbf{y}}+z_7c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S III
$\mathbf{B}_{26}$	$=$	$-x_7\,\mathbf{a}_1+\left(\frac{1}{2}+y_7\right)\mathbf{a}_2+\left(\frac{1}{2}-z_7\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}c\cos\beta-x_7a-z_7c\cos\beta\right)\hat{\mathbf{x}}+\left(\frac{1}{2}+y_7\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_7\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S III
$\mathbf{B}_{27}$	$=$	$-x_7\,\mathbf{a}_1-y_7\,\mathbf{a}_2-z_7\,\mathbf{a}_3$	$=$	$(-x_7a-z_7c\cos\beta)\,\hat{\mathbf{x}}-y_7b\,\hat{\mathbf{y}}-z_7c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S III
$\mathbf{B}_{28}$	$=$	$x_7\,\mathbf{a}_1+\left(\frac{1}{2}-y_7\right)\mathbf{a}_2+\left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}c\cos\beta+x_7a+z_7c\cos\beta\right)\hat{\mathbf{x}}+\left(\frac{1}{2}-y_7\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_7\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S III
$\mathbf{B}_{29}$	$=$	$x_8\,\mathbf{a}_1+y_8\,\mathbf{a}_2+z_8\,\mathbf{a}_3$	$=$	$(x_8a+z_8c\cos\beta)\,\hat{\mathbf{x}}+y_8b\,\hat{\mathbf{y}}+z_8c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S IV
$\mathbf{B}_{30}$	$=$	$-x_8\,\mathbf{a}_1+\left(\frac{1}{2}+y_8\right)\mathbf{a}_2+\left(\frac{1}{2}-z_8\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}c\cos\beta-x_8a-z_8c\cos\beta\right)\hat{\mathbf{x}}+\left(\frac{1}{2}+y_8\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_8\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S IV
$\mathbf{B}_{31}$	$=$	$-x_8\,\mathbf{a}_1-y_8\,\mathbf{a}_2-z_8\,\mathbf{a}_3$	$=$	$(-x_8a-z_8c\cos\beta)\,\hat{\mathbf{x}}-y_8b\,\hat{\mathbf{y}}-z_8c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S IV
$\mathbf{B}_{32}$	$=$	$x_8\,\mathbf{a}_1+\left(\frac{1}{2}-y_8\right)\mathbf{a}_2+\left(\frac{1}{2}+z_8\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}c\cos\beta+x_8a+z_8c\cos\beta\right)\hat{\mathbf{x}}+\left(\frac{1}{2}-y_8\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_8\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S IV

References:

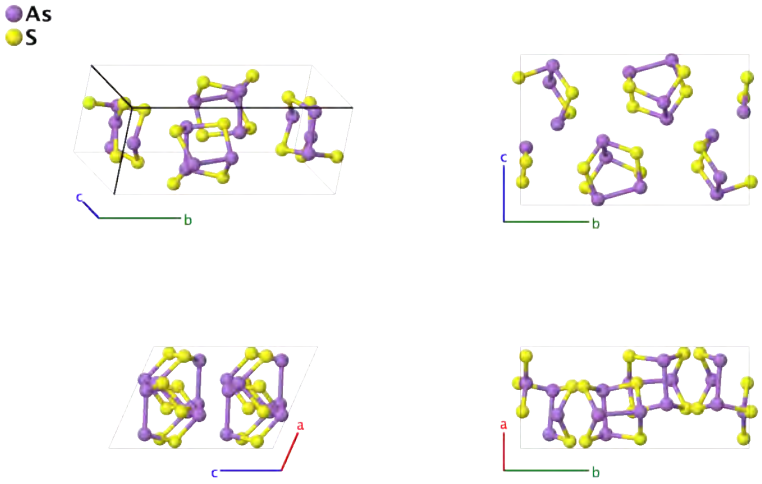
- P. Bonazzi, S. Menchetti, and G. Pratesi, *The crystal structure of pararealgar, As_4S_4*, Am. Mineral. **80**, 400–403 (1995), doi:10.2138/am-1995-3-422.
- D. L. Douglass, C. Shing, and G. Wang, *The light-induced alteration of realgar to pararealgar*, Am. Mineral. **77**, 1266–1274 (1992).

Geometry files:

- CIF: pp. 1557
- POSCAR: pp. 1558

Realgar (AsS, B_l) Structure: AB_mP32_14_4e_4e

http://aflow.org/prototype-encyclopedia/AB_mP32_14_4e_4e.realgar



Prototype	:	AsS
AFLOW prototype label	:	AB_mP32_14_4e_4e
***Strukturbericht* designation**	:	B_l
Pearson symbol	:	mP32
Space group number	:	14
Space group symbol	:	$P2_1/c$
AFLOW prototype command	:	`aflow --proto=AB_mP32_14_4e_4e --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8$

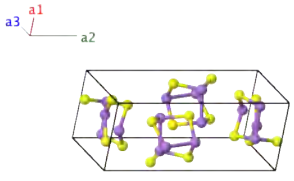
Other compounds with this structure

- AsSe

- Realgar transforms to pararealgar on exposure to visible light with wavelengths in the range 500-670 nm (Douglass, 1992).
- (Ito, 1952) give the crystal structure of realgar in the $P2_1/n$ setting of space group #14. We have used FINDSYM to transform this to the standard $P2_1/c$ setting. This involved both rotations and an origin shift.

Simple Monoclinic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As I
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2}+y_1\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_1 a - z_1 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As I
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$(-x_1 a - z_1 c\cos\beta)\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} - z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As I
$\mathbf{B}_4$	=	$x_1\,\mathbf{a}_1 + \left(\frac{1}{2}-y_1\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_1 a + z_1 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As I
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As II
$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2}+y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_2 a - z_2 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As II
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} - z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As II
$\mathbf{B}_8$	=	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2}-y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_2 a + z_2 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As II
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As III
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2}+y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_3 a - z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As III
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As III
$\mathbf{B}_{12}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2}-y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As III
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As IV
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2}+y_4\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_4 a - z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As IV
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As IV
$\mathbf{B}_{16}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2}-y_4\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	As IV
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S I
$\mathbf{B}_{18}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2}+y_5\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_5 a - z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}+y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S I
$\mathbf{B}_{19}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S I
$\mathbf{B}_{20}$	=	$x_5\,\mathbf{a}_1 + \left(\frac{1}{2}-y_5\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\hat{\mathbf{x}} + \left(\frac{1}{2}-y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S I
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S II

$$\mathbf{B}_{22} = -x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_6 a - z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_6\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_6\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{S II}$$

$$\mathbf{B}_{23} = -x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3 = (-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{S II}$$

$$\mathbf{B}_{24} = x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_6\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{S II}$$

$$\mathbf{B}_{25} = x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 = (x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{S III}$$

$$\mathbf{B}_{26} = -x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_7 a - z_7 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_7\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_7\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{S III}$$

$$\mathbf{B}_{27} = -x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3 = (-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{S III}$$

$$\mathbf{B}_{28} = x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_7 a + z_7 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_7\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_7\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{S III}$$

$$\mathbf{B}_{29} = x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 = (x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{S IV}$$

$$\mathbf{B}_{30} = -x_8\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_8 a - z_8 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_8\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_8\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{S IV}$$

$$\mathbf{B}_{31} = -x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3 = (-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{S IV}$$

$$\mathbf{B}_{32} = x_8\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_8 a + z_8 c\cos\beta\right)\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_8\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_8\right)c\sin\beta\,\hat{\mathbf{z}} \quad (4e) \quad \text{S IV}$$

References:

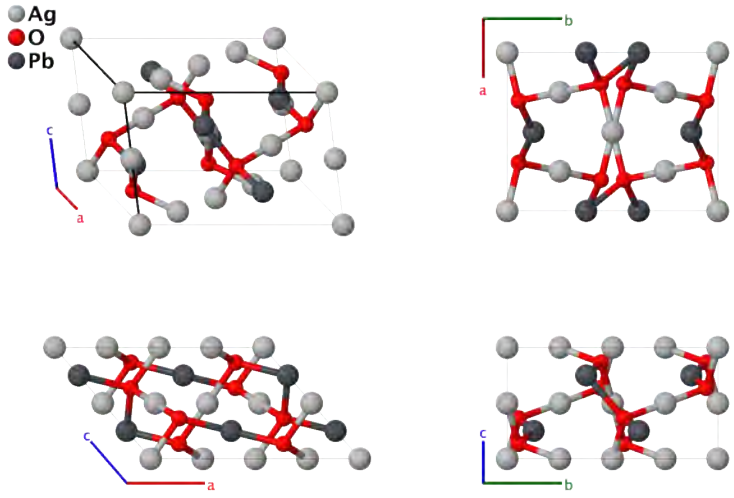
- T. Ito, N. Morimoto, and R. Sadanaga, *The Crystal Structure of Realgar*, Acta Cryst. **5**, 755–782 (1952), doi:10.1107/S0365110X52002112.
- D. L. Douglass, C. Shing, and G. Wang, *The light-induced alteration of realgar to pararealgar*, Am. Mineral. **77**, 1266–1274 (1992).

Geometry files:

- CIF: pp. 1558
- POSCAR: pp. 1558

Ag_2PbO_2 Structure: A2B2C_mC20_15_ad_f_e

http://aflow.org/prototype-encyclopedia/A2B2C_mC20_15_ad_f_e

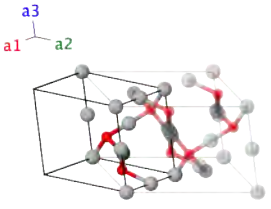


Prototype	:	Ag_2O_2Pb
AFLOW prototype label	:	A2B2C_mC20_15_ad_f_e
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC20
Space group number	:	15
Space group symbol	:	$C2/c$
AFLOW prototype command	:	`aflow --proto=A2B2C_mC20_15_ad_f_e --params=`$a, b/a, c/a, \beta, y_3, x_4, y_4, z_4$

- (Byström, 1950) gives the structure in the $I2/c$ setting of space group #15. We have used FINDSYM to change this to the standard $C2/c$ setting. This conversion involved a rotation of the axis, and placed the origin on what had been an Ag ($4d$) site, tranforming it to Ag ($4a$).

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($4a$)	Ag I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	($4a$)	Ag I
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{4}a + \frac{1}{2}c\cos\beta\right)\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	($4d$)	Ag II

$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} - \frac{1}{4}b\,\hat{\mathbf{y}}$	$(4d)$	Ag II
$\mathbf{B}_5$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Pb
$\mathbf{B}_6$	=	$y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Pb
$\mathbf{B}_7$	=	$(x_4 - y_4)\,\mathbf{a}_1 + (x_4 + y_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O
$\mathbf{B}_8$	=	$(-x_4 - y_4)\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_4 a - z_4 c\cos\beta\right)\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O
$\mathbf{B}_9$	=	$(-x_4 + y_4)\,\mathbf{a}_1 + (-x_4 - y_4)\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O
$\mathbf{B}_{10}$	=	$(x_4 + y_4)\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O

References:

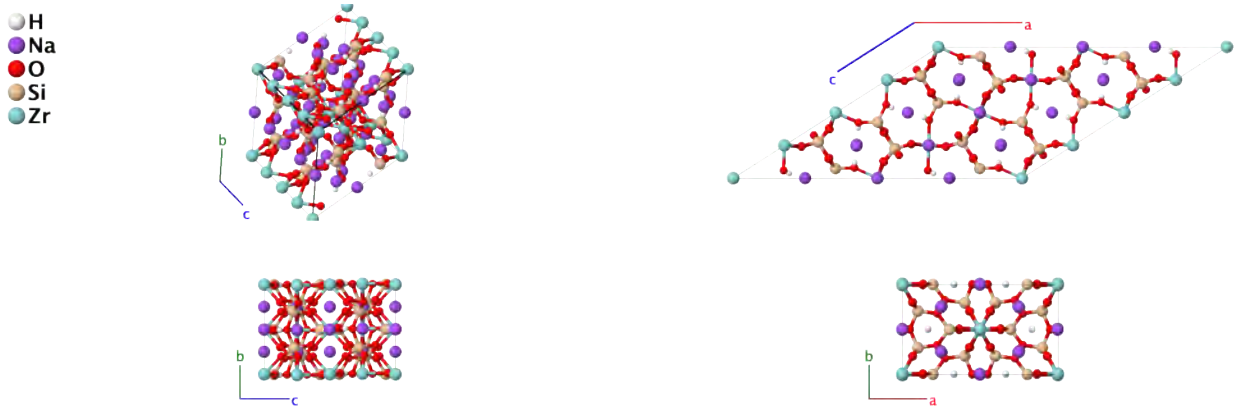
- A. Byström and L. Evers, *The Crystal Structures of Ag_2PbO_2 and $Ag5Pb_2O_6$*, Acta Chem. Scand. **4**, 613–627 (1950), doi:10.3891/acta.chem.scand.04-0613.

Geometry files:

- CIF: pp. 1559
- POSCAR: pp. 1559

Catapleiite ($Na_2ZrSi_3O_9{\cdot}2H_2O$) Structure: A2B3C9D3E_mC144_15_2f_bcdef_9f_3f_ae

http://aflow.org/prototype-encyclopedia/A2B3C9D3E_mC144_15_2f_bcdef_9f_3f_ae

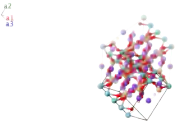


Prototype	:	$(H_2O)_2Na_2O_9Si_3Zr$
AFLOW prototype label	:	A2B3C9D3E_mC144_15_2f_bcdef_9f_3f_ae
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC144
Space group number	:	15
Space group symbol	:	$C2/c$
AFLOW prototype command	:	`aflow --proto=A2B3C9D3E_mC144_15_2f_bcdef_9f_3f_ae --params=`$a, b/a, c/a, \beta, y_5, y_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}, x_{17}, y_{17}, z_{17}, x_{18}, y_{18}, z_{18}, x_{19}, y_{19}, z_{19}, x_{20}, y_{20}, z_{20}, x_{21}, y_{21}, z_{21}$

- This is a refinement of the crystal structure of catapleiite. The original hexagonal structure was given the *Strukturbericht* designation $S\,3_4$ by (Gottfried, 1937).
- The sodium atom sites are only 66.7% occupied in this structure. However, the stoichiometry in the AFLOW label treats the Na sites as fully occupied (*i.e.*, a stoichiometric ratio of three as opposed to two).
- (Ilyushin, 1981) gave the lattice parameters and Wyckoff positions in terms of the $B2/b$ setting of space group #15. We used FINDSYM to transform this to the standard $C2/c$ setting. This required the $y-$ and $z-$axes to be swapped.

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(4a)$	Zr I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	$(4a)$	Zr I

$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	(4*b*)	Na I
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(a + c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	(4*b*)	Na I
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}}$	(4*c*)	Na II
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{4}a + \frac{1}{2}c\cos\beta\right)\,\hat{\mathbf{x}} - \frac{1}{4}b\,\hat{\mathbf{y}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	(4*c*)	Na II
$\mathbf{B}_7$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{4}a + \frac{1}{2}c\cos\beta\right)\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	(4*d*)	Na III
$\mathbf{B}_8$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} - \frac{1}{4}b\,\hat{\mathbf{y}}$	(4*d*)	Na III
$\mathbf{B}_9$	=	$-y_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Na IV
$\mathbf{B}_{10}$	=	$y_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Na IV
$\mathbf{B}_{11}$	=	$-y_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Zr II
$\mathbf{B}_{12}$	=	$y_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Zr II
$\mathbf{B}_{13}$	=	$(x_7 - y_7)\,\mathbf{a}_1 + (x_7 + y_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	H_2O I
$\mathbf{B}_{14}$	=	$(-x_7 - y_7)\,\mathbf{a}_1 + (-x_7 + y_7)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_7 a - z_7 c\cos\beta\right)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right)c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	H_2O I
$\mathbf{B}_{15}$	=	$(-x_7 + y_7)\,\mathbf{a}_1 + (-x_7 - y_7)\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	H_2O I
$\mathbf{B}_{16}$	=	$(x_7 + y_7)\,\mathbf{a}_1 + (x_7 - y_7)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_7 a + z_7 c\cos\beta\right)\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right)c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	H_2O I
$\mathbf{B}_{17}$	=	$(x_8 - y_8)\,\mathbf{a}_1 + (x_8 + y_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	H_2O II
$\mathbf{B}_{18}$	=	$(-x_8 - y_8)\,\mathbf{a}_1 + (-x_8 + y_8)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_8 a - z_8 c\cos\beta\right)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right)c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	H_2O II
$\mathbf{B}_{19}$	=	$(-x_8 + y_8)\,\mathbf{a}_1 + (-x_8 - y_8)\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	H_2O II
$\mathbf{B}_{20}$	=	$(x_8 + y_8)\,\mathbf{a}_1 + (x_8 - y_8)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_8 a + z_8 c\cos\beta\right)\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right)c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	H_2O II
$\mathbf{B}_{21}$	=	$(x_9 - y_9)\,\mathbf{a}_1 + (x_9 + y_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$(x_9 a + z_9 c\cos\beta)\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	Na V
$\mathbf{B}_{22}$	=	$(-x_9 - y_9)\,\mathbf{a}_1 + (-x_9 + y_9)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_9\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_9 a - z_9 c\cos\beta\right)\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_9\right)c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	Na V
$\mathbf{B}_{23}$	=	$(-x_9 + y_9)\,\mathbf{a}_1 + (-x_9 - y_9)\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$(-x_9 a - z_9 c\cos\beta)\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} - z_9 c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	Na V
$\mathbf{B}_{24}$	=	$(x_9 + y_9)\,\mathbf{a}_1 + (x_9 - y_9)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_9 a + z_9 c\cos\beta\right)\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right)c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	Na V
$\mathbf{B}_{25}$	=	$(x_{10} - y_{10})\,\mathbf{a}_1 + (x_{10} + y_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$(x_{10} a + z_{10} c\cos\beta)\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	O I
$\mathbf{B}_{26}$	=	$(-x_{10} - y_{10})\,\mathbf{a}_1 + (-x_{10} + y_{10})\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{10}\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_{10} a - z_{10} c\cos\beta\right)\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{10}\right)c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	O I
$\mathbf{B}_{27}$	=	$(-x_{10} + y_{10})\,\mathbf{a}_1 + (-x_{10} - y_{10})\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3$	=	$(-x_{10} a - z_{10} c\cos\beta)\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} - z_{10} c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	O I
$\mathbf{B}_{28}$	=	$(x_{10} + y_{10})\,\mathbf{a}_1 + (x_{10} - y_{10})\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{10}\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{10} a + z_{10} c\cos\beta\right)\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{10}\right)c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	O I

$$\mathbf{B}_{29} = (x_{11}-y_{11})\,\mathbf{a}_1+(x_{11}+y_{11})\,\mathbf{a}_2+z_{11}\,\mathbf{a}_3 = (x_{11}a+z_{11}c\cos\beta)\,\hat{\mathbf{x}}+y_{11}b\,\hat{\mathbf{y}}+z_{11}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O II}$$

$$\mathbf{B}_{30} = (-x_{11}-y_{11})\,\mathbf{a}_1+(-x_{11}+y_{11})\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta-x_{11}a-z_{11}c\cos\beta\right)\hat{\mathbf{x}}+y_{11}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{11}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O II}$$

$$\mathbf{B}_{31} = (-x_{11}+y_{11})\,\mathbf{a}_1+(-x_{11}-y_{11})\,\mathbf{a}_2-z_{11}\,\mathbf{a}_3 = (-x_{11}a-z_{11}c\cos\beta)\,\hat{\mathbf{x}}-y_{11}b\,\hat{\mathbf{y}}-z_{11}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O II}$$

$$\mathbf{B}_{32} = (x_{11}+y_{11})\,\mathbf{a}_1+(x_{11}-y_{11})\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta+x_{11}a+z_{11}c\cos\beta\right)\hat{\mathbf{x}}-y_{11}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{11}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O II}$$

$$\mathbf{B}_{33} = (x_{12}-y_{12})\,\mathbf{a}_1+(x_{12}+y_{12})\,\mathbf{a}_2+z_{12}\,\mathbf{a}_3 = (x_{12}a+z_{12}c\cos\beta)\,\hat{\mathbf{x}}+y_{12}b\,\hat{\mathbf{y}}+z_{12}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O III}$$

$$\mathbf{B}_{34} = (-x_{12}-y_{12})\,\mathbf{a}_1+(-x_{12}+y_{12})\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta-x_{12}a-z_{12}c\cos\beta\right)\hat{\mathbf{x}}+y_{12}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{12}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O III}$$

$$\mathbf{B}_{35} = (-x_{12}+y_{12})\,\mathbf{a}_1+(-x_{12}-y_{12})\,\mathbf{a}_2-z_{12}\,\mathbf{a}_3 = (-x_{12}a-z_{12}c\cos\beta)\,\hat{\mathbf{x}}-y_{12}b\,\hat{\mathbf{y}}-z_{12}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O III}$$

$$\mathbf{B}_{36} = (x_{12}+y_{12})\,\mathbf{a}_1+(x_{12}-y_{12})\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta+x_{12}a+z_{12}c\cos\beta\right)\hat{\mathbf{x}}-y_{12}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{12}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O III}$$

$$\mathbf{B}_{37} = (x_{13}-y_{13})\,\mathbf{a}_1+(x_{13}+y_{13})\,\mathbf{a}_2+z_{13}\,\mathbf{a}_3 = (x_{13}a+z_{13}c\cos\beta)\,\hat{\mathbf{x}}+y_{13}b\,\hat{\mathbf{y}}+z_{13}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O IV}$$

$$\mathbf{B}_{38} = (-x_{13}-y_{13})\,\mathbf{a}_1+(-x_{13}+y_{13})\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta-x_{13}a-z_{13}c\cos\beta\right)\hat{\mathbf{x}}+y_{13}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{13}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O IV}$$

$$\mathbf{B}_{39} = (-x_{13}+y_{13})\,\mathbf{a}_1+(-x_{13}-y_{13})\,\mathbf{a}_2-z_{13}\,\mathbf{a}_3 = (-x_{13}a-z_{13}c\cos\beta)\,\hat{\mathbf{x}}-y_{13}b\,\hat{\mathbf{y}}-z_{13}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O IV}$$

$$\mathbf{B}_{40} = (x_{13}+y_{13})\,\mathbf{a}_1+(x_{13}-y_{13})\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta+x_{13}a+z_{13}c\cos\beta\right)\hat{\mathbf{x}}-y_{13}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{13}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O IV}$$

$$\mathbf{B}_{41} = (x_{14}-y_{14})\,\mathbf{a}_1+(x_{14}+y_{14})\,\mathbf{a}_2+z_{14}\,\mathbf{a}_3 = (x_{14}a+z_{14}c\cos\beta)\,\hat{\mathbf{x}}+y_{14}b\,\hat{\mathbf{y}}+z_{14}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O V}$$

$$\mathbf{B}_{42} = (-x_{14}-y_{14})\,\mathbf{a}_1+(-x_{14}+y_{14})\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_{14}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta-x_{14}a-z_{14}c\cos\beta\right)\hat{\mathbf{x}}+y_{14}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{14}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O V}$$

$$\mathbf{B}_{43} = (-x_{14}+y_{14})\,\mathbf{a}_1+(-x_{14}-y_{14})\,\mathbf{a}_2-z_{14}\,\mathbf{a}_3 = (-x_{14}a-z_{14}c\cos\beta)\,\hat{\mathbf{x}}-y_{14}b\,\hat{\mathbf{y}}-z_{14}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O V}$$

$$\mathbf{B}_{44} = (x_{14}+y_{14})\,\mathbf{a}_1+(x_{14}-y_{14})\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_{14}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta+x_{14}a+z_{14}c\cos\beta\right)\hat{\mathbf{x}}-y_{14}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{14}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O V}$$

$$\mathbf{B}_{45} = (x_{15}-y_{15})\,\mathbf{a}_1+(x_{15}+y_{15})\,\mathbf{a}_2+z_{15}\,\mathbf{a}_3 = (x_{15}a+z_{15}c\cos\beta)\,\hat{\mathbf{x}}+y_{15}b\,\hat{\mathbf{y}}+z_{15}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VI}$$

$$\mathbf{B}_{46} = (-x_{15}-y_{15})\,\mathbf{a}_1+(-x_{15}+y_{15})\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_{15}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta-x_{15}a-z_{15}c\cos\beta\right)\hat{\mathbf{x}}+y_{15}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{15}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VI}$$

$$\mathbf{B}_{47} = (-x_{15}+y_{15})\,\mathbf{a}_1+(-x_{15}-y_{15})\,\mathbf{a}_2-z_{15}\,\mathbf{a}_3 = (-x_{15}a-z_{15}c\cos\beta)\,\hat{\mathbf{x}}-y_{15}b\,\hat{\mathbf{y}}-z_{15}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VI}$$

$$\mathbf{B}_{48} = (x_{15}+y_{15})\,\mathbf{a}_1+(x_{15}-y_{15})\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_{15}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta+x_{15}a+z_{15}c\cos\beta\right)\hat{\mathbf{x}}-y_{15}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{15}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VI}$$

$$\mathbf{B}_{49} = (x_{16}-y_{16})\,\mathbf{a}_1+(x_{16}+y_{16})\,\mathbf{a}_2+z_{16}\,\mathbf{a}_3 = (x_{16}a+z_{16}c\cos\beta)\,\hat{\mathbf{x}}+y_{16}b\,\hat{\mathbf{y}}+z_{16}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VII}$$

$$\mathbf{B}_{50} = (-x_{16}-y_{16})\,\mathbf{a}_1+(-x_{16}+y_{16})\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_{16}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta-x_{16}a-z_{16}c\cos\beta\right)\hat{\mathbf{x}}+y_{16}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{16}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VII}$$

$$\mathbf{B}_{51} = (-x_{16}+y_{16})\,\mathbf{a}_1 + (-x_{16}-y_{16})\,\mathbf{a}_2 - z_{16}\,\mathbf{a}_3 = (-x_{16}a - z_{16}c\cos\beta)\,\hat{\mathbf{x}} - y_{16}b\,\hat{\mathbf{y}} - z_{16}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VII}$$

$$\mathbf{B}_{52} = (x_{16}+y_{16})\,\mathbf{a}_1 + (x_{16}-y_{16})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{16}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{16}a + z_{16}c\cos\beta\right)\hat{\mathbf{x}} - y_{16}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{16}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VII}$$

$$\mathbf{B}_{53} = (x_{17}-y_{17})\,\mathbf{a}_1 + (x_{17}+y_{17})\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 = (x_{17}a + z_{17}c\cos\beta)\,\hat{\mathbf{x}} + y_{17}b\,\hat{\mathbf{y}} + z_{17}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VIII}$$

$$\mathbf{B}_{54} = (-x_{17}-y_{17})\,\mathbf{a}_1 + (-x_{17}+y_{17})\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_{17}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{17}a - z_{17}c\cos\beta\right)\hat{\mathbf{x}} + y_{17}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_{17}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VIII}$$

$$\mathbf{B}_{55} = (-x_{17}+y_{17})\,\mathbf{a}_1 + (-x_{17}-y_{17})\,\mathbf{a}_2 - z_{17}\,\mathbf{a}_3 = (-x_{17}a - z_{17}c\cos\beta)\,\hat{\mathbf{x}} - y_{17}b\,\hat{\mathbf{y}} - z_{17}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VIII}$$

$$\mathbf{B}_{56} = (x_{17}+y_{17})\,\mathbf{a}_1 + (x_{17}-y_{17})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{17}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{17}a + z_{17}c\cos\beta\right)\hat{\mathbf{x}} - y_{17}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{17}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VIII}$$

$$\mathbf{B}_{57} = (x_{18}-y_{18})\,\mathbf{a}_1 + (x_{18}+y_{18})\,\mathbf{a}_2 + z_{18}\,\mathbf{a}_3 = (x_{18}a + z_{18}c\cos\beta)\,\hat{\mathbf{x}} + y_{18}b\,\hat{\mathbf{y}} + z_{18}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O IX}$$

$$\mathbf{B}_{58} = (-x_{18}-y_{18})\,\mathbf{a}_1 + (-x_{18}+y_{18})\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_{18}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{18}a - z_{18}c\cos\beta\right)\hat{\mathbf{x}} + y_{18}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_{18}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O IX}$$

$$\mathbf{B}_{59} = (-x_{18}+y_{18})\,\mathbf{a}_1 + (-x_{18}-y_{18})\,\mathbf{a}_2 - z_{18}\,\mathbf{a}_3 = (-x_{18}a - z_{18}c\cos\beta)\,\hat{\mathbf{x}} - y_{18}b\,\hat{\mathbf{y}} - z_{18}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O IX}$$

$$\mathbf{B}_{60} = (x_{18}+y_{18})\,\mathbf{a}_1 + (x_{18}-y_{18})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{18}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{18}a + z_{18}c\cos\beta\right)\hat{\mathbf{x}} - y_{18}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{18}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O IX}$$

$$\mathbf{B}_{61} = (x_{19}-y_{19})\,\mathbf{a}_1 + (x_{19}+y_{19})\,\mathbf{a}_2 + z_{19}\,\mathbf{a}_3 = (x_{19}a + z_{19}c\cos\beta)\,\hat{\mathbf{x}} + y_{19}b\,\hat{\mathbf{y}} + z_{19}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{Si I}$$

$$\mathbf{B}_{62} = (-x_{19}-y_{19})\,\mathbf{a}_1 + (-x_{19}+y_{19})\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_{19}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{19}a - z_{19}c\cos\beta\right)\hat{\mathbf{x}} + y_{19}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_{19}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{Si I}$$

$$\mathbf{B}_{63} = (-x_{19}+y_{19})\,\mathbf{a}_1 + (-x_{19}-y_{19})\,\mathbf{a}_2 - z_{19}\,\mathbf{a}_3 = (-x_{19}a - z_{19}c\cos\beta)\,\hat{\mathbf{x}} - y_{19}b\,\hat{\mathbf{y}} - z_{19}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{Si I}$$

$$\mathbf{B}_{64} = (x_{19}+y_{19})\,\mathbf{a}_1 + (x_{19}-y_{19})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{19}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{19}a + z_{19}c\cos\beta\right)\hat{\mathbf{x}} - y_{19}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{19}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{Si I}$$

$$\mathbf{B}_{65} = (x_{20}-y_{20})\,\mathbf{a}_1 + (x_{20}+y_{20})\,\mathbf{a}_2 + z_{20}\,\mathbf{a}_3 = (x_{20}a + z_{20}c\cos\beta)\,\hat{\mathbf{x}} + y_{20}b\,\hat{\mathbf{y}} + z_{20}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{Si II}$$

$$\mathbf{B}_{66} = (-x_{20}-y_{20})\,\mathbf{a}_1 + (-x_{20}+y_{20})\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_{20}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{20}a - z_{20}c\cos\beta\right)\hat{\mathbf{x}} + y_{20}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_{20}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{Si II}$$

$$\mathbf{B}_{67} = (-x_{20}+y_{20})\,\mathbf{a}_1 + (-x_{20}-y_{20})\,\mathbf{a}_2 - z_{20}\,\mathbf{a}_3 = (-x_{20}a - z_{20}c\cos\beta)\,\hat{\mathbf{x}} - y_{20}b\,\hat{\mathbf{y}} - z_{20}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{Si II}$$

$$\mathbf{B}_{68} = (x_{20}+y_{20})\,\mathbf{a}_1 + (x_{20}-y_{20})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{20}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{20}a + z_{20}c\cos\beta\right)\hat{\mathbf{x}} - y_{20}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{20}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{Si II}$$

$$\mathbf{B}_{69} = (x_{21}-y_{21})\,\mathbf{a}_1 + (x_{21}+y_{21})\,\mathbf{a}_2 + z_{21}\,\mathbf{a}_3 = (x_{21}a + z_{21}c\cos\beta)\,\hat{\mathbf{x}} + y_{21}b\,\hat{\mathbf{y}} + z_{21}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{Si III}$$

$$\mathbf{B}_{70} = (-x_{21}-y_{21})\,\mathbf{a}_1 + (-x_{21}+y_{21})\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_{21}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{21}a - z_{21}c\cos\beta\right)\hat{\mathbf{x}} + y_{21}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_{21}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{Si III}$$

$$\mathbf{B}_{71} = (-x_{21}+y_{21})\,\mathbf{a}_1 + (-x_{21}-y_{21})\,\mathbf{a}_2 - z_{21}\,\mathbf{a}_3 = (-x_{21}a - z_{21}c\cos\beta)\,\hat{\mathbf{x}} - y_{21}b\,\hat{\mathbf{y}} - z_{21}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{Si III}$$

$$\mathbf{B}_{72} = (x_{21}+y_{21})\,\mathbf{a}_1 + (x_{21}-y_{21})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{21}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{21}a + z_{21}c\cos\beta\right)\hat{\mathbf{x}} - y_{21}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{21}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{Si III}$$

References:

- G. D. Ilyushin, A. A. Voronkov, V. V. Ilyukhin, N. N. Nevskii, and N. V. Belov, *Crystal structure of natural monoclinic catapleiite, $Na_2ZrSi_3O_9 \cdot 2H_2O$*, Doklady Akademii Nauk SSSR **260**, 623–627 (1981).
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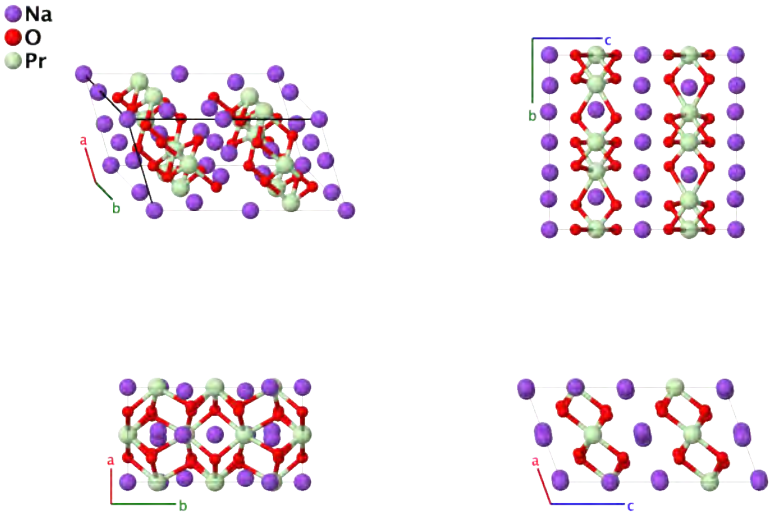
Geometry files:

- CIF: pp. 1559
- POSCAR: pp. 1560

Na_2PrO_3 Structure: A2B3C_mC48_15_aef_3f_2e

http://aflow.org/prototype-encyclopedia/A2B3C_mC48_15_aef_3f_2e



Prototype	:	Na_2O_3Pr
AFLOW prototype label	:	A2B3C_mC48_15_aef_3f_2e
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC48
Space group number	:	15
Space group symbol	:	$C2/c$
AFLOW prototype command	:	`aflow --proto=A2B3C_mC48_15_aef_3f_2e --params=`$a, b/a, c/a, \beta, y_2, y_3, y_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8$

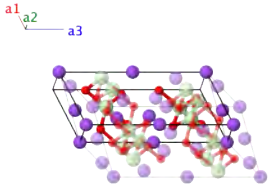
Other compounds with this structure

- Na_2CrO_3 and Na_2TbO_3

- (Hinatsu, 2006) found that the site we have labeled Na-II is actually 2/3 sodium and 1/3 praseodymium, statistically distributed, while the site we label Pr-I is 1/3 sodium and 2/3 praseodymium.

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(4*a*)	Na I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	(4*a*)	Na I
$\mathbf{B}_3$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Na II
$\mathbf{B}_4$	=	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Na II
$\mathbf{B}_5$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Pr I
$\mathbf{B}_6$	=	$y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Pr I
$\mathbf{B}_7$	=	$-y_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Pr II
$\mathbf{B}_8$	=	$y_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Pr II
$\mathbf{B}_9$	=	$(x_5 - y_5)\,\mathbf{a}_1 + (x_5 + y_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	Na III
$\mathbf{B}_{10}$	=	$(-x_5 - y_5)\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_5 a - z_5 c\cos\beta\right)\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	Na III
$\mathbf{B}_{11}$	=	$(-x_5 + y_5)\,\mathbf{a}_1 + (-x_5 - y_5)\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	Na III
$\mathbf{B}_{12}$	=	$(x_5 + y_5)\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	Na III
$\mathbf{B}_{13}$	=	$(x_6 - y_6)\,\mathbf{a}_1 + (x_6 + y_6)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	O I
$\mathbf{B}_{14}$	=	$(-x_6 - y_6)\,\mathbf{a}_1 + (-x_6 + y_6)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_6 a - z_6 c\cos\beta\right)\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	O I
$\mathbf{B}_{15}$	=	$(-x_6 + y_6)\,\mathbf{a}_1 + (-x_6 - y_6)\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	O I
$\mathbf{B}_{16}$	=	$(x_6 + y_6)\,\mathbf{a}_1 + (x_6 - y_6)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	O I
$\mathbf{B}_{17}$	=	$(x_7 - y_7)\,\mathbf{a}_1 + (x_7 + y_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	O II
$\mathbf{B}_{18}$	=	$(-x_7 - y_7)\,\mathbf{a}_1 + (-x_7 + y_7)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_7 a - z_7 c\cos\beta\right)\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right)c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	O II
$\mathbf{B}_{19}$	=	$(-x_7 + y_7)\,\mathbf{a}_1 + (-x_7 - y_7)\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	O II
$\mathbf{B}_{20}$	=	$(x_7 + y_7)\,\mathbf{a}_1 + (x_7 - y_7)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_7 a + z_7 c\cos\beta\right)\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right)c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	O II
$\mathbf{B}_{21}$	=	$(x_8 - y_8)\,\mathbf{a}_1 + (x_8 + y_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	O III
$\mathbf{B}_{22}$	=	$(-x_8 - y_8)\,\mathbf{a}_1 + (-x_8 + y_8)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_8 a - z_8 c\cos\beta\right)\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right)c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	O III
$\mathbf{B}_{23}$	=	$(-x_8 + y_8)\,\mathbf{a}_1 + (-x_8 - y_8)\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	O III
$\mathbf{B}_{24}$	=	$(x_8 + y_8)\,\mathbf{a}_1 + (x_8 - y_8)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_8 a + z_8 c\cos\beta\right)\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right)c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	O III

References:

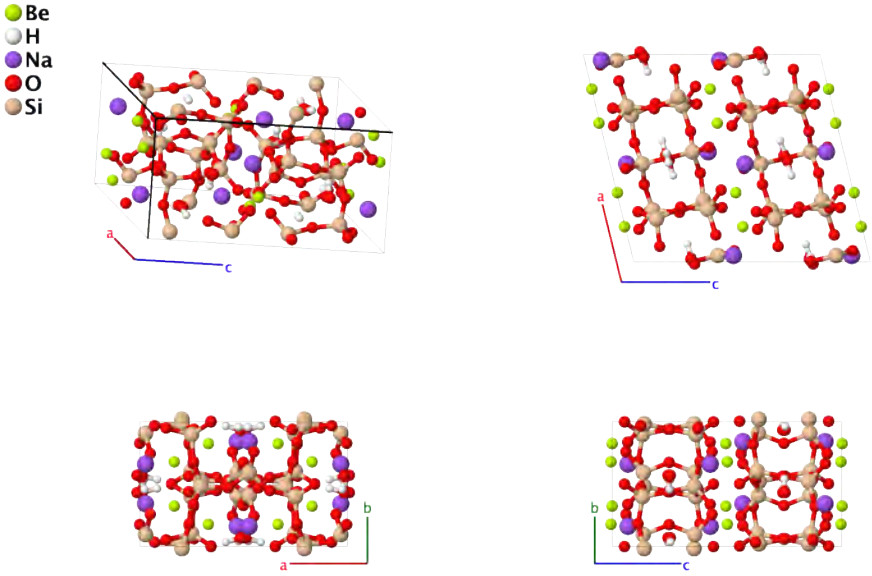
- Y. Hinatsu and Y. Doi, *Crystal structures and magnetic properties of alkali-metal lanthanide oxides A_2LnO_3 (A = Li, Na;*

Ln = Ce, Pr, Tb), J. Alloys Compd. **418**, 155–160 (2006), doi:10.1016/j.jallcom.2005.08.100.

Geometry files:
- CIF: pp. 1560
- POSCAR: pp. 1560

Eudidymite ($BeHNaO_8Si_3$) Structure: A2B4C2D17E6_mC124_15_f_2f_f_e8f_3f

http://aflow.org/prototype-encyclopedia/A2B4C2D17E6_mC124_15_f_2f_f_e8f_3f

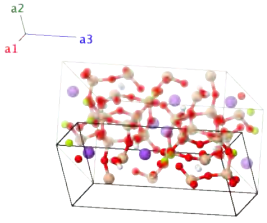


Prototype	:	$BeHNaO_8Si_3$
AFLOW prototype label	:	A2B4C2D17E6_mC124_15_f_2f_f_e8f_3f
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC124
Space group number	:	15
Space group symbol	:	$C2/c$
AFLOW prototype command	:	`aflow --proto=A2B4C2D17E6_mC124_15_f_2f_f_e8f_3f --params=`$a, b/a, c/a, \beta, y_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}$

- Eudidymite and its dimorph, epididymite ($S4_7$) are two forms of hydrated sodium beryllium silicate which are stable under ambient conditions. (Diego Gatta, 2008)
- The occupancy of the sites making up the water molecule (H-I, H-II and O-IX) is 50%. In the picture on this page the atoms forming the water molecules appear doubled because of this.

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_2$	=	$y_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_3$	=	$(x_2 - y_2)\,\mathbf{a}_1 + (x_2 + y_2)\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Be
$\mathbf{B}_4$	=	$(-x_2 - y_2)\,\mathbf{a}_1 + (-x_2 + y_2)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_2 a - z_2 c\cos\beta\right)\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Be
$\mathbf{B}_5$	=	$(-x_2 + y_2)\,\mathbf{a}_1 + (-x_2 - y_2)\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} - z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Be
$\mathbf{B}_6$	=	$(x_2 + y_2)\,\mathbf{a}_1 + (x_2 - y_2)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_2 a + z_2 c\cos\beta\right)\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Be
$\mathbf{B}_7$	=	$(x_3 - y_3)\,\mathbf{a}_1 + (x_3 + y_3)\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H I
$\mathbf{B}_8$	=	$(-x_3 - y_3)\,\mathbf{a}_1 + (-x_3 + y_3)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_3 a - z_3 c\cos\beta\right)\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H I
$\mathbf{B}_9$	=	$(-x_3 + y_3)\,\mathbf{a}_1 + (-x_3 - y_3)\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H I
$\mathbf{B}_{10}$	=	$(x_3 + y_3)\,\mathbf{a}_1 + (x_3 - y_3)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H I
$\mathbf{B}_{11}$	=	$(x_4 - y_4)\,\mathbf{a}_1 + (x_4 + y_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H II
$\mathbf{B}_{12}$	=	$(-x_4 - y_4)\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_4 a - z_4 c\cos\beta\right)\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H II
$\mathbf{B}_{13}$	=	$(-x_4 + y_4)\,\mathbf{a}_1 + (-x_4 - y_4)\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H II
$\mathbf{B}_{14}$	=	$(x_4 + y_4)\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H II
$\mathbf{B}_{15}$	=	$(x_5 - y_5)\,\mathbf{a}_1 + (x_5 + y_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Na
$\mathbf{B}_{16}$	=	$(-x_5 - y_5)\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_5 a - z_5 c\cos\beta\right)\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Na
$\mathbf{B}_{17}$	=	$(-x_5 + y_5)\,\mathbf{a}_1 + (-x_5 - y_5)\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Na
$\mathbf{B}_{18}$	=	$(x_5 + y_5)\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Na
$\mathbf{B}_{19}$	=	$(x_6 - y_6)\,\mathbf{a}_1 + (x_6 + y_6)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{20}$	=	$(-x_6 - y_6)\,\mathbf{a}_1 + (-x_6 + y_6)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_6 a - z_6 c\cos\beta\right)\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{21}$	=	$(-x_6 + y_6)\,\mathbf{a}_1 + (-x_6 - y_6)\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II

$$\mathbf{B}_{22} = (x_6 + y_6)\,\mathbf{a}_1 + (x_6 - y_6)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O II}$$

$$\mathbf{B}_{23} = (x_7 - y_7)\,\mathbf{a}_1 + (x_7 + y_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 = (x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O III}$$

$$\mathbf{B}_{24} = (-x_7 - y_7)\,\mathbf{a}_1 + (-x_7 + y_7)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_7 a - z_7 c\cos\beta\right)\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_7\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O III}$$

$$\mathbf{B}_{25} = (-x_7 + y_7)\,\mathbf{a}_1 + (-x_7 - y_7)\,\mathbf{a}_2 - z_7\,\mathbf{a}_3 = (-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O III}$$

$$\mathbf{B}_{26} = (x_7 + y_7)\,\mathbf{a}_1 + (x_7 - y_7)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_7 a + z_7 c\cos\beta\right)\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_7\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O III}$$

$$\mathbf{B}_{27} = (x_8 - y_8)\,\mathbf{a}_1 + (x_8 + y_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 = (x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O IV}$$

$$\mathbf{B}_{28} = (-x_8 - y_8)\,\mathbf{a}_1 + (-x_8 + y_8)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_8 a - z_8 c\cos\beta\right)\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_8\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O IV}$$

$$\mathbf{B}_{29} = (-x_8 + y_8)\,\mathbf{a}_1 + (-x_8 - y_8)\,\mathbf{a}_2 - z_8\,\mathbf{a}_3 = (-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O IV}$$

$$\mathbf{B}_{30} = (x_8 + y_8)\,\mathbf{a}_1 + (x_8 - y_8)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_8 a + z_8 c\cos\beta\right)\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_8\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O IV}$$

$$\mathbf{B}_{31} = (x_9 - y_9)\,\mathbf{a}_1 + (x_9 + y_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 = (x_9 a + z_9 c\cos\beta)\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O V}$$

$$\mathbf{B}_{32} = (-x_9 - y_9)\,\mathbf{a}_1 + (-x_9 + y_9)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_9 a - z_9 c\cos\beta\right)\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_9\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O V}$$

$$\mathbf{B}_{33} = (-x_9 + y_9)\,\mathbf{a}_1 + (-x_9 - y_9)\,\mathbf{a}_2 - z_9\,\mathbf{a}_3 = (-x_9 a - z_9 c\cos\beta)\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} - z_9 c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O V}$$

$$\mathbf{B}_{34} = (x_9 + y_9)\,\mathbf{a}_1 + (x_9 - y_9)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_9 a + z_9 c\cos\beta\right)\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_9\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O V}$$

$$\mathbf{B}_{35} = (x_{10} - y_{10})\,\mathbf{a}_1 + (x_{10} + y_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 = (x_{10} a + z_{10} c\cos\beta)\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VI}$$

$$\mathbf{B}_{36} = (-x_{10} - y_{10})\,\mathbf{a}_1 + (-x_{10} + y_{10})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{10} a - z_{10} c\cos\beta\right)\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{10}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VI}$$

$$\mathbf{B}_{37} = (-x_{10} + y_{10})\,\mathbf{a}_1 + (-x_{10} - y_{10})\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 = (-x_{10} a - z_{10} c\cos\beta)\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} - z_{10} c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VI}$$

$$\mathbf{B}_{38} = (x_{10} + y_{10})\,\mathbf{a}_1 + (x_{10} - y_{10})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{10} a + z_{10} c\cos\beta\right)\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{10}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VI}$$

$$\mathbf{B}_{39} = (x_{11} - y_{11})\,\mathbf{a}_1 + (x_{11} + y_{11})\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 = (x_{11} a + z_{11} c\cos\beta)\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + z_{11} c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VII}$$

$$\mathbf{B}_{40} = (-x_{11} - y_{11})\,\mathbf{a}_1 + (-x_{11} + y_{11})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{11} a - z_{11} c\cos\beta\right)\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{11}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VII}$$

$$\mathbf{B}_{41} = (-x_{11} + y_{11})\,\mathbf{a}_1 + (-x_{11} - y_{11})\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3 = (-x_{11} a - z_{11} c\cos\beta)\,\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} - z_{11} c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VII}$$

$$\mathbf{B}_{42} = (x_{11} + y_{11})\,\mathbf{a}_1 + (x_{11} - y_{11})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{11} a + z_{11} c\cos\beta\right)\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{11}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VII}$$

$$\mathbf{B}_{43} = (x_{12} - y_{12})\,\mathbf{a}_1 + (x_{12} + y_{12})\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 = (x_{12} a + z_{12} c\cos\beta)\,\hat{\mathbf{x}} + y_{12} b\,\hat{\mathbf{y}} + z_{12} c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VIII}$$

$$\mathbf{B}_{44} = (-x_{12} - y_{12})\,\mathbf{a}_1 + (-x_{12} + y_{12})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{12}a - z_{12}c\cos\beta\right)\hat{\mathbf{x}} + y_{12}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{12}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VIII}$$

$$\mathbf{B}_{45} = (-x_{12} + y_{12})\,\mathbf{a}_1 + (-x_{12} - y_{12})\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3 = (-x_{12}a - z_{12}c\cos\beta)\,\hat{\mathbf{x}} - y_{12}b\,\hat{\mathbf{y}} - z_{12}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VIII}$$

$$\mathbf{B}_{46} = (x_{12} + y_{12})\,\mathbf{a}_1 + (x_{12} - y_{12})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{12}a + z_{12}c\cos\beta\right)\hat{\mathbf{x}} - y_{12}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{12}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VIII}$$

$$\mathbf{B}_{47} = (x_{13} - y_{13})\,\mathbf{a}_1 + (x_{13} + y_{13})\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3 = (x_{13}a + z_{13}c\cos\beta)\,\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} + z_{13}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O IX}$$

$$\mathbf{B}_{48} = (-x_{13} - y_{13})\,\mathbf{a}_1 + (-x_{13} + y_{13})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{13}a - z_{13}c\cos\beta\right)\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{13}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O IX}$$

$$\mathbf{B}_{49} = (-x_{13} + y_{13})\,\mathbf{a}_1 + (-x_{13} - y_{13})\,\mathbf{a}_2 - z_{13}\,\mathbf{a}_3 = (-x_{13}a - z_{13}c\cos\beta)\,\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} - z_{13}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O IX}$$

$$\mathbf{B}_{50} = (x_{13} + y_{13})\,\mathbf{a}_1 + (x_{13} - y_{13})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{13}a + z_{13}c\cos\beta\right)\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{13}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O IX}$$

$$\mathbf{B}_{51} = (x_{14} - y_{14})\,\mathbf{a}_1 + (x_{14} + y_{14})\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3 = (x_{14}a + z_{14}c\cos\beta)\,\hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} + z_{14}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{Si I}$$

$$\mathbf{B}_{52} = (-x_{14} - y_{14})\,\mathbf{a}_1 + (-x_{14} + y_{14})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{14}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{14}a - z_{14}c\cos\beta\right)\hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{14}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{Si I}$$

$$\mathbf{B}_{53} = (-x_{14} + y_{14})\,\mathbf{a}_1 + (-x_{14} - y_{14})\,\mathbf{a}_2 - z_{14}\,\mathbf{a}_3 = (-x_{14}a - z_{14}c\cos\beta)\,\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} - z_{14}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{Si I}$$

$$\mathbf{B}_{54} = (x_{14} + y_{14})\,\mathbf{a}_1 + (x_{14} - y_{14})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{14}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{14}a + z_{14}c\cos\beta\right)\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{14}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{Si I}$$

$$\mathbf{B}_{55} = (x_{15} - y_{15})\,\mathbf{a}_1 + (x_{15} + y_{15})\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 = (x_{15}a + z_{15}c\cos\beta)\,\hat{\mathbf{x}} + y_{15}b\,\hat{\mathbf{y}} + z_{15}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{Si II}$$

$$\mathbf{B}_{56} = (-x_{15} - y_{15})\,\mathbf{a}_1 + (-x_{15} + y_{15})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{15}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{15}a - z_{15}c\cos\beta\right)\hat{\mathbf{x}} + y_{15}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{15}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{Si II}$$

$$\mathbf{B}_{57} = (-x_{15} + y_{15})\,\mathbf{a}_1 + (-x_{15} - y_{15})\,\mathbf{a}_2 - z_{15}\,\mathbf{a}_3 = (-x_{15}a - z_{15}c\cos\beta)\,\hat{\mathbf{x}} - y_{15}b\,\hat{\mathbf{y}} - z_{15}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{Si II}$$

$$\mathbf{B}_{58} = (x_{15} + y_{15})\,\mathbf{a}_1 + (x_{15} - y_{15})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{15}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{15}a + z_{15}c\cos\beta\right)\hat{\mathbf{x}} - y_{15}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{15}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{Si II}$$

$$\mathbf{B}_{59} = (x_{16} - y_{16})\,\mathbf{a}_1 + (x_{16} + y_{16})\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 = (x_{16}a + z_{16}c\cos\beta)\,\hat{\mathbf{x}} + y_{16}b\,\hat{\mathbf{y}} + z_{16}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{Si III}$$

$$\mathbf{B}_{60} = (-x_{16} - y_{16})\,\mathbf{a}_1 + (-x_{16} + y_{16})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{16}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{16}a - z_{16}c\cos\beta\right)\hat{\mathbf{x}} + y_{16}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{16}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{Si III}$$

$$\mathbf{B}_{61} = (-x_{16} + y_{16})\,\mathbf{a}_1 + (-x_{16} - y_{16})\,\mathbf{a}_2 - z_{16}\,\mathbf{a}_3 = (-x_{16}a - z_{16}c\cos\beta)\,\hat{\mathbf{x}} - y_{16}b\,\hat{\mathbf{y}} - z_{16}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{Si III}$$

$$\mathbf{B}_{62} = (x_{16} + y_{16})\,\mathbf{a}_1 + (x_{16} - y_{16})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{16}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{16}a + z_{16}c\cos\beta\right)\hat{\mathbf{x}} - y_{16}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{16}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{Si III}$$

References:

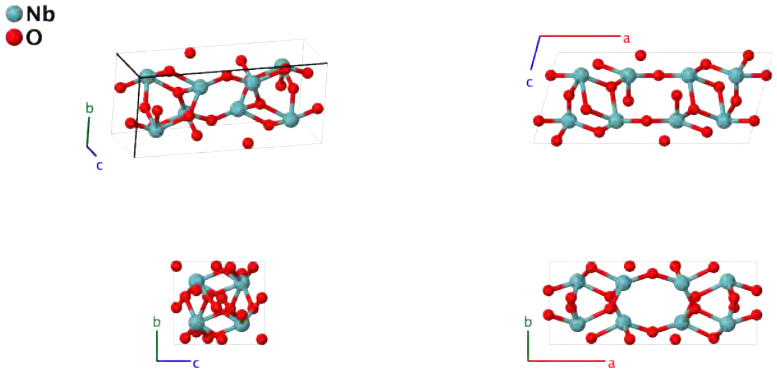
- G. Diego Gatta, N. Rotiroti, G. J. McIntyre, A. Guastoni, and F. Nestola, *New insights into the crystal chemistry of epididymite and eudidymite from Malosa, Malawi: A single-crystal neutron diffraction study*, Am. Mineral. **93**, 1158–1165 (2008), doi:10.2138/am.2008.2965.

Geometry files:

- CIF: pp. 1561
- POSCAR: pp. 1561

ζ-Nb_2O_5 (B-Nb_2O_5) Structure: A2B5_mC28_15_f_e2f

http://aflow.org/prototype-encyclopedia/A2B5_mC28_15_f_e2f.Nb2O5



Prototype	:	Nb_2O_5
AFLOW prototype label	:	A2B5_mC28_15_f_e2f
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC28
Space group number	:	15
Space group symbol	:	$C2/c$
AFLOW prototype command	:	`aflow --proto=A2B5_mC28_15_f_e2f --params=`$a, b/a, c/a, \beta, y_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4$

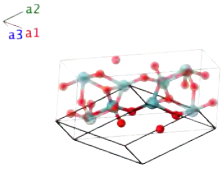
Other compounds with this structure

- Sb_2O_5

- This structure is referred to as both ζ-Nb_2O_5 and B-Nb_2O_5.
- Nb_2O_5 and B_2Pd_5 share the same AFLOW prototype label, A2B5_mC28_15_f_e2f, but have substantially different environments around each atom. The structures are generated by the same symmetry operations with different sets of parameters (`--params`) specified in their corresponding CIF files.

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_2$	=	$y_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_3$	=	$(x_2 - y_2)\,\mathbf{a}_1 + (x_2 + y_2)\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Nb

$\mathbf{B}_4$	=	$(-x_2 - y_2)\,\mathbf{a}_1 + (-x_2 + y_2)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_2 a - z_2 c\cos\beta\right)\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Nb
$\mathbf{B}_5$	=	$(-x_2 + y_2)\,\mathbf{a}_1 + (-x_2 - y_2)\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} - z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Nb
$\mathbf{B}_6$	=	$(x_2 + y_2)\,\mathbf{a}_1 + (x_2 - y_2)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_2 a + z_2 c\cos\beta\right)\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Nb
$\mathbf{B}_7$	=	$(x_3 - y_3)\,\mathbf{a}_1 + (x_3 + y_3)\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_8$	=	$(-x_3 - y_3)\,\mathbf{a}_1 + (-x_3 + y_3)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_3 a - z_3 c\cos\beta\right)\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_9$	=	$(-x_3 + y_3)\,\mathbf{a}_1 + (-x_3 - y_3)\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{10}$	=	$(x_3 + y_3)\,\mathbf{a}_1 + (x_3 - y_3)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{11}$	=	$(x_4 - y_4)\,\mathbf{a}_1 + (x_4 + y_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{12}$	=	$(-x_4 - y_4)\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_4 a - z_4 c\cos\beta\right)\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{13}$	=	$(-x_4 + y_4)\,\mathbf{a}_1 + (-x_4 - y_4)\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{14}$	=	$(x_4 + y_4)\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III

References:

- T. S. Ercit, *Refinement of the structure of ζ-Nb_2O_5 and its relationship to the rutile and thoreaulite structures*, Mineral. Petrol. **43**, 217–223 (1991), doi:10.1007/BF01166893.

Found in:

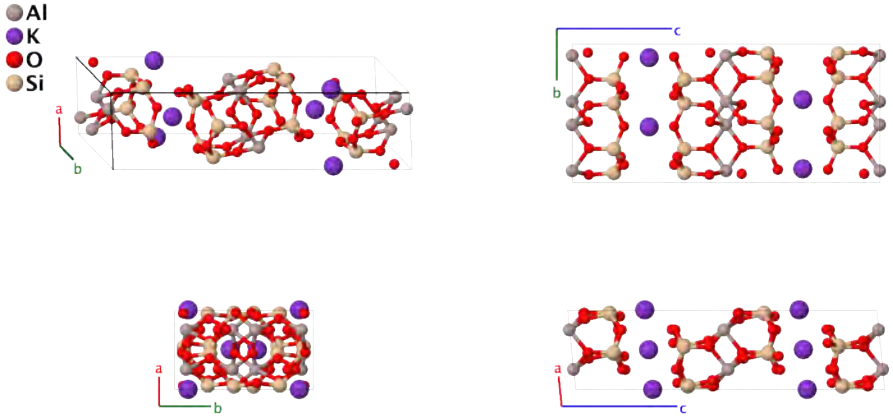
- T. S. Ercit, *ResearchGate*, http://www.researchgate.net/publication/252542874_Refinement_of_the_structure_of_-Nb2O5_and_its_relationship_to_the_rutile_and_thoreaulite_structures. .

Geometry files:

- CIF: pp. 1561
- POSCAR: pp. 1562

Muscovite ($KH_2Al_3Si_3O_{12}$, $S5_1$) Structure: A2BC10D2E4_mC76_15_f_e_5f_f_2f

http://aflow.org/prototype-encyclopedia/A2BC10D2E4_mC76_15_f_e_5f_f_2f

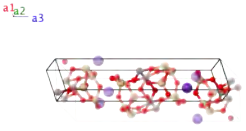


Prototype	:	$Al_3KO_{10}(OH)_2Si_3$
AFLOW prototype label	:	A2BC10D2E4_mC76_15_f_e_5f_f_2f
***Strukturbericht* designation**	:	$S5_1$
Pearson symbol	:	mC76
Space group number	:	15
Space group symbol	:	$C2/c$
AFLOW prototype command	:	`aflow --proto=A2BC10D2E4_mC76_15_f_e_5f_f_2f --params=`$a, b/a, c/a, \beta, y_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}$

- The sites we label Si-I and Si-II are actually approximately 75% silicon and 25% aluminum, but trace elements like manganese can appear on both sites. In addition, the potassium site (K) can be alloyed with small amounts of sodium, and the aluminum site (Al) with iron. These trace elements give muscovite a variety of colors. The sample studied by (Richardson, 1982) was pink.

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	K
$\mathbf{B}_2$	=	$y_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	K
$\mathbf{B}_3$	=	$(x_2 - y_2)\,\mathbf{a}_1 + (x_2 + y_2)\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Al

$$\mathbf{B}_4 = (-x_2 - y_2)\,\mathbf{a}_1 + (-x_2 + y_2)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_2\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_2 a - z_2 c\cos\beta\right)\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_2\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{Al}$$

$$\mathbf{B}_5 = (-x_2 + y_2)\,\mathbf{a}_1 + (-x_2 - y_2)\,\mathbf{a}_2 - z_2\,\mathbf{a}_3 = (-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} - z_2 c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{Al}$$

$$\mathbf{B}_6 = (x_2 + y_2)\,\mathbf{a}_1 + (x_2 - y_2)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_2\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_2 a + z_2 c\cos\beta\right)\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_2\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{Al}$$

$$\mathbf{B}_7 = (x_3 - y_3)\,\mathbf{a}_1 + (x_3 + y_3)\,\mathbf{a}_2 + z_3\,\mathbf{a}_3 = (x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O I}$$

$$\mathbf{B}_8 = (-x_3 - y_3)\,\mathbf{a}_1 + (-x_3 + y_3)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_3 a - z_3 c\cos\beta\right)\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_3\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O I}$$

$$\mathbf{B}_9 = (-x_3 + y_3)\,\mathbf{a}_1 + (-x_3 - y_3)\,\mathbf{a}_2 - z_3\,\mathbf{a}_3 = (-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O I}$$

$$\mathbf{B}_{10} = (x_3 + y_3)\,\mathbf{a}_1 + (x_3 - y_3)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_3\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O I}$$

$$\mathbf{B}_{11} = (x_4 - y_4)\,\mathbf{a}_1 + (x_4 + y_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3 = (x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O II}$$

$$\mathbf{B}_{12} = (-x_4 - y_4)\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_4 a - z_4 c\cos\beta\right)\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_4\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O II}$$

$$\mathbf{B}_{13} = (-x_4 + y_4)\,\mathbf{a}_1 + (-x_4 - y_4)\,\mathbf{a}_2 - z_4\,\mathbf{a}_3 = (-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O II}$$

$$\mathbf{B}_{14} = (x_4 + y_4)\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_4\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O II}$$

$$\mathbf{B}_{15} = (x_5 - y_5)\,\mathbf{a}_1 + (x_5 + y_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3 = (x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O III}$$

$$\mathbf{B}_{16} = (-x_5 - y_5)\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_5 a - z_5 c\cos\beta\right)\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_5\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O III}$$

$$\mathbf{B}_{17} = (-x_5 + y_5)\,\mathbf{a}_1 + (-x_5 - y_5)\,\mathbf{a}_2 - z_5\,\mathbf{a}_3 = (-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O III}$$

$$\mathbf{B}_{18} = (x_5 + y_5)\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_5\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O III}$$

$$\mathbf{B}_{19} = (x_6 - y_6)\,\mathbf{a}_1 + (x_6 + y_6)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 = (x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O IV}$$

$$\mathbf{B}_{20} = (-x_6 - y_6)\,\mathbf{a}_1 + (-x_6 + y_6)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_6 a - z_6 c\cos\beta\right)\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_6\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O IV}$$

$$\mathbf{B}_{21} = (-x_6 + y_6)\,\mathbf{a}_1 + (-x_6 - y_6)\,\mathbf{a}_2 - z_6\,\mathbf{a}_3 = (-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O IV}$$

$$\mathbf{B}_{22} = (x_6 + y_6)\,\mathbf{a}_1 + (x_6 - y_6)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O IV}$$

$$\mathbf{B}_{23} = (x_7 - y_7)\,\mathbf{a}_1 + (x_7 + y_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 = (x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O V}$$

$$\mathbf{B}_{24} = (-x_7 - y_7)\,\mathbf{a}_1 + (-x_7 + y_7)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_7 a - z_7 c\cos\beta\right)\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_7\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O V}$$

$$\mathbf{B}_{25} = (-x_7 + y_7)\,\mathbf{a}_1 + (-x_7 - y_7)\,\mathbf{a}_2 - z_7\,\mathbf{a}_3 = (-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O V}$$

$\mathbf{B}_{26}$	=	$(x_7 + y_7)\,\mathbf{a}_1 + (x_7 - y_7)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_7 a + z_7 c\cos\beta\right)\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right)c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	O V
$\mathbf{B}_{27}$	=	$(x_8 - y_8)\,\mathbf{a}_1 + (x_8 + y_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	OH
$\mathbf{B}_{28}$	=	$(-x_8 - y_8)\,\mathbf{a}_1 + (-x_8 + y_8)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_8 a - z_8 c\cos\beta\right)\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right)c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	OH
$\mathbf{B}_{29}$	=	$(-x_8 + y_8)\,\mathbf{a}_1 + (-x_8 - y_8)\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	OH
$\mathbf{B}_{30}$	=	$(x_8 + y_8)\,\mathbf{a}_1 + (x_8 - y_8)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_8 a + z_8 c\cos\beta\right)\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right)c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	OH
$\mathbf{B}_{31}$	=	$(x_9 - y_9)\,\mathbf{a}_1 + (x_9 + y_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$(x_9 a + z_9 c\cos\beta)\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	Si I
$\mathbf{B}_{32}$	=	$(-x_9 - y_9)\,\mathbf{a}_1 + (-x_9 + y_9)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_9 a - z_9 c\cos\beta\right)\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_9\right)c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	Si I
$\mathbf{B}_{33}$	=	$(-x_9 + y_9)\,\mathbf{a}_1 + (-x_9 - y_9)\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$(-x_9 a - z_9 c\cos\beta)\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} - z_9 c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	Si I
$\mathbf{B}_{34}$	=	$(x_9 + y_9)\,\mathbf{a}_1 + (x_9 - y_9)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_9 a + z_9 c\cos\beta\right)\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right)c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	Si I
$\mathbf{B}_{35}$	=	$(x_{10} - y_{10})\,\mathbf{a}_1 + (x_{10} + y_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$(x_{10} a + z_{10} c\cos\beta)\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	Si II
$\mathbf{B}_{36}$	=	$(-x_{10} - y_{10})\,\mathbf{a}_1 + (-x_{10} + y_{10})\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{10}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_{10} a - z_{10} c\cos\beta\right)\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{10}\right)c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	Si II
$\mathbf{B}_{37}$	=	$(-x_{10} + y_{10})\,\mathbf{a}_1 + (-x_{10} - y_{10})\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3$	=	$(-x_{10} a - z_{10} c\cos\beta)\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} - z_{10} c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	Si II
$\mathbf{B}_{38}$	=	$(x_{10} + y_{10})\,\mathbf{a}_1 + (x_{10} - y_{10})\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{10}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{10} a + z_{10} c\cos\beta\right)\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{10}\right)c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	Si II

References:

- S. M. Richardson and J. W. Richardson, Jr., *Crystal structure of a pink muscovite from Archer's Post, Kenya: Implications for reverse pleochroism in dioctahedral micas*, Am. Mineral. **67**, 69–75 (1982).

Found in:

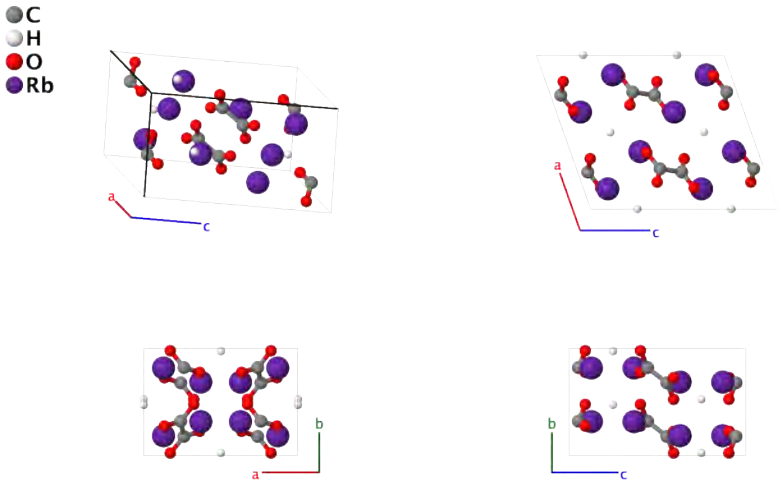
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1562
- POSCAR: pp. 1562

$Rb_2C_2O_4{\cdot}H_2O$ Structure: A2BC4D2_mC36_15_f_e_2f_f

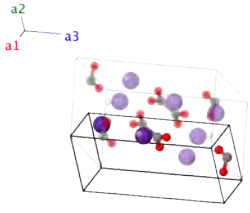
http://aflow.org/prototype-encyclopedia/A2BC4D2_mC36_15_f_e_2f_f



Prototype	:	$C_2(H_2O)O_4Rb_2$
AFLOW prototype label	:	A2BC4D2_mC36_15_f_e_2f_f
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC36
Space group number	:	15
Space group symbol	:	$C2/c$
AFLOW prototype command	:	`aflow --proto=A2BC4D2_mC36_15_f_e_2f_f --params=`$a, b/a, c/a, \beta, y_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5$

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H_2O
$\mathbf{B}_2$	=	$y_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	H_2O
$\mathbf{B}_3$	=	$(x_2 - y_2)\,\mathbf{a}_1 + (x_2 + y_2)\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	C
$\mathbf{B}_4$	=	$(-x_2 - y_2)\,\mathbf{a}_1 + (-x_2 + y_2)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_2 a - z_2 c\cos\beta\right)\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	C

$\mathbf{B}_5$	=	$(-x_2+y_2)\,\mathbf{a}_1+(-x_2-y_2)\,\mathbf{a}_2-z_2\,\mathbf{a}_3$	=	$(-x_2a-z_2c\cos\beta)\,\hat{\mathbf{x}}-y_2b\,\hat{\mathbf{y}}-z_2c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	C
$\mathbf{B}_6$	=	$(x_2+y_2)\,\mathbf{a}_1+(x_2-y_2)\,\mathbf{a}_2+\left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta+x_2a+z_2c\cos\beta\right)\hat{\mathbf{x}}-y_2b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	C
$\mathbf{B}_7$	=	$(x_3-y_3)\,\mathbf{a}_1+(x_3+y_3)\,\mathbf{a}_2+z_3\,\mathbf{a}_3$	=	$(x_3a+z_3c\cos\beta)\,\hat{\mathbf{x}}+y_3b\,\hat{\mathbf{y}}+z_3c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O I
$\mathbf{B}_8$	=	$(-x_3-y_3)\,\mathbf{a}_1+(-x_3+y_3)\,\mathbf{a}_2+\left(\frac{1}{2}-z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta-x_3a-z_3c\cos\beta\right)\hat{\mathbf{x}}+y_3b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O I
$\mathbf{B}_9$	=	$(-x_3+y_3)\,\mathbf{a}_1+(-x_3-y_3)\,\mathbf{a}_2-z_3\,\mathbf{a}_3$	=	$(-x_3a-z_3c\cos\beta)\,\hat{\mathbf{x}}-y_3b\,\hat{\mathbf{y}}-z_3c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O I
$\mathbf{B}_{10}$	=	$(x_3+y_3)\,\mathbf{a}_1+(x_3-y_3)\,\mathbf{a}_2+\left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta+x_3a+z_3c\cos\beta\right)\hat{\mathbf{x}}-y_3b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O I
$\mathbf{B}_{11}$	=	$(x_4-y_4)\,\mathbf{a}_1+(x_4+y_4)\,\mathbf{a}_2+z_4\,\mathbf{a}_3$	=	$(x_4a+z_4c\cos\beta)\,\hat{\mathbf{x}}+y_4b\,\hat{\mathbf{y}}+z_4c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{12}$	=	$(-x_4-y_4)\,\mathbf{a}_1+(-x_4+y_4)\,\mathbf{a}_2+\left(\frac{1}{2}-z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta-x_4a-z_4c\cos\beta\right)\hat{\mathbf{x}}+y_4b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{13}$	=	$(-x_4+y_4)\,\mathbf{a}_1+(-x_4-y_4)\,\mathbf{a}_2-z_4\,\mathbf{a}_3$	=	$(-x_4a-z_4c\cos\beta)\,\hat{\mathbf{x}}-y_4b\,\hat{\mathbf{y}}-z_4c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{14}$	=	$(x_4+y_4)\,\mathbf{a}_1+(x_4-y_4)\,\mathbf{a}_2+\left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta+x_4a+z_4c\cos\beta\right)\hat{\mathbf{x}}-y_4b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{15}$	=	$(x_5-y_5)\,\mathbf{a}_1+(x_5+y_5)\,\mathbf{a}_2+z_5\,\mathbf{a}_3$	=	$(x_5a+z_5c\cos\beta)\,\hat{\mathbf{x}}+y_5b\,\hat{\mathbf{y}}+z_5c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Rb
$\mathbf{B}_{16}$	=	$(-x_5-y_5)\,\mathbf{a}_1+(-x_5+y_5)\,\mathbf{a}_2+\left(\frac{1}{2}-z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta-x_5a-z_5c\cos\beta\right)\hat{\mathbf{x}}+y_5b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Rb
$\mathbf{B}_{17}$	=	$(-x_5+y_5)\,\mathbf{a}_1+(-x_5-y_5)\,\mathbf{a}_2-z_5\,\mathbf{a}_3$	=	$(-x_5a-z_5c\cos\beta)\,\hat{\mathbf{x}}-y_5b\,\hat{\mathbf{y}}-z_5c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Rb
$\mathbf{B}_{18}$	=	$(x_5+y_5)\,\mathbf{a}_1+(x_5-y_5)\,\mathbf{a}_2+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta+x_5a+z_5c\cos\beta\right)\hat{\mathbf{x}}-y_5b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Rb

References:

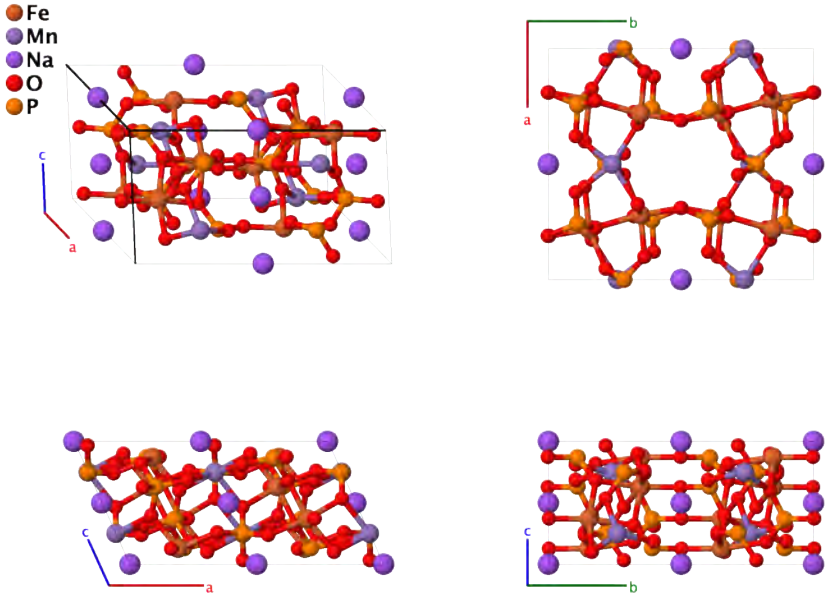
- B. F. Pedersen, *The Crystal Structure of Rubidium. Oxalate Monohydrate, $Rb_2C_2O_4{\cdot}H_2O$*, Acta Chem. Scand. **19**, 1815–1818 (1965), doi:10.3891/acta.chem.scand.19-1815.

Geometry files:

- CIF: pp. 1563
- POSCAR: pp. 1563

Alluaudite [$NaMnFe_2(PO_4)_3$] Structure: A2BCD12E3_mC76_15_f_e_b_6f_ef

http://aflow.org/prototype-encyclopedia/A2BCD12E3_mC76_15_f_e_b_6f_ef



Prototype	:	$Fe_2MnNaO_{12}P_3$
AFLOW prototype label	:	A2BCD12E3_mC76_15_f_e_b_6f_ef
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC76
Space group number	:	15
Space group symbol	:	$C2/c$
AFLOW prototype command	:	`aflow --proto=A2BCD12E3_mC76_15_f_e_b_6f_ef --params=`$a, b/a, c/a, \beta, y_2, y_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}$

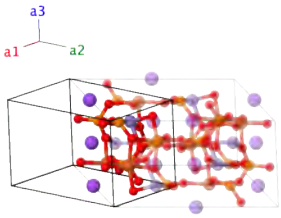
Other compounds with this structure

- $Ag_xNa_{1-x}Mn_3(PO_4)_3$, $Li_xNa_{1-x}CdIn_2(PO_4)_3$, $Li_xNa_{1-x}MnFe_2(PO_4)_3$, $Na_2(Fe,Co)Fe(VO_4)_3$, $Na_2Co_2Cr(PO_4)_3$, $Na_2Fe_2V(PO_4)_3$, $Na_2Zn_2Fe(VO_4)_3$, $Na_3Bi_2(AsO_4)_3$, $Na_4Co(MnO_4)_3$, and $Na_xMn_yAl_{5-x-y}(PO_4)_3$

- This is a rather idealized version of alluaudite. (Moore, 1971) gives the composition of the sodium site (Na) as $Na_{0.625}Mn_{0.175}Ca_{0.125}$, with vacancies on the remaining sites; the manganese (Mn) site is $Mn_{0.950}Mg_{0.025}Li_{0.025}$; and the iron (Fe) site as $Fe_{0.988}Mg_{0.012}$. (Hatert, 2005) puts some sodium atoms on another ($4e$) site, with $y \approx 0$; the exact placement and concentration depends upon the manganese content. We may expect similar variations in other compounds having this structure.

Base-centered Monoclinic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2}\,b\,\hat{\mathbf{y}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,b\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	$(4b)$	Na
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(a + c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	$(4b)$	Na
$\mathbf{B}_3$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Mn
$\mathbf{B}_4$	=	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Mn
$\mathbf{B}_5$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	P I
$\mathbf{B}_6$	=	$y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	P I
$\mathbf{B}_7$	=	$(x_4 - y_4)\,\mathbf{a}_1 + (x_4 + y_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Fe
$\mathbf{B}_8$	=	$(-x_4 - y_4)\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_4 a - z_4 c\cos\beta\right)\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Fe
$\mathbf{B}_9$	=	$(-x_4 + y_4)\,\mathbf{a}_1 + (-x_4 - y_4)\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Fe
$\mathbf{B}_{10}$	=	$(x_4 + y_4)\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Fe
$\mathbf{B}_{11}$	=	$(x_5 - y_5)\,\mathbf{a}_1 + (x_5 + y_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O I
$\mathbf{B}_{12}$	=	$(-x_5 - y_5)\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_5 a - z_5 c\cos\beta\right)\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O I
$\mathbf{B}_{13}$	=	$(-x_5 + y_5)\,\mathbf{a}_1 + (-x_5 - y_5)\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O I
$\mathbf{B}_{14}$	=	$(x_5 + y_5)\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O I
$\mathbf{B}_{15}$	=	$(x_6 - y_6)\,\mathbf{a}_1 + (x_6 + y_6)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{16}$	=	$(-x_6 - y_6)\,\mathbf{a}_1 + (-x_6 + y_6)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_6 a - z_6 c\cos\beta\right)\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{17}$	=	$(-x_6 + y_6)\,\mathbf{a}_1 + (-x_6 - y_6)\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{18}$	=	$(x_6 + y_6)\,\mathbf{a}_1 + (x_6 - y_6)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{19}$	=	$(x_7 - y_7)\,\mathbf{a}_1 + (x_7 + y_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III

$\mathbf{B}_{20}$	=	$(-x_7 - y_7)\,\mathbf{a}_1 + (-x_7 + y_7)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_7 a - z_7 c\cos\beta\right)\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{21}$	=	$(-x_7 + y_7)\,\mathbf{a}_1 + (-x_7 - y_7)\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{22}$	=	$(x_7 + y_7)\,\mathbf{a}_1 + (x_7 - y_7)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_7 a + z_7 c\cos\beta\right)\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{23}$	=	$(x_8 - y_8)\,\mathbf{a}_1 + (x_8 + y_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O IV
$\mathbf{B}_{24}$	=	$(-x_8 - y_8)\,\mathbf{a}_1 + (-x_8 + y_8)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_8 a - z_8 c\cos\beta\right)\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O IV
$\mathbf{B}_{25}$	=	$(-x_8 + y_8)\,\mathbf{a}_1 + (-x_8 - y_8)\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O IV
$\mathbf{B}_{26}$	=	$(x_8 + y_8)\,\mathbf{a}_1 + (x_8 - y_8)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_8 a + z_8 c\cos\beta\right)\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O IV
$\mathbf{B}_{27}$	=	$(x_9 - y_9)\,\mathbf{a}_1 + (x_9 + y_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$(x_9 a + z_9 c\cos\beta)\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O V
$\mathbf{B}_{28}$	=	$(-x_9 - y_9)\,\mathbf{a}_1 + (-x_9 + y_9)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_9 a - z_9 c\cos\beta\right)\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_9\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O V
$\mathbf{B}_{29}$	=	$(-x_9 + y_9)\,\mathbf{a}_1 + (-x_9 - y_9)\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$(-x_9 a - z_9 c\cos\beta)\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} - z_9 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O V
$\mathbf{B}_{30}$	=	$(x_9 + y_9)\,\mathbf{a}_1 + (x_9 - y_9)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_9 a + z_9 c\cos\beta\right)\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O V
$\mathbf{B}_{31}$	=	$(x_{10} - y_{10})\,\mathbf{a}_1 + (x_{10} + y_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$(x_{10} a + z_{10} c\cos\beta)\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O VI
$\mathbf{B}_{32}$	=	$(-x_{10} - y_{10})\,\mathbf{a}_1 + (-x_{10} + y_{10})\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{10}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_{10} a - z_{10} c\cos\beta\right)\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{10}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O VI
$\mathbf{B}_{33}$	=	$(-x_{10} + y_{10})\,\mathbf{a}_1 + (-x_{10} - y_{10})\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3$	=	$(-x_{10} a - z_{10} c\cos\beta)\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} - z_{10} c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O VI
$\mathbf{B}_{34}$	=	$(x_{10} + y_{10})\,\mathbf{a}_1 + (x_{10} - y_{10})\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{10}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{10} a + z_{10} c\cos\beta\right)\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{10}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O VI
$\mathbf{B}_{35}$	=	$(x_{11} - y_{11})\,\mathbf{a}_1 + (x_{11} + y_{11})\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$(x_{11} a + z_{11} c\cos\beta)\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + z_{11} c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	P II
$\mathbf{B}_{36}$	=	$(-x_{11} - y_{11})\,\mathbf{a}_1 + (-x_{11} + y_{11})\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{11}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_{11} a - z_{11} c\cos\beta\right)\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{11}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	P II
$\mathbf{B}_{37}$	=	$(-x_{11} + y_{11})\,\mathbf{a}_1 + (-x_{11} - y_{11})\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3$	=	$(-x_{11} a - z_{11} c\cos\beta)\,\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} - z_{11} c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	P II
$\mathbf{B}_{38}$	=	$(x_{11} + y_{11})\,\mathbf{a}_1 + (x_{11} - y_{11})\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{11}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{11} a + z_{11} c\cos\beta\right)\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{11}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	P II

References:

- P. B. Moore, *Crystal Chemistry of the Alluaudite Structure Type: Contribution to the Paragenesis of Pegmatite Phosphate Giant Crystals*, Am. Mineral. **56**, 1955–1975 (1971).
- F. Hatert, L. Rebbouh, R. P. Hermann, A.-M. Fransolet, G. J. Long, and F. Grandjean, *Crystal chemistry of the hydrothermally synthesized* $Na_2(Mn_{1-x}Fe_x^{2+})_2Fe^{3+}(PO_4)_3$ *alluaudite-type solid solution*, Am. Mineral. **90**, 653–662 (2005), doi:10.2138/am.2005.1551.

Found in:

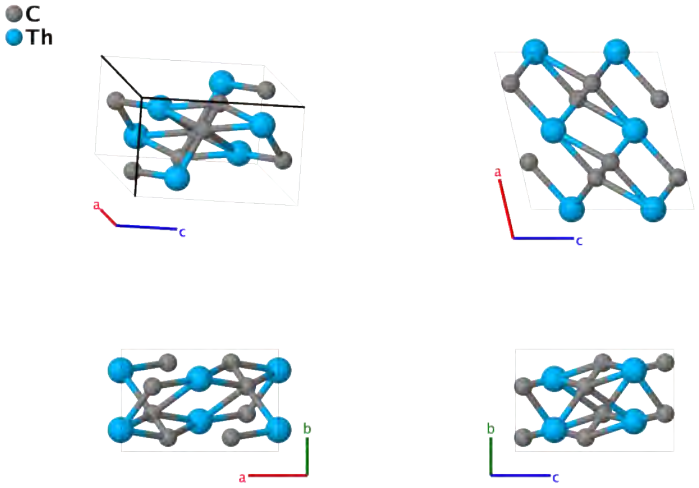
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1563
- POSCAR: pp. 1564

ThC_2 (C_g) Structure: A2B_mC12_15_f_e

http://aflow.org/prototype-encyclopedia/A2B_mC12_15_f_e



Prototype	:	C_2Th
AFLOW prototype label	:	A2B_mC12_15_f_e
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC12
Space group number	:	15
Space group symbol	:	$C2/c$
AFLOW prototype command	:	`aflow --proto=A2B_mC12_15_f_e --params=`$a, b/a, c/a, \beta, y_1, x_2, y_2, z_2$

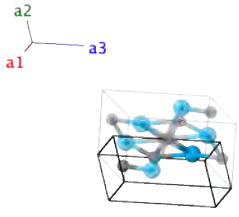
Other compounds with this structure

- YbS_2 and BaS_2

- This is a metastable phase of CaC_2. The stable room-temperature phase is the $C11_a$ (A2B_tI6_139_e_a) structure.

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$ =	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Th

$\mathbf{B}_2$	=	$y_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Th
$\mathbf{B}_3$	=	$(x_2 - y_2)\,\mathbf{a}_1 + (x_2 + y_2)\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	C
$\mathbf{B}_4$	=	$(-x_2 - y_2)\,\mathbf{a}_1 + (-x_2 + y_2)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_2 a - z_2 c\cos\beta\right)\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	C
$\mathbf{B}_5$	=	$(-x_2 + y_2)\,\mathbf{a}_1 + (-x_2 - y_2)\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} - z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	C
$\mathbf{B}_6$	=	$(x_2 + y_2)\,\mathbf{a}_1 + (x_2 - y_2)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_2 a + z_2 c\cos\beta\right)\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	C

References:

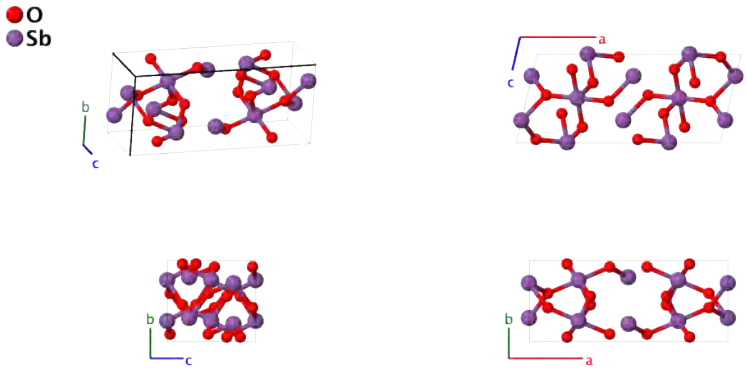
- A. L. Bowman, N. H. Krikorian, G. P. Arnold, T. C. Wallace, and N. G. Nereson, *The Crystal Structure of ThC_2*, Acta Crystallogr. Sect. B Struct. Sci. **24**, 1121–1123 (1968), doi:10.1107/S056774086800378X.

Geometry files:

- CIF: pp. 1564
- POSCAR: pp. 1564

Clinocervantite (β-Sb_2O_4) Structure: A2B_mC24_15_2f_ce

http://aflow.org/prototype-encyclopedia/A2B_mC24_15_2f_ce

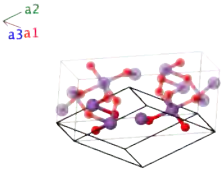


Prototype	:	O_2Sb
AFLOW prototype label	:	A2B_mC24_15_2f_ce
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC24
Space group number	:	15
Space group symbol	:	$C2/c$
AFLOW prototype command	:	`aflow --proto=A2B_mC24_15_2f_ce --params=`$a, b/a, c/a, \beta, y_2, x_3, y_3, z_3, x_4, y_4, z_4$

- This is *not* the $D6_2$ structure of SbO_2, which was found to be erroneous. Clinocervantite is a naturally occurring monoclinic modification of cervantite (α-Sb_2O_4)

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}}$	$(4c)$	Sb I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{4}a + \frac{1}{2}c\cos\beta\right)\hat{\mathbf{x}} - \frac{1}{4}b\,\hat{\mathbf{y}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	$(4c)$	Sb I
$\mathbf{B}_3$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Sb II
$\mathbf{B}_4$	=	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Sb II
$\mathbf{B}_5$	=	$(x_3 - y_3)\,\mathbf{a}_1 + (x_3 + y_3)\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O I

$\mathbf{B}_6$	=	$(-x_3 - y_3)\,\mathbf{a}_1 + (-x_3 + y_3)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_3 a - z_3 c\cos\beta\right)\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O I
$\mathbf{B}_7$	=	$(-x_3 + y_3)\,\mathbf{a}_1 + (-x_3 - y_3)\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O I
$\mathbf{B}_8$	=	$(x_3 + y_3)\,\mathbf{a}_1 + (x_3 - y_3)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O I
$\mathbf{B}_9$	=	$(x_4 - y_4)\,\mathbf{a}_1 + (x_4 + y_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{10}$	=	$(-x_4 - y_4)\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_4 a - z_4 c\cos\beta\right)\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{11}$	=	$(-x_4 + y_4)\,\mathbf{a}_1 + (-x_4 - y_4)\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{12}$	=	$(x_4 + y_4)\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II

References:

- R. Basso, G. Lucchetti, L. Zefiro, and A. Palenzona, *Clinocervantite, β-Sb_2O_4, the natural monoclinic polymorph of cervantite from the Cetine mine, Siena, Italy*, Eur. J. Mineral. **11**, 95–100 (1999), doi:10.1127/ejm/11/1/0095.

Found in:

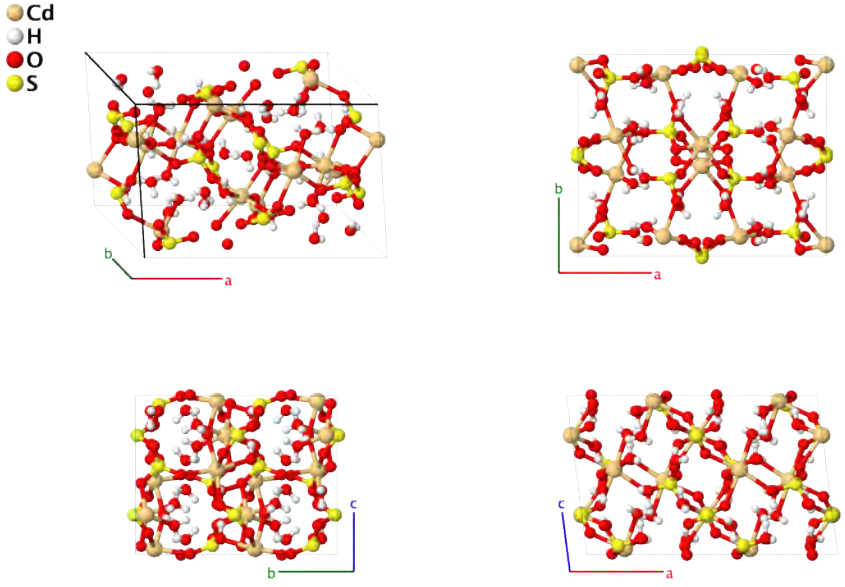
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1564
- POSCAR: pp. 1565

$(CdSO_4)_3 \cdot 8H_2O$ ($H4_{20}$) Structure: A3B16C20D3_mC168_15_ef_8f_10f_ef

http://aflow.org/prototype-encyclopedia/A3B16C20D3_mC168_15_ef_8f_10f_ef

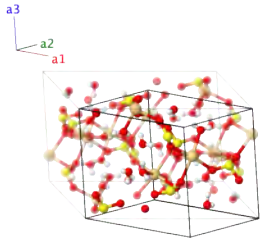


Prototype	:	$Cd_3H_{16}O_{20}S_3$
AFLOW prototype label	:	A3B16C20D3_mC168_15_ef_8f_10f_ef
***Strukturbericht* designation**	:	$H4_{20}$
Pearson symbol	:	mC168
Space group number	:	15
Space group symbol	:	$C2/c$
AFLOW prototype command	:	`aflow --proto=A3B16C20D3_mC168_15_ef_8f_10f_ef --params=`$a, b/a, c/a, \beta, y_1, y_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}, x_{17}, y_{17}, z_{17}, x_{18}, y_{18}, z_{18}, x_{19}, y_{19}, z_{19}, x_{20}, y_{20}, z_{20}, x_{21}, y_{21}, z_{21}, x_{22}, y_{22}, z_{22}$

- Some of the H-O distances appear to be very small. For example, the O-VIII – H-III distance is only 0.43Å. We have checked our inputs compared to (Caminiti, 1981) and found no error. Unfortunately they do not provide H-O distances for comparison.

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cd I
$\mathbf{B}_2$	=	$y_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Cd I
$\mathbf{B}_3$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S I
$\mathbf{B}_4$	=	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S I
$\mathbf{B}_5$	=	$(x_3 - y_3)\,\mathbf{a}_1 + (x_3 + y_3)\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Cd II
$\mathbf{B}_6$	=	$(-x_3 - y_3)\,\mathbf{a}_1 + (-x_3 + y_3)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_3 a - z_3 c\cos\beta\right)\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Cd II
$\mathbf{B}_7$	=	$(-x_3 + y_3)\,\mathbf{a}_1 + (-x_3 - y_3)\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Cd II
$\mathbf{B}_8$	=	$(x_3 + y_3)\,\mathbf{a}_1 + (x_3 - y_3)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Cd II
$\mathbf{B}_9$	=	$(x_4 - y_4)\,\mathbf{a}_1 + (x_4 + y_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H I
$\mathbf{B}_{10}$	=	$(-x_4 - y_4)\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_4 a - z_4 c\cos\beta\right)\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H I
$\mathbf{B}_{11}$	=	$(-x_4 + y_4)\,\mathbf{a}_1 + (-x_4 - y_4)\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H I
$\mathbf{B}_{12}$	=	$(x_4 + y_4)\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H I
$\mathbf{B}_{13}$	=	$(x_5 - y_5)\,\mathbf{a}_1 + (x_5 + y_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H II
$\mathbf{B}_{14}$	=	$(-x_5 - y_5)\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_5 a - z_5 c\cos\beta\right)\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H II
$\mathbf{B}_{15}$	=	$(-x_5 + y_5)\,\mathbf{a}_1 + (-x_5 - y_5)\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H II
$\mathbf{B}_{16}$	=	$(x_5 + y_5)\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H II
$\mathbf{B}_{17}$	=	$(x_6 - y_6)\,\mathbf{a}_1 + (x_6 + y_6)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H III
$\mathbf{B}_{18}$	=	$(-x_6 - y_6)\,\mathbf{a}_1 + (-x_6 + y_6)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_6 a - z_6 c\cos\beta\right)\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H III
$\mathbf{B}_{19}$	=	$(-x_6 + y_6)\,\mathbf{a}_1 + (-x_6 - y_6)\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H III
$\mathbf{B}_{20}$	=	$(x_6 + y_6)\,\mathbf{a}_1 + (x_6 - y_6)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H III
$\mathbf{B}_{21}$	=	$(x_7 - y_7)\,\mathbf{a}_1 + (x_7 + y_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H IV
$\mathbf{B}_{22}$	=	$(-x_7 - y_7)\,\mathbf{a}_1 + (-x_7 + y_7)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_7 a - z_7 c\cos\beta\right)\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H IV

$$\mathbf{B}_{23} = (-x_7 + y_7)\,\mathbf{a}_1 + (-x_7 - y_7)\,\mathbf{a}_2 - z_7\,\mathbf{a}_3 = (-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H IV}$$

$$\mathbf{B}_{24} = (x_7 + y_7)\,\mathbf{a}_1 + (x_7 - y_7)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_7 a + z_7 c\cos\beta\right)\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_7\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H IV}$$

$$\mathbf{B}_{25} = (x_8 - y_8)\,\mathbf{a}_1 + (x_8 + y_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 = (x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H V}$$

$$\mathbf{B}_{26} = (-x_8 - y_8)\,\mathbf{a}_1 + (-x_8 + y_8)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_8 a - z_8 c\cos\beta\right)\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_8\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H V}$$

$$\mathbf{B}_{27} = (-x_8 + y_8)\,\mathbf{a}_1 + (-x_8 - y_8)\,\mathbf{a}_2 - z_8\,\mathbf{a}_3 = (-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H V}$$

$$\mathbf{B}_{28} = (x_8 + y_8)\,\mathbf{a}_1 + (x_8 - y_8)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_8 a + z_8 c\cos\beta\right)\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_8\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H V}$$

$$\mathbf{B}_{29} = (x_9 - y_9)\,\mathbf{a}_1 + (x_9 + y_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 = (x_9 a + z_9 c\cos\beta)\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H VI}$$

$$\mathbf{B}_{30} = (-x_9 - y_9)\,\mathbf{a}_1 + (-x_9 + y_9)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_9 a - z_9 c\cos\beta\right)\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_9\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H VI}$$

$$\mathbf{B}_{31} = (-x_9 + y_9)\,\mathbf{a}_1 + (-x_9 - y_9)\,\mathbf{a}_2 - z_9\,\mathbf{a}_3 = (-x_9 a - z_9 c\cos\beta)\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} - z_9 c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H VI}$$

$$\mathbf{B}_{32} = (x_9 + y_9)\,\mathbf{a}_1 + (x_9 - y_9)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_9 a + z_9 c\cos\beta\right)\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_9\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H VI}$$

$$\mathbf{B}_{33} = (x_{10} - y_{10})\,\mathbf{a}_1 + (x_{10} + y_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 = (x_{10} a + z_{10} c\cos\beta)\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H VII}$$

$$\mathbf{B}_{34} = (-x_{10} - y_{10})\,\mathbf{a}_1 + (-x_{10} + y_{10})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{10} a - z_{10} c\cos\beta\right)\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{10}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H VII}$$

$$\mathbf{B}_{35} = (-x_{10} + y_{10})\,\mathbf{a}_1 + (-x_{10} - y_{10})\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 = (-x_{10} a - z_{10} c\cos\beta)\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} - z_{10} c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H VII}$$

$$\mathbf{B}_{36} = (x_{10} + y_{10})\,\mathbf{a}_1 + (x_{10} - y_{10})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{10} a + z_{10} c\cos\beta\right)\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{10}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H VII}$$

$$\mathbf{B}_{37} = (x_{11} - y_{11})\,\mathbf{a}_1 + (x_{11} + y_{11})\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 = (x_{11} a + z_{11} c\cos\beta)\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + z_{11} c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H VIII}$$

$$\mathbf{B}_{38} = (-x_{11} - y_{11})\,\mathbf{a}_1 + (-x_{11} + y_{11})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{11} a - z_{11} c\cos\beta\right)\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{11}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H VIII}$$

$$\mathbf{B}_{39} = (-x_{11} + y_{11})\,\mathbf{a}_1 + (-x_{11} - y_{11})\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3 = (-x_{11} a - z_{11} c\cos\beta)\,\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} - z_{11} c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H VIII}$$

$$\mathbf{B}_{40} = (x_{11} + y_{11})\,\mathbf{a}_1 + (x_{11} - y_{11})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{11} a + z_{11} c\cos\beta\right)\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{11}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H VIII}$$

$$\mathbf{B}_{41} = (x_{12} - y_{12})\,\mathbf{a}_1 + (x_{12} + y_{12})\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 = (x_{12} a + z_{12} c\cos\beta)\,\hat{\mathbf{x}} + y_{12} b\,\hat{\mathbf{y}} + z_{12} c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O I}$$

$$\mathbf{B}_{42} = (-x_{12} - y_{12})\,\mathbf{a}_1 + (-x_{12} + y_{12})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{12} a - z_{12} c\cos\beta\right)\hat{\mathbf{x}} + y_{12} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{12}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O I}$$

$$\mathbf{B}_{43} = (-x_{12} + y_{12})\,\mathbf{a}_1 + (-x_{12} - y_{12})\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3 = (-x_{12} a - z_{12} c\cos\beta)\,\hat{\mathbf{x}} - y_{12} b\,\hat{\mathbf{y}} - z_{12} c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O I}$$

$$\mathbf{B}_{44} = (x_{12} + y_{12})\,\mathbf{a}_1 + (x_{12} - y_{12})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{12} a + z_{12} c\cos\beta\right)\hat{\mathbf{x}} - y_{12} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{12}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O I}$$

$$\mathbf{B}_{45} = (x_{13} - y_{13})\,\mathbf{a}_1 + (x_{13} + y_{13})\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3 = (x_{13}a + z_{13}c\cos\beta)\,\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} + z_{13}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O II}$$

$$\mathbf{B}_{46} = (-x_{13} - y_{13})\,\mathbf{a}_1 + (-x_{13} + y_{13})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{13}a - z_{13}c\cos\beta\right)\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{13}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O II}$$

$$\mathbf{B}_{47} = (-x_{13} + y_{13})\,\mathbf{a}_1 + (-x_{13} - y_{13})\,\mathbf{a}_2 - z_{13}\,\mathbf{a}_3 = (-x_{13}a - z_{13}c\cos\beta)\,\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} - z_{13}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O II}$$

$$\mathbf{B}_{48} = (x_{13} + y_{13})\,\mathbf{a}_1 + (x_{13} - y_{13})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{13}a + z_{13}c\cos\beta\right)\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{13}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O II}$$

$$\mathbf{B}_{49} = (x_{14} - y_{14})\,\mathbf{a}_1 + (x_{14} + y_{14})\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3 = (x_{14}a + z_{14}c\cos\beta)\,\hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} + z_{14}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O III}$$

$$\mathbf{B}_{50} = (-x_{14} - y_{14})\,\mathbf{a}_1 + (-x_{14} + y_{14})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{14}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{14}a - z_{14}c\cos\beta\right)\hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{14}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O III}$$

$$\mathbf{B}_{51} = (-x_{14} + y_{14})\,\mathbf{a}_1 + (-x_{14} - y_{14})\,\mathbf{a}_2 - z_{14}\,\mathbf{a}_3 = (-x_{14}a - z_{14}c\cos\beta)\,\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} - z_{14}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O III}$$

$$\mathbf{B}_{52} = (x_{14} + y_{14})\,\mathbf{a}_1 + (x_{14} - y_{14})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{14}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{14}a + z_{14}c\cos\beta\right)\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{14}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O III}$$

$$\mathbf{B}_{53} = (x_{15} - y_{15})\,\mathbf{a}_1 + (x_{15} + y_{15})\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 = (x_{15}a + z_{15}c\cos\beta)\,\hat{\mathbf{x}} + y_{15}b\,\hat{\mathbf{y}} + z_{15}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O IV}$$

$$\mathbf{B}_{54} = (-x_{15} - y_{15})\,\mathbf{a}_1 + (-x_{15} + y_{15})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{15}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{15}a - z_{15}c\cos\beta\right)\hat{\mathbf{x}} + y_{15}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{15}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O IV}$$

$$\mathbf{B}_{55} = (-x_{15} + y_{15})\,\mathbf{a}_1 + (-x_{15} - y_{15})\,\mathbf{a}_2 - z_{15}\,\mathbf{a}_3 = (-x_{15}a - z_{15}c\cos\beta)\,\hat{\mathbf{x}} - y_{15}b\,\hat{\mathbf{y}} - z_{15}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O IV}$$

$$\mathbf{B}_{56} = (x_{15} + y_{15})\,\mathbf{a}_1 + (x_{15} - y_{15})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{15}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{15}a + z_{15}c\cos\beta\right)\hat{\mathbf{x}} - y_{15}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{15}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O IV}$$

$$\mathbf{B}_{57} = (x_{16} - y_{16})\,\mathbf{a}_1 + (x_{16} + y_{16})\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 = (x_{16}a + z_{16}c\cos\beta)\,\hat{\mathbf{x}} + y_{16}b\,\hat{\mathbf{y}} + z_{16}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O V}$$

$$\mathbf{B}_{58} = (-x_{16} - y_{16})\,\mathbf{a}_1 + (-x_{16} + y_{16})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{16}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{16}a - z_{16}c\cos\beta\right)\hat{\mathbf{x}} + y_{16}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{16}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O V}$$

$$\mathbf{B}_{59} = (-x_{16} + y_{16})\,\mathbf{a}_1 + (-x_{16} - y_{16})\,\mathbf{a}_2 - z_{16}\,\mathbf{a}_3 = (-x_{16}a - z_{16}c\cos\beta)\,\hat{\mathbf{x}} - y_{16}b\,\hat{\mathbf{y}} - z_{16}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O V}$$

$$\mathbf{B}_{60} = (x_{16} + y_{16})\,\mathbf{a}_1 + (x_{16} - y_{16})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{16}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{16}a + z_{16}c\cos\beta\right)\hat{\mathbf{x}} - y_{16}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{16}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O V}$$

$$\mathbf{B}_{61} = (x_{17} - y_{17})\,\mathbf{a}_1 + (x_{17} + y_{17})\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 = (x_{17}a + z_{17}c\cos\beta)\,\hat{\mathbf{x}} + y_{17}b\,\hat{\mathbf{y}} + z_{17}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VI}$$

$$\mathbf{B}_{62} = (-x_{17} - y_{17})\,\mathbf{a}_1 + (-x_{17} + y_{17})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{17}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{17}a - z_{17}c\cos\beta\right)\hat{\mathbf{x}} + y_{17}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{17}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VI}$$

$$\mathbf{B}_{63} = (-x_{17} + y_{17})\,\mathbf{a}_1 + (-x_{17} - y_{17})\,\mathbf{a}_2 - z_{17}\,\mathbf{a}_3 = (-x_{17}a - z_{17}c\cos\beta)\,\hat{\mathbf{x}} - y_{17}b\,\hat{\mathbf{y}} - z_{17}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VI}$$

$$\mathbf{B}_{64} = (x_{17} + y_{17})\,\mathbf{a}_1 + (x_{17} - y_{17})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{17}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{17}a + z_{17}c\cos\beta\right)\hat{\mathbf{x}} - y_{17}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{17}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VI}$$

$$\mathbf{B}_{65} = (x_{18} - y_{18})\,\mathbf{a}_1 + (x_{18} + y_{18})\,\mathbf{a}_2 + z_{18}\,\mathbf{a}_3 = (x_{18}a + z_{18}c\cos\beta)\,\hat{\mathbf{x}} + y_{18}b\,\hat{\mathbf{y}} + z_{18}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VII}$$

$$\mathbf{B}_{66} = (-x_{18} - y_{18})\,\mathbf{a}_1 + (-x_{18} + y_{18})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{18}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{18}a - z_{18}c\cos\beta\right)\hat{\mathbf{x}} + y_{18}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{18}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VII}$$

$\mathbf{B}_{67}$	=	$(-x_{18}+y_{18})\,\mathbf{a}_1 + (-x_{18}-y_{18})\,\mathbf{a}_2 - z_{18}\,\mathbf{a}_3$	=	$(-x_{18}a - z_{18}c\cos\beta)\,\hat{\mathbf{x}} - y_{18}b\,\hat{\mathbf{y}} - z_{18}c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O VII
$\mathbf{B}_{68}$	=	$(x_{18}+y_{18})\,\mathbf{a}_1 + (x_{18}-y_{18})\,\mathbf{a}_2 + \left(\frac{1}{2}+z_{18}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{18}a + z_{18}c\cos\beta\right)\hat{\mathbf{x}} - y_{18}b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_{18}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O VII
$\mathbf{B}_{69}$	=	$(x_{19}-y_{19})\,\mathbf{a}_1 + (x_{19}+y_{19})\,\mathbf{a}_2 + z_{19}\,\mathbf{a}_3$	=	$(x_{19}a + z_{19}c\cos\beta)\,\hat{\mathbf{x}} + y_{19}b\,\hat{\mathbf{y}} + z_{19}c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O VIII
$\mathbf{B}_{70}$	=	$(-x_{19}-y_{19})\,\mathbf{a}_1 + (-x_{19}+y_{19})\,\mathbf{a}_2 + \left(\frac{1}{2}-z_{19}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_{19}a - z_{19}c\cos\beta\right)\hat{\mathbf{x}} + y_{19}b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_{19}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O VIII
$\mathbf{B}_{71}$	=	$(-x_{19}+y_{19})\,\mathbf{a}_1 + (-x_{19}-y_{19})\,\mathbf{a}_2 - z_{19}\,\mathbf{a}_3$	=	$(-x_{19}a - z_{19}c\cos\beta)\,\hat{\mathbf{x}} - y_{19}b\,\hat{\mathbf{y}} - z_{19}c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O VIII
$\mathbf{B}_{72}$	=	$(x_{19}+y_{19})\,\mathbf{a}_1 + (x_{19}-y_{19})\,\mathbf{a}_2 + \left(\frac{1}{2}+z_{19}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{19}a + z_{19}c\cos\beta\right)\hat{\mathbf{x}} - y_{19}b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_{19}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O VIII
$\mathbf{B}_{73}$	=	$(x_{20}-y_{20})\,\mathbf{a}_1 + (x_{20}+y_{20})\,\mathbf{a}_2 + z_{20}\,\mathbf{a}_3$	=	$(x_{20}a + z_{20}c\cos\beta)\,\hat{\mathbf{x}} + y_{20}b\,\hat{\mathbf{y}} + z_{20}c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O IX
$\mathbf{B}_{74}$	=	$(-x_{20}-y_{20})\,\mathbf{a}_1 + (-x_{20}+y_{20})\,\mathbf{a}_2 + \left(\frac{1}{2}-z_{20}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_{20}a - z_{20}c\cos\beta\right)\hat{\mathbf{x}} + y_{20}b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_{20}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O IX
$\mathbf{B}_{75}$	=	$(-x_{20}+y_{20})\,\mathbf{a}_1 + (-x_{20}-y_{20})\,\mathbf{a}_2 - z_{20}\,\mathbf{a}_3$	=	$(-x_{20}a - z_{20}c\cos\beta)\,\hat{\mathbf{x}} - y_{20}b\,\hat{\mathbf{y}} - z_{20}c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O IX
$\mathbf{B}_{76}$	=	$(x_{20}+y_{20})\,\mathbf{a}_1 + (x_{20}-y_{20})\,\mathbf{a}_2 + \left(\frac{1}{2}+z_{20}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{20}a + z_{20}c\cos\beta\right)\hat{\mathbf{x}} - y_{20}b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_{20}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O IX
$\mathbf{B}_{77}$	=	$(x_{21}-y_{21})\,\mathbf{a}_1 + (x_{21}+y_{21})\,\mathbf{a}_2 + z_{21}\,\mathbf{a}_3$	=	$(x_{21}a + z_{21}c\cos\beta)\,\hat{\mathbf{x}} + y_{21}b\,\hat{\mathbf{y}} + z_{21}c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O X
$\mathbf{B}_{78}$	=	$(-x_{21}-y_{21})\,\mathbf{a}_1 + (-x_{21}+y_{21})\,\mathbf{a}_2 + \left(\frac{1}{2}-z_{21}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_{21}a - z_{21}c\cos\beta\right)\hat{\mathbf{x}} + y_{21}b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_{21}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O X
$\mathbf{B}_{79}$	=	$(-x_{21}+y_{21})\,\mathbf{a}_1 + (-x_{21}-y_{21})\,\mathbf{a}_2 - z_{21}\,\mathbf{a}_3$	=	$(-x_{21}a - z_{21}c\cos\beta)\,\hat{\mathbf{x}} - y_{21}b\,\hat{\mathbf{y}} - z_{21}c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O X
$\mathbf{B}_{80}$	=	$(x_{21}+y_{21})\,\mathbf{a}_1 + (x_{21}-y_{21})\,\mathbf{a}_2 + \left(\frac{1}{2}+z_{21}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{21}a + z_{21}c\cos\beta\right)\hat{\mathbf{x}} - y_{21}b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_{21}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O X
$\mathbf{B}_{81}$	=	$(x_{22}-y_{22})\,\mathbf{a}_1 + (x_{22}+y_{22})\,\mathbf{a}_2 + z_{22}\,\mathbf{a}_3$	=	$(x_{22}a + z_{22}c\cos\beta)\,\hat{\mathbf{x}} + y_{22}b\,\hat{\mathbf{y}} + z_{22}c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	S II
$\mathbf{B}_{82}$	=	$(-x_{22}-y_{22})\,\mathbf{a}_1 + (-x_{22}+y_{22})\,\mathbf{a}_2 + \left(\frac{1}{2}-z_{22}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_{22}a - z_{22}c\cos\beta\right)\hat{\mathbf{x}} + y_{22}b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_{22}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	S II
$\mathbf{B}_{83}$	=	$(-x_{22}+y_{22})\,\mathbf{a}_1 + (-x_{22}-y_{22})\,\mathbf{a}_2 - z_{22}\,\mathbf{a}_3$	=	$(-x_{22}a - z_{22}c\cos\beta)\,\hat{\mathbf{x}} - y_{22}b\,\hat{\mathbf{y}} - z_{22}c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	S II
$\mathbf{B}_{84}$	=	$(x_{22}+y_{22})\,\mathbf{a}_1 + (x_{22}-y_{22})\,\mathbf{a}_2 + \left(\frac{1}{2}+z_{22}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_{22}a + z_{22}c\cos\beta\right)\hat{\mathbf{x}} - y_{22}b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_{22}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	S II

References:

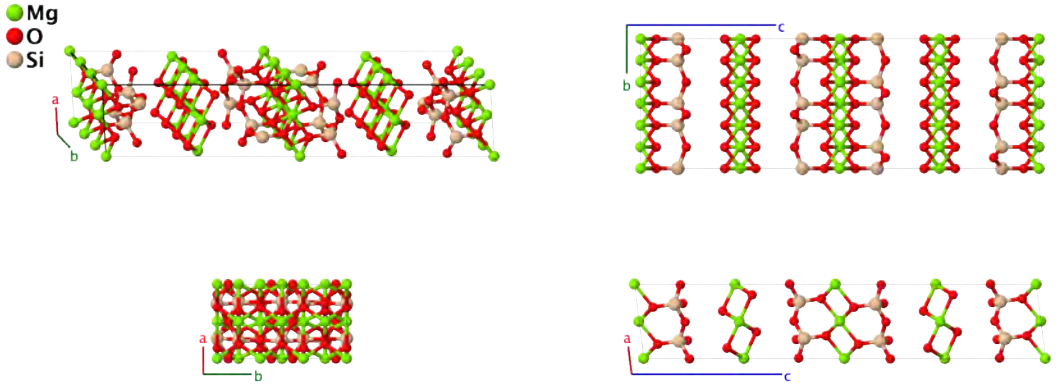
- R. Caminiti and G. Johansson, *A refinement of the Crystal Structure of the Cadmium Sulfate $3CdSO_4{\cdot}8H_2O$*, Acta Chem. Scand. **35a**, 451–455 (1981), doi:10.3891/acta.chem.scand.35a-0451.

Geometry files:

- CIF: pp. 1565
- POSCAR: pp. 1565

Chlorite [$Mg_3(Mg_2Al)(Si_3Al)O_{10}(OH)_8$, $S5_5$] Structure: A3B5C4D2_mC112_15_a3ef_5f_4f_2f

http://aflow.org/prototype-encyclopedia/A3B5C4D2_mC112_15_a3ef_5f_4f_2f

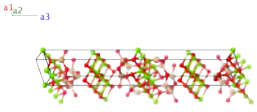


Prototype	:	$Mg_3O_5(OH)_4Si_2$
AFLOW prototype label	:	A3B5C4D2_mC112_15_a3ef_5f_4f_2f
***Strukturbericht* designation**	:	$S5_5$
Pearson symbol	:	mC112
Space group number	:	15
Space group symbol	:	$C2/c$
AFLOW prototype command	:	`aflow --proto=A3B5C4D2_mC112_15_a3ef_5f_4f_2f --params=`$a, b/a, c/a, \beta, y_2, y_3, y_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}$

- We follow the interpretation of (McMurchy, 1934) as given by (Downs, 2003): The sites we have labeled "Mg (4*e*)" are actually 2/3 magnesium and 1/3 aluminum, and the sites we have labeled "Si (8*f*)" are 3/4 silicon and 1/4 aluminum.

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(4*a*)	Mg I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	(4*a*)	Mg I
$\mathbf{B}_3$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Mg II
$\mathbf{B}_4$	=	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Mg II
$\mathbf{B}_5$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Mg III

$\mathbf{B}_6$	=	$y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Mg III
$\mathbf{B}_7$	=	$-y_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Mg IV
$\mathbf{B}_8$	=	$y_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Mg IV
$\mathbf{B}_9$	=	$(x_5 - y_5)\,\mathbf{a}_1 + (x_5 + y_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Mg V
$\mathbf{B}_{10}$	=	$(-x_5 - y_5)\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_5 a - z_5 c\cos\beta\right)\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Mg V
$\mathbf{B}_{11}$	=	$(-x_5 + y_5)\,\mathbf{a}_1 + (-x_5 - y_5)\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Mg V
$\mathbf{B}_{12}$	=	$(x_5 + y_5)\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Mg V
$\mathbf{B}_{13}$	=	$(x_6 - y_6)\,\mathbf{a}_1 + (x_6 + y_6)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O I
$\mathbf{B}_{14}$	=	$(-x_6 - y_6)\,\mathbf{a}_1 + (-x_6 + y_6)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_6 a - z_6 c\cos\beta\right)\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O I
$\mathbf{B}_{15}$	=	$(-x_6 + y_6)\,\mathbf{a}_1 + (-x_6 - y_6)\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O I
$\mathbf{B}_{16}$	=	$(x_6 + y_6)\,\mathbf{a}_1 + (x_6 - y_6)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O I
$\mathbf{B}_{17}$	=	$(x_7 - y_7)\,\mathbf{a}_1 + (x_7 + y_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{18}$	=	$(-x_7 - y_7)\,\mathbf{a}_1 + (-x_7 + y_7)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_7 a - z_7 c\cos\beta\right)\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{19}$	=	$(-x_7 + y_7)\,\mathbf{a}_1 + (-x_7 - y_7)\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{20}$	=	$(x_7 + y_7)\,\mathbf{a}_1 + (x_7 - y_7)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_7 a + z_7 c\cos\beta\right)\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{21}$	=	$(x_8 - y_8)\,\mathbf{a}_1 + (x_8 + y_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{22}$	=	$(-x_8 - y_8)\,\mathbf{a}_1 + (-x_8 + y_8)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_8 a - z_8 c\cos\beta\right)\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{23}$	=	$(-x_8 + y_8)\,\mathbf{a}_1 + (-x_8 - y_8)\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{24}$	=	$(x_8 + y_8)\,\mathbf{a}_1 + (x_8 - y_8)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_8 a + z_8 c\cos\beta\right)\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{25}$	=	$(x_9 - y_9)\,\mathbf{a}_1 + (x_9 + y_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$(x_9 a + z_9 c\cos\beta)\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O IV
$\mathbf{B}_{26}$	=	$(-x_9 - y_9)\,\mathbf{a}_1 + (-x_9 + y_9)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_9 a - z_9 c\cos\beta\right)\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_9\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O IV
$\mathbf{B}_{27}$	=	$(-x_9 + y_9)\,\mathbf{a}_1 + (-x_9 - y_9)\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$(-x_9 a - z_9 c\cos\beta)\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} - z_9 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O IV
$\mathbf{B}_{28}$	=	$(x_9 + y_9)\,\mathbf{a}_1 + (x_9 - y_9)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_9 a + z_9 c\cos\beta\right)\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O IV

$$\mathbf{B}_{29} = (x_{10} - y_{10})\,\mathbf{a}_1 + (x_{10} + y_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 = (x_{10}a + z_{10}c\cos\beta)\,\hat{\mathbf{x}} + y_{10}b\,\hat{\mathbf{y}} + z_{10}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O V}$$

$$\mathbf{B}_{30} = (-x_{10} - y_{10})\,\mathbf{a}_1 + (-x_{10} + y_{10})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{10}a - z_{10}c\cos\beta\right)\hat{\mathbf{x}} + y_{10}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{10}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O V}$$

$$\mathbf{B}_{31} = (-x_{10} + y_{10})\,\mathbf{a}_1 + (-x_{10} - y_{10})\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 = (-x_{10}a - z_{10}c\cos\beta)\,\hat{\mathbf{x}} - y_{10}b\,\hat{\mathbf{y}} - z_{10}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O V}$$

$$\mathbf{B}_{32} = (x_{10} + y_{10})\,\mathbf{a}_1 + (x_{10} - y_{10})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{10}a + z_{10}c\cos\beta\right)\hat{\mathbf{x}} - y_{10}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{10}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O V}$$

$$\mathbf{B}_{33} = (x_{11} - y_{11})\,\mathbf{a}_1 + (x_{11} + y_{11})\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 = (x_{11}a + z_{11}c\cos\beta)\,\hat{\mathbf{x}} + y_{11}b\,\hat{\mathbf{y}} + z_{11}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{OH I}$$

$$\mathbf{B}_{34} = (-x_{11} - y_{11})\,\mathbf{a}_1 + (-x_{11} + y_{11})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{11}a - z_{11}c\cos\beta\right)\hat{\mathbf{x}} + y_{11}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{11}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{OH I}$$

$$\mathbf{B}_{35} = (-x_{11} + y_{11})\,\mathbf{a}_1 + (-x_{11} - y_{11})\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3 = (-x_{11}a - z_{11}c\cos\beta)\,\hat{\mathbf{x}} - y_{11}b\,\hat{\mathbf{y}} - z_{11}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{OH I}$$

$$\mathbf{B}_{36} = (x_{11} + y_{11})\,\mathbf{a}_1 + (x_{11} - y_{11})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{11}a + z_{11}c\cos\beta\right)\hat{\mathbf{x}} - y_{11}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{11}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{OH I}$$

$$\mathbf{B}_{37} = (x_{12} - y_{12})\,\mathbf{a}_1 + (x_{12} + y_{12})\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 = (x_{12}a + z_{12}c\cos\beta)\,\hat{\mathbf{x}} + y_{12}b\,\hat{\mathbf{y}} + z_{12}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{OH II}$$

$$\mathbf{B}_{38} = (-x_{12} - y_{12})\,\mathbf{a}_1 + (-x_{12} + y_{12})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{12}a - z_{12}c\cos\beta\right)\hat{\mathbf{x}} + y_{12}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{12}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{OH II}$$

$$\mathbf{B}_{39} = (-x_{12} + y_{12})\,\mathbf{a}_1 + (-x_{12} - y_{12})\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3 = (-x_{12}a - z_{12}c\cos\beta)\,\hat{\mathbf{x}} - y_{12}b\,\hat{\mathbf{y}} - z_{12}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{OH II}$$

$$\mathbf{B}_{40} = (x_{12} + y_{12})\,\mathbf{a}_1 + (x_{12} - y_{12})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{12}a + z_{12}c\cos\beta\right)\hat{\mathbf{x}} - y_{12}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{12}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{OH II}$$

$$\mathbf{B}_{41} = (x_{13} - y_{13})\,\mathbf{a}_1 + (x_{13} + y_{13})\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3 = (x_{13}a + z_{13}c\cos\beta)\,\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} + z_{13}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{OH III}$$

$$\mathbf{B}_{42} = (-x_{13} - y_{13})\,\mathbf{a}_1 + (-x_{13} + y_{13})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{13}a - z_{13}c\cos\beta\right)\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{13}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{OH III}$$

$$\mathbf{B}_{43} = (-x_{13} + y_{13})\,\mathbf{a}_1 + (-x_{13} - y_{13})\,\mathbf{a}_2 - z_{13}\,\mathbf{a}_3 = (-x_{13}a - z_{13}c\cos\beta)\,\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} - z_{13}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{OH III}$$

$$\mathbf{B}_{44} = (x_{13} + y_{13})\,\mathbf{a}_1 + (x_{13} - y_{13})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{13}a + z_{13}c\cos\beta\right)\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{13}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{OH III}$$

$$\mathbf{B}_{45} = (x_{14} - y_{14})\,\mathbf{a}_1 + (x_{14} + y_{14})\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3 = (x_{14}a + z_{14}c\cos\beta)\,\hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} + z_{14}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{OH IV}$$

$$\mathbf{B}_{46} = (-x_{14} - y_{14})\,\mathbf{a}_1 + (-x_{14} + y_{14})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{14}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{14}a - z_{14}c\cos\beta\right)\hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{14}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{OH IV}$$

$$\mathbf{B}_{47} = (-x_{14} + y_{14})\,\mathbf{a}_1 + (-x_{14} - y_{14})\,\mathbf{a}_2 - z_{14}\,\mathbf{a}_3 = (-x_{14}a - z_{14}c\cos\beta)\,\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} - z_{14}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{OH IV}$$

$$\mathbf{B}_{48} = (x_{14} + y_{14})\,\mathbf{a}_1 + (x_{14} - y_{14})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{14}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{14}a + z_{14}c\cos\beta\right)\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{14}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{OH IV}$$

$$\mathbf{B}_{49} = (x_{15} - y_{15})\,\mathbf{a}_1 + (x_{15} + y_{15})\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 = (x_{15}a + z_{15}c\cos\beta)\,\hat{\mathbf{x}} + y_{15}b\,\hat{\mathbf{y}} + z_{15}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{Si I}$$

$$\mathbf{B}_{50} = (-x_{15} - y_{15})\,\mathbf{a}_1 + (-x_{15} + y_{15})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{15}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{15}a - z_{15}c\cos\beta\right)\hat{\mathbf{x}} + y_{15}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{15}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{Si I}$$

$\mathbf{B}_{51}$	=	$(-x_{15}+y_{15})\,\mathbf{a}_1+(-x_{15}-y_{15})\,\mathbf{a}_2-z_{15}\,\mathbf{a}_3$	=	$(-x_{15}a-z_{15}c\cos\beta)\,\hat{\mathbf{x}}-y_{15}b\,\hat{\mathbf{y}}-z_{15}c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Si I
$\mathbf{B}_{52}$	=	$(x_{15}+y_{15})\,\mathbf{a}_1+(x_{15}-y_{15})\,\mathbf{a}_2+\left(\frac{1}{2}+z_{15}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta+x_{15}a+z_{15}c\cos\beta\right)\hat{\mathbf{x}}-y_{15}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{15}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Si I
$\mathbf{B}_{53}$	=	$(x_{16}-y_{16})\,\mathbf{a}_1+(x_{16}+y_{16})\,\mathbf{a}_2+z_{16}\,\mathbf{a}_3$	=	$(x_{16}a+z_{16}c\cos\beta)\,\hat{\mathbf{x}}+y_{16}b\,\hat{\mathbf{y}}+z_{16}c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Si II
$\mathbf{B}_{54}$	=	$(-x_{16}-y_{16})\,\mathbf{a}_1+(-x_{16}+y_{16})\,\mathbf{a}_2+\left(\frac{1}{2}-z_{16}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta-x_{16}a-z_{16}c\cos\beta\right)\hat{\mathbf{x}}+y_{16}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{16}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Si II
$\mathbf{B}_{55}$	=	$(-x_{16}+y_{16})\,\mathbf{a}_1+(-x_{16}-y_{16})\,\mathbf{a}_2-z_{16}\,\mathbf{a}_3$	=	$(-x_{16}a-z_{16}c\cos\beta)\,\hat{\mathbf{x}}-y_{16}b\,\hat{\mathbf{y}}-z_{16}c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Si II
$\mathbf{B}_{56}$	=	$(x_{16}+y_{16})\,\mathbf{a}_1+(x_{16}-y_{16})\,\mathbf{a}_2+\left(\frac{1}{2}+z_{16}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta+x_{16}a+z_{16}c\cos\beta\right)\hat{\mathbf{x}}-y_{16}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{16}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Si II

References:

- R. C. McMurchy, *The Crystal Structure of the Chlorite Minerals*, Zeitschrift für Kristallographie - Crystalline Materials **88**, 420–432 (1934), doi:10.1524/zkri.1934.88.1.420.

Found in:

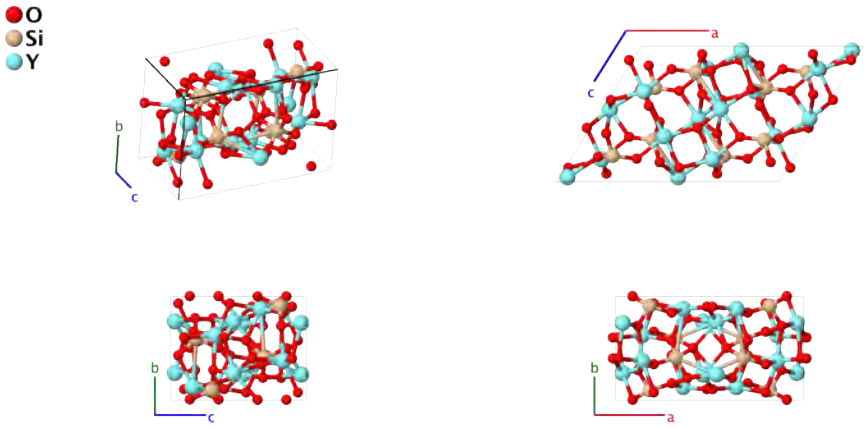
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1566
- POSCAR: pp. 1566

Y_2SiO_5 (RE_2SiO_5 X2) Structure: A5BC2_mC64_15_5f_f_2f

http://aflow.org/prototype-encyclopedia/A5BC2_mC64_15_5f_f_2f



Prototype	:	O_5SiY_2
AFLOW prototype label	:	A5BC2_mC64_15_5f_f_2f
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC64
Space group number	:	15
Space group symbol	:	$C2/c$
AFLOW prototype command	:	`aflow --proto=A5BC2_mC64_15_5f_f_2f --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8$

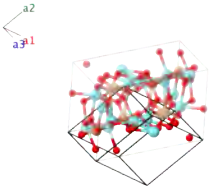
Other compounds with this structure

- Dy_2SiO_5, Er_2SiO_5, Ho_2SiO_5, Lu_2SiO_5, Tb_2SiO_5, Tm_2SiO_5, and Yb_2SiO_5

- Compounds of the form $RESiO_5$ (RE = Rare Earth and related elements) crystallize in one of two forms (Wang, 2001 and Tian, 2016): X1, space group $P2_1/c$ #14, for rare earths between lanthanum and gadolinium, and X2, this structure, for the later rare earths, with the dysprosium, yttrium and ytterbium compounds existing in both structures.
- This is the high temperature phase of Y_2SiO_5. Below 1190 °C it transforms into the X1 phase.

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$\mathbf{B}_1$	=	$(x_1 - y_1)\,\mathbf{a}_1 + (x_1 + y_1)\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$(x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O I
$\mathbf{B}_2$	=	$(-x_1 - y_1)\,\mathbf{a}_1 + (-x_1 + y_1)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_1 a - z_1 c\cos\beta\right)\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O I
$\mathbf{B}_3$	=	$(-x_1 + y_1)\,\mathbf{a}_1 + (-x_1 - y_1)\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$(-x_1 a - z_1 c\cos\beta)\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} - z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O I
$\mathbf{B}_4$	=	$(x_1 + y_1)\,\mathbf{a}_1 + (x_1 - y_1)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_1 a + z_1 c\cos\beta\right)\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O I
$\mathbf{B}_5$	=	$(x_2 - y_2)\,\mathbf{a}_1 + (x_2 + y_2)\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_6$	=	$(-x_2 - y_2)\,\mathbf{a}_1 + (-x_2 + y_2)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_2 a - z_2 c\cos\beta\right)\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_7$	=	$(-x_2 + y_2)\,\mathbf{a}_1 + (-x_2 - y_2)\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} - z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_8$	=	$(x_2 + y_2)\,\mathbf{a}_1 + (x_2 - y_2)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_2 a + z_2 c\cos\beta\right)\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_9$	=	$(x_3 - y_3)\,\mathbf{a}_1 + (x_3 + y_3)\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{10}$	=	$(-x_3 - y_3)\,\mathbf{a}_1 + (-x_3 + y_3)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_3 a - z_3 c\cos\beta\right)\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{11}$	=	$(-x_3 + y_3)\,\mathbf{a}_1 + (-x_3 - y_3)\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{12}$	=	$(x_3 + y_3)\,\mathbf{a}_1 + (x_3 - y_3)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{13}$	=	$(x_4 - y_4)\,\mathbf{a}_1 + (x_4 + y_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O IV
$\mathbf{B}_{14}$	=	$(-x_4 - y_4)\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_4 a - z_4 c\cos\beta\right)\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O IV
$\mathbf{B}_{15}$	=	$(-x_4 + y_4)\,\mathbf{a}_1 + (-x_4 - y_4)\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O IV
$\mathbf{B}_{16}$	=	$(x_4 + y_4)\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O IV
$\mathbf{B}_{17}$	=	$(x_5 - y_5)\,\mathbf{a}_1 + (x_5 + y_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O V
$\mathbf{B}_{18}$	=	$(-x_5 - y_5)\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_5 a - z_5 c\cos\beta\right)\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O V
$\mathbf{B}_{19}$	=	$(-x_5 + y_5)\,\mathbf{a}_1 + (-x_5 - y_5)\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O V
$\mathbf{B}_{20}$	=	$(x_5 + y_5)\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O V
$\mathbf{B}_{21}$	=	$(x_6 - y_6)\,\mathbf{a}_1 + (x_6 + y_6)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Si
$\mathbf{B}_{22}$	=	$(-x_6 - y_6)\,\mathbf{a}_1 + (-x_6 + y_6)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_6 a - z_6 c\cos\beta\right)\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Si

$\mathbf{B}_{23}$	=	$(-x_6 + y_6)\,\mathbf{a}_1 + (-x_6 - y_6)\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Si
$\mathbf{B}_{24}$	=	$(x_6 + y_6)\,\mathbf{a}_1 + (x_6 - y_6)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Si
$\mathbf{B}_{25}$	=	$(x_7 - y_7)\,\mathbf{a}_1 + (x_7 + y_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Y I
$\mathbf{B}_{26}$	=	$(-x_7 - y_7)\,\mathbf{a}_1 + (-x_7 + y_7)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_7 a - z_7 c\cos\beta\right)\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Y I
$\mathbf{B}_{27}$	=	$(-x_7 + y_7)\,\mathbf{a}_1 + (-x_7 - y_7)\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Y I
$\mathbf{B}_{28}$	=	$(x_7 + y_7)\,\mathbf{a}_1 + (x_7 - y_7)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_7 a + z_7 c\cos\beta\right)\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Y I
$\mathbf{B}_{29}$	=	$(x_8 - y_8)\,\mathbf{a}_1 + (x_8 + y_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Y II
$\mathbf{B}_{30}$	=	$(-x_8 - y_8)\,\mathbf{a}_1 + (-x_8 + y_8)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_8 a - z_8 c\cos\beta\right)\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Y II
$\mathbf{B}_{31}$	=	$(-x_8 + y_8)\,\mathbf{a}_1 + (-x_8 - y_8)\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Y II
$\mathbf{B}_{32}$	=	$(x_8 + y_8)\,\mathbf{a}_1 + (x_8 - y_8)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_8 a + z_8 c\cos\beta\right)\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Y II

References:

- G. V. Anan'eva, A. M. Korovkin, T. I. Merkulyaeva, A. M. Morozova, M. V. Petrov, I. R. Savinova, V. R. Startsev, and P. P. Feofilov, *Growth of lanthanide oxyorthosilicate single crystals, and their structural and optical characteristics*, Inorg. Mat. **17**, 754–758 (1981). Translated from *Neorganicheskie Materialy*.
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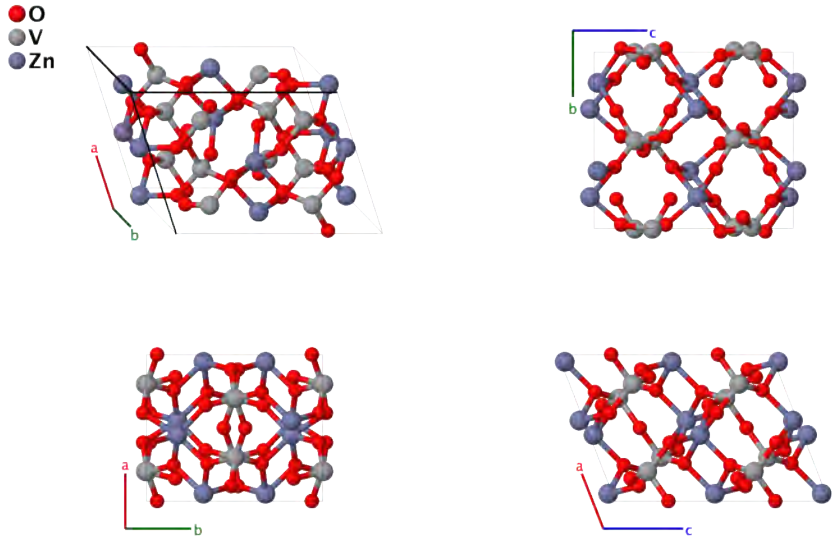
Geometry files:

- CIF: pp. 1567
- POSCAR: pp. 1567

α-$Zn_2V_2O_7$ Structure: A7B2C2_mC44_15_e3f_f_f

http://aflow.org/prototype-encyclopedia/A7B2C2_mC44_15_e3f_f_f



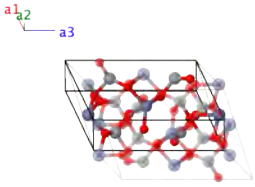
Prototype	:	$O_7V_2Zn_2$
AFLOW prototype label	:	A7B2C2_mC44_15_e3f_f_f
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC44
Space group number	:	15
Space group symbol	:	$C2/c$
AFLOW prototype command	:	`aflow --proto=A7B2C2_mC44_15_e3f_f_f --params=`$a, b/a, c/a, \beta, y_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6$

Other compounds with this structure

- β-$Cu_2V_2O_7$

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_2$	=	$y_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_3$	=	$(x_2 - y_2)\,\mathbf{a}_1 + (x_2 + y_2)\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II

$\mathbf{B}_4$	=	$(-x_2 - y_2)\,\mathbf{a}_1 + (-x_2 + y_2)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_2 a - z_2 c\cos\beta\right)\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_5$	=	$(-x_2 + y_2)\,\mathbf{a}_1 + (-x_2 - y_2)\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$(-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} - z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_6$	=	$(x_2 + y_2)\,\mathbf{a}_1 + (x_2 - y_2)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_2 a + z_2 c\cos\beta\right)\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_7$	=	$(x_3 - y_3)\,\mathbf{a}_1 + (x_3 + y_3)\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_8$	=	$(-x_3 - y_3)\,\mathbf{a}_1 + (-x_3 + y_3)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_3 a - z_3 c\cos\beta\right)\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_9$	=	$(-x_3 + y_3)\,\mathbf{a}_1 + (-x_3 - y_3)\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{10}$	=	$(x_3 + y_3)\,\mathbf{a}_1 + (x_3 - y_3)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{11}$	=	$(x_4 - y_4)\,\mathbf{a}_1 + (x_4 + y_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O IV
$\mathbf{B}_{12}$	=	$(-x_4 - y_4)\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_4 a - z_4 c\cos\beta\right)\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O IV
$\mathbf{B}_{13}$	=	$(-x_4 + y_4)\,\mathbf{a}_1 + (-x_4 - y_4)\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O IV
$\mathbf{B}_{14}$	=	$(x_4 + y_4)\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O IV
$\mathbf{B}_{15}$	=	$(x_5 - y_5)\,\mathbf{a}_1 + (x_5 + y_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	V
$\mathbf{B}_{16}$	=	$(-x_5 - y_5)\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_5 a - z_5 c\cos\beta\right)\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	V
$\mathbf{B}_{17}$	=	$(-x_5 + y_5)\,\mathbf{a}_1 + (-x_5 - y_5)\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	V
$\mathbf{B}_{18}$	=	$(x_5 + y_5)\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	V
$\mathbf{B}_{19}$	=	$(x_6 - y_6)\,\mathbf{a}_1 + (x_6 + y_6)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Zn
$\mathbf{B}_{20}$	=	$(-x_6 - y_6)\,\mathbf{a}_1 + (-x_6 + y_6)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_6 a - z_6 c\cos\beta\right)\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Zn
$\mathbf{B}_{21}$	=	$(-x_6 + y_6)\,\mathbf{a}_1 + (-x_6 - y_6)\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Zn
$\mathbf{B}_{22}$	=	$(x_6 + y_6)\,\mathbf{a}_1 + (x_6 - y_6)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Zn

References:

- R. Gopal and C. Calvo, *Crystal Structure of α-$Zn_2V_2O_7$*, Can. J. Chem. **51**, 1004–1009 (1973), doi:10.1139/v73-149.

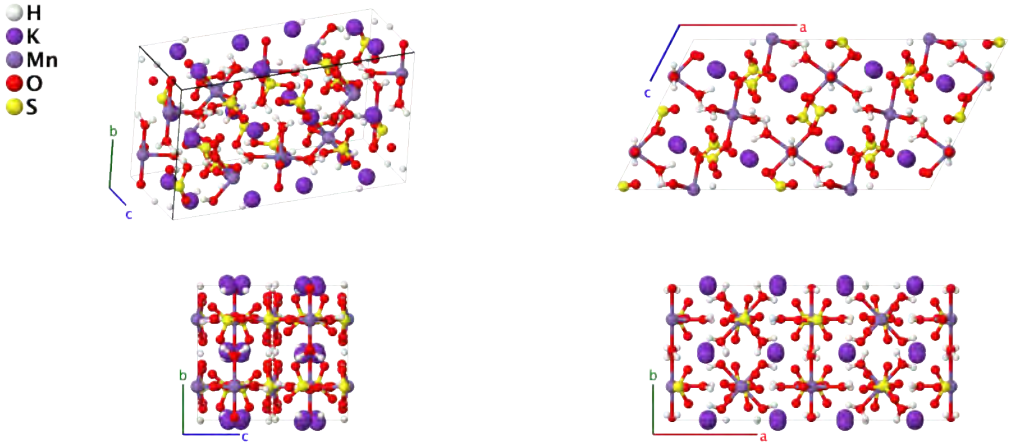
Found in:

- C. Calvo and R. Faggiani, *α Cupric Divanadate*, Acta Crystallogr. Sect. B Struct. Sci. **31**, 603–605 (1975), doi:10.1107/S0567740875003354.

Geometry files:
- CIF: pp. 1567
- POSCAR: pp. 1568

Manganese-leonite 185 K [$K_2Mn(SO_4)_2{\cdot}4H_2O$] Structure: A8B2CD12E2_mC200_15_8f_2f_ce_2e11f_2f

http://aflow.org/prototype-encyclopedia/A8B2CD12E2_mC200_15_8f_2f_ce_2e11f_2f



Prototype	:	$H_8K_2MnO_{12}S_2$
AFLOW prototype label	:	A8B2CD12E2_mC200_15_8f_2f_ce_2e11f_2f
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC200
Space group number	:	15
Space group symbol	:	$C2/c$
AFLOW prototype command	:	`aflow --proto=A8B2CD12E2_mC200_15_8f_2f_ce_2e11f_2f --params=`$a, b/a, c/a, \beta, y_2, y_3, y_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}, x_{17}, y_{17}, z_{17}, x_{18}, y_{18}, z_{18}, x_{19}, y_{19}, z_{19}, x_{20}, y_{20}, z_{20}, x_{21}, y_{21}, z_{21}, x_{22}, y_{22}, z_{22}, x_{23}, y_{23}, z_{23}, x_{24}, y_{24}, z_{24}, x_{25}, y_{25}, z_{25}, x_{26}, y_{26}, z_{26}, x_{27}, y_{27}, z_{27}$

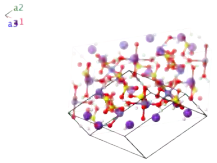
Other compounds with this structure

- $K_2Mg(SO_4)_2{\cdot}4H_2O$ (leonite) and $K_2Fe(SO_4)_2{\cdot}4H_2O$ (mereiterite)

- This is the intermediate-low temperature structure of leonite. For Mn-leonite, discussed here, it is stable between 168 K and 205 K, and we show the structure at 185 K.
- The room temperature structure is *Strukturbericht* $H4_{23}$, space group $C2/m$ #13. The current structure doubles the size of the unit cell and orders all of the SO_4 radicals.
- The low temperature structure has space group $P2_1/c$ #14.
- (Hertweck, 2001) give crystallographic information of the 185 K phase in the $I2/a$ setting of space group #15, with the origin supposedly shifted by $(1/4, 1/4, 1/2)$ from the -1 point on the a-glide plane (the symmetry operations for this setting may be found here). We were unable to use their data to construct a realistic crystal structure. Instead, we used the interpretation of their results by (Villars, 2016) to put the structure in the standard $C2/c$ setting of space group #15.

Base-centered Monoclinic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2}\,b\,\hat{\mathbf{y}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,b\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}}$	$(4c)$	Mn I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{4}a + \frac{1}{2}c\cos\beta\right)\,\hat{\mathbf{x}} - \frac{1}{4}b\,\hat{\mathbf{y}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	$(4c)$	Mn I
$\mathbf{B}_3$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Mn II
$\mathbf{B}_4$	=	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Mn II
$\mathbf{B}_5$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_6$	=	$y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_7$	=	$-y_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_8$	=	$y_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_9$	=	$(x_5 - y_5)\,\mathbf{a}_1 + (x_5 + y_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H I
$\mathbf{B}_{10}$	=	$(-x_5 - y_5)\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_5 a - z_5 c\cos\beta\right)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H I
$\mathbf{B}_{11}$	=	$(-x_5 + y_5)\,\mathbf{a}_1 + (-x_5 - y_5)\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H I
$\mathbf{B}_{12}$	=	$(x_5 + y_5)\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H I
$\mathbf{B}_{13}$	=	$(x_6 - y_6)\,\mathbf{a}_1 + (x_6 + y_6)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H II
$\mathbf{B}_{14}$	=	$(-x_6 - y_6)\,\mathbf{a}_1 + (-x_6 + y_6)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_6 a - z_6 c\cos\beta\right)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H II
$\mathbf{B}_{15}$	=	$(-x_6 + y_6)\,\mathbf{a}_1 + (-x_6 - y_6)\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H II
$\mathbf{B}_{16}$	=	$(x_6 + y_6)\,\mathbf{a}_1 + (x_6 - y_6)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H II
$\mathbf{B}_{17}$	=	$(x_7 - y_7)\,\mathbf{a}_1 + (x_7 + y_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H III
$\mathbf{B}_{18}$	=	$(-x_7 - y_7)\,\mathbf{a}_1 + (-x_7 + y_7)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_7 a - z_7 c\cos\beta\right)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H III
$\mathbf{B}_{19}$	=	$(-x_7 + y_7)\,\mathbf{a}_1 + (-x_7 - y_7)\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H III
$\mathbf{B}_{20}$	=	$(x_7 + y_7)\,\mathbf{a}_1 + (x_7 - y_7)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_7 a + z_7 c\cos\beta\right)\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H III
$\mathbf{B}_{21}$	=	$(x_8 - y_8)\,\mathbf{a}_1 + (x_8 + y_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H IV

$$\mathbf{B}_{22} = (-x_8 - y_8)\,\mathbf{a}_1 + (-x_8 + y_8)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_8 a - z_8 c\cos\beta\right)\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_8\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H IV}$$

$$\mathbf{B}_{23} = (-x_8 + y_8)\,\mathbf{a}_1 + (-x_8 - y_8)\,\mathbf{a}_2 - z_8\,\mathbf{a}_3 = (-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H IV}$$

$$\mathbf{B}_{24} = (x_8 + y_8)\,\mathbf{a}_1 + (x_8 - y_8)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_8 a + z_8 c\cos\beta\right)\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_8\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H IV}$$

$$\mathbf{B}_{25} = (x_9 - y_9)\,\mathbf{a}_1 + (x_9 + y_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 = (x_9 a + z_9 c\cos\beta)\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H V}$$

$$\mathbf{B}_{26} = (-x_9 - y_9)\,\mathbf{a}_1 + (-x_9 + y_9)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_9 a - z_9 c\cos\beta\right)\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_9\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H V}$$

$$\mathbf{B}_{27} = (-x_9 + y_9)\,\mathbf{a}_1 + (-x_9 - y_9)\,\mathbf{a}_2 - z_9\,\mathbf{a}_3 = (-x_9 a - z_9 c\cos\beta)\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} - z_9 c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H V}$$

$$\mathbf{B}_{28} = (x_9 + y_9)\,\mathbf{a}_1 + (x_9 - y_9)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_9 a + z_9 c\cos\beta\right)\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_9\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H V}$$

$$\mathbf{B}_{29} = (x_{10} - y_{10})\,\mathbf{a}_1 + (x_{10} + y_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 = (x_{10} a + z_{10} c\cos\beta)\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H VI}$$

$$\mathbf{B}_{30} = (-x_{10} - y_{10})\,\mathbf{a}_1 + (-x_{10} + y_{10})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{10} a - z_{10} c\cos\beta\right)\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{10}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H VI}$$

$$\mathbf{B}_{31} = (-x_{10} + y_{10})\,\mathbf{a}_1 + (-x_{10} - y_{10})\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 = (-x_{10} a - z_{10} c\cos\beta)\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} - z_{10} c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H VI}$$

$$\mathbf{B}_{32} = (x_{10} + y_{10})\,\mathbf{a}_1 + (x_{10} - y_{10})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{10} a + z_{10} c\cos\beta\right)\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{10}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H VI}$$

$$\mathbf{B}_{33} = (x_{11} - y_{11})\,\mathbf{a}_1 + (x_{11} + y_{11})\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 = (x_{11} a + z_{11} c\cos\beta)\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + z_{11} c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H VII}$$

$$\mathbf{B}_{34} = (-x_{11} - y_{11})\,\mathbf{a}_1 + (-x_{11} + y_{11})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{11} a - z_{11} c\cos\beta\right)\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{11}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H VII}$$

$$\mathbf{B}_{35} = (-x_{11} + y_{11})\,\mathbf{a}_1 + (-x_{11} - y_{11})\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3 = (-x_{11} a - z_{11} c\cos\beta)\,\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} - z_{11} c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H VII}$$

$$\mathbf{B}_{36} = (x_{11} + y_{11})\,\mathbf{a}_1 + (x_{11} - y_{11})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{11} a + z_{11} c\cos\beta\right)\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{11}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H VII}$$

$$\mathbf{B}_{37} = (x_{12} - y_{12})\,\mathbf{a}_1 + (x_{12} + y_{12})\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 = (x_{12} a + z_{12} c\cos\beta)\,\hat{\mathbf{x}} + y_{12} b\,\hat{\mathbf{y}} + z_{12} c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H VIII}$$

$$\mathbf{B}_{38} = (-x_{12} - y_{12})\,\mathbf{a}_1 + (-x_{12} + y_{12})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{12} a - z_{12} c\cos\beta\right)\hat{\mathbf{x}} + y_{12} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{12}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H VIII}$$

$$\mathbf{B}_{39} = (-x_{12} + y_{12})\,\mathbf{a}_1 + (-x_{12} - y_{12})\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3 = (-x_{12} a - z_{12} c\cos\beta)\,\hat{\mathbf{x}} - y_{12} b\,\hat{\mathbf{y}} - z_{12} c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H VIII}$$

$$\mathbf{B}_{40} = (x_{12} + y_{12})\,\mathbf{a}_1 + (x_{12} - y_{12})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{12} a + z_{12} c\cos\beta\right)\hat{\mathbf{x}} - y_{12} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{12}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{H VIII}$$

$$\mathbf{B}_{41} = (x_{13} - y_{13})\,\mathbf{a}_1 + (x_{13} + y_{13})\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3 = (x_{13} a + z_{13} c\cos\beta)\,\hat{\mathbf{x}} + y_{13} b\,\hat{\mathbf{y}} + z_{13} c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{K I}$$

$$\mathbf{B}_{42} = (-x_{13} - y_{13})\,\mathbf{a}_1 + (-x_{13} + y_{13})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{13} a - z_{13} c\cos\beta\right)\hat{\mathbf{x}} + y_{13} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{13}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{K I}$$

$$\mathbf{B}_{43} = (-x_{13} + y_{13})\,\mathbf{a}_1 + (-x_{13} - y_{13})\,\mathbf{a}_2 - z_{13}\,\mathbf{a}_3 = (-x_{13} a - z_{13} c\cos\beta)\,\hat{\mathbf{x}} - y_{13} b\,\hat{\mathbf{y}} - z_{13} c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{K I}$$

$\mathbf{B}_{44}$	=	$(x_{13}+y_{13})\,\mathbf{a}_1+(x_{13}-y_{13})\,\mathbf{a}_2+\left(\frac{1}{2}+z_{13}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta+x_{13}a+z_{13}c\cos\beta\right)\hat{\mathbf{x}}-y_{13}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{13}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	K I
$\mathbf{B}_{45}$	=	$(x_{14}-y_{14})\,\mathbf{a}_1+(x_{14}+y_{14})\,\mathbf{a}_2+z_{14}\,\mathbf{a}_3$	=	$(x_{14}a+z_{14}c\cos\beta)\,\hat{\mathbf{x}}+y_{14}b\,\hat{\mathbf{y}}+z_{14}c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	K II
$\mathbf{B}_{46}$	=	$(-x_{14}-y_{14})\,\mathbf{a}_1+(-x_{14}+y_{14})\,\mathbf{a}_2+\left(\frac{1}{2}-z_{14}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta-x_{14}a-z_{14}c\cos\beta\right)\hat{\mathbf{x}}+y_{14}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{14}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	K II
$\mathbf{B}_{47}$	=	$(-x_{14}+y_{14})\,\mathbf{a}_1+(-x_{14}-y_{14})\,\mathbf{a}_2-z_{14}\,\mathbf{a}_3$	=	$(-x_{14}a-z_{14}c\cos\beta)\,\hat{\mathbf{x}}-y_{14}b\,\hat{\mathbf{y}}-z_{14}c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	K II
$\mathbf{B}_{48}$	=	$(x_{14}+y_{14})\,\mathbf{a}_1+(x_{14}-y_{14})\,\mathbf{a}_2+\left(\frac{1}{2}+z_{14}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta+x_{14}a+z_{14}c\cos\beta\right)\hat{\mathbf{x}}-y_{14}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{14}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	K II
$\mathbf{B}_{49}$	=	$(x_{15}-y_{15})\,\mathbf{a}_1+(x_{15}+y_{15})\,\mathbf{a}_2+z_{15}\,\mathbf{a}_3$	=	$(x_{15}a+z_{15}c\cos\beta)\,\hat{\mathbf{x}}+y_{15}b\,\hat{\mathbf{y}}+z_{15}c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{50}$	=	$(-x_{15}-y_{15})\,\mathbf{a}_1+(-x_{15}+y_{15})\,\mathbf{a}_2+\left(\frac{1}{2}-z_{15}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta-x_{15}a-z_{15}c\cos\beta\right)\hat{\mathbf{x}}+y_{15}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{15}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{51}$	=	$(-x_{15}+y_{15})\,\mathbf{a}_1+(-x_{15}-y_{15})\,\mathbf{a}_2-z_{15}\,\mathbf{a}_3$	=	$(-x_{15}a-z_{15}c\cos\beta)\,\hat{\mathbf{x}}-y_{15}b\,\hat{\mathbf{y}}-z_{15}c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{52}$	=	$(x_{15}+y_{15})\,\mathbf{a}_1+(x_{15}-y_{15})\,\mathbf{a}_2+\left(\frac{1}{2}+z_{15}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta+x_{15}a+z_{15}c\cos\beta\right)\hat{\mathbf{x}}-y_{15}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{15}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{53}$	=	$(x_{16}-y_{16})\,\mathbf{a}_1+(x_{16}+y_{16})\,\mathbf{a}_2+z_{16}\,\mathbf{a}_3$	=	$(x_{16}a+z_{16}c\cos\beta)\,\hat{\mathbf{x}}+y_{16}b\,\hat{\mathbf{y}}+z_{16}c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O IV
$\mathbf{B}_{54}$	=	$(-x_{16}-y_{16})\,\mathbf{a}_1+(-x_{16}+y_{16})\,\mathbf{a}_2+\left(\frac{1}{2}-z_{16}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta-x_{16}a-z_{16}c\cos\beta\right)\hat{\mathbf{x}}+y_{16}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{16}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O IV
$\mathbf{B}_{55}$	=	$(-x_{16}+y_{16})\,\mathbf{a}_1+(-x_{16}-y_{16})\,\mathbf{a}_2-z_{16}\,\mathbf{a}_3$	=	$(-x_{16}a-z_{16}c\cos\beta)\,\hat{\mathbf{x}}-y_{16}b\,\hat{\mathbf{y}}-z_{16}c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O IV
$\mathbf{B}_{56}$	=	$(x_{16}+y_{16})\,\mathbf{a}_1+(x_{16}-y_{16})\,\mathbf{a}_2+\left(\frac{1}{2}+z_{16}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta+x_{16}a+z_{16}c\cos\beta\right)\hat{\mathbf{x}}-y_{16}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{16}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O IV
$\mathbf{B}_{57}$	=	$(x_{17}-y_{17})\,\mathbf{a}_1+(x_{17}+y_{17})\,\mathbf{a}_2+z_{17}\,\mathbf{a}_3$	=	$(x_{17}a+z_{17}c\cos\beta)\,\hat{\mathbf{x}}+y_{17}b\,\hat{\mathbf{y}}+z_{17}c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O V
$\mathbf{B}_{58}$	=	$(-x_{17}-y_{17})\,\mathbf{a}_1+(-x_{17}+y_{17})\,\mathbf{a}_2+\left(\frac{1}{2}-z_{17}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta-x_{17}a-z_{17}c\cos\beta\right)\hat{\mathbf{x}}+y_{17}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{17}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O V
$\mathbf{B}_{59}$	=	$(-x_{17}+y_{17})\,\mathbf{a}_1+(-x_{17}-y_{17})\,\mathbf{a}_2-z_{17}\,\mathbf{a}_3$	=	$(-x_{17}a-z_{17}c\cos\beta)\,\hat{\mathbf{x}}-y_{17}b\,\hat{\mathbf{y}}-z_{17}c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O V
$\mathbf{B}_{60}$	=	$(x_{17}+y_{17})\,\mathbf{a}_1+(x_{17}-y_{17})\,\mathbf{a}_2+\left(\frac{1}{2}+z_{17}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta+x_{17}a+z_{17}c\cos\beta\right)\hat{\mathbf{x}}-y_{17}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{17}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O V
$\mathbf{B}_{61}$	=	$(x_{18}-y_{18})\,\mathbf{a}_1+(x_{18}+y_{18})\,\mathbf{a}_2+z_{18}\,\mathbf{a}_3$	=	$(x_{18}a+z_{18}c\cos\beta)\,\hat{\mathbf{x}}+y_{18}b\,\hat{\mathbf{y}}+z_{18}c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O VI
$\mathbf{B}_{62}$	=	$(-x_{18}-y_{18})\,\mathbf{a}_1+(-x_{18}+y_{18})\,\mathbf{a}_2+\left(\frac{1}{2}-z_{18}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta-x_{18}a-z_{18}c\cos\beta\right)\hat{\mathbf{x}}+y_{18}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{18}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O VI
$\mathbf{B}_{63}$	=	$(-x_{18}+y_{18})\,\mathbf{a}_1+(-x_{18}-y_{18})\,\mathbf{a}_2-z_{18}\,\mathbf{a}_3$	=	$(-x_{18}a-z_{18}c\cos\beta)\,\hat{\mathbf{x}}-y_{18}b\,\hat{\mathbf{y}}-z_{18}c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O VI
$\mathbf{B}_{64}$	=	$(x_{18}+y_{18})\,\mathbf{a}_1+(x_{18}-y_{18})\,\mathbf{a}_2+\left(\frac{1}{2}+z_{18}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta+x_{18}a+z_{18}c\cos\beta\right)\hat{\mathbf{x}}-y_{18}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{18}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O VI
$\mathbf{B}_{65}$	=	$(x_{19}-y_{19})\,\mathbf{a}_1+(x_{19}+y_{19})\,\mathbf{a}_2+z_{19}\,\mathbf{a}_3$	=	$(x_{19}a+z_{19}c\cos\beta)\,\hat{\mathbf{x}}+y_{19}b\,\hat{\mathbf{y}}+z_{19}c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O VII

$$\mathbf{B}_{66} = (-x_{19} - y_{19})\,\mathbf{a}_1 + (-x_{19} + y_{19})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{19}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{19}a - z_{19}c\cos\beta\right)\hat{\mathbf{x}} + y_{19}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{19}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VII}$$

$$\mathbf{B}_{67} = (-x_{19} + y_{19})\,\mathbf{a}_1 + (-x_{19} - y_{19})\,\mathbf{a}_2 - z_{19}\,\mathbf{a}_3 = (-x_{19}a - z_{19}c\cos\beta)\,\hat{\mathbf{x}} - y_{19}b\,\hat{\mathbf{y}} - z_{19}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VII}$$

$$\mathbf{B}_{68} = (x_{19} + y_{19})\,\mathbf{a}_1 + (x_{19} - y_{19})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{19}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{19}a + z_{19}c\cos\beta\right)\hat{\mathbf{x}} - y_{19}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{19}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VII}$$

$$\mathbf{B}_{69} = (x_{20} - y_{20})\,\mathbf{a}_1 + (x_{20} + y_{20})\,\mathbf{a}_2 + z_{20}\,\mathbf{a}_3 = (x_{20}a + z_{20}c\cos\beta)\,\hat{\mathbf{x}} + y_{20}b\,\hat{\mathbf{y}} + z_{20}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VIII}$$

$$\mathbf{B}_{70} = (-x_{20} - y_{20})\,\mathbf{a}_1 + (-x_{20} + y_{20})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{20}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{20}a - z_{20}c\cos\beta\right)\hat{\mathbf{x}} + y_{20}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{20}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VIII}$$

$$\mathbf{B}_{71} = (-x_{20} + y_{20})\,\mathbf{a}_1 + (-x_{20} - y_{20})\,\mathbf{a}_2 - z_{20}\,\mathbf{a}_3 = (-x_{20}a - z_{20}c\cos\beta)\,\hat{\mathbf{x}} - y_{20}b\,\hat{\mathbf{y}} - z_{20}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VIII}$$

$$\mathbf{B}_{72} = (x_{20} + y_{20})\,\mathbf{a}_1 + (x_{20} - y_{20})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{20}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{20}a + z_{20}c\cos\beta\right)\hat{\mathbf{x}} - y_{20}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{20}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O VIII}$$

$$\mathbf{B}_{73} = (x_{21} - y_{21})\,\mathbf{a}_1 + (x_{21} + y_{21})\,\mathbf{a}_2 + z_{21}\,\mathbf{a}_3 = (x_{21}a + z_{21}c\cos\beta)\,\hat{\mathbf{x}} + y_{21}b\,\hat{\mathbf{y}} + z_{21}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O IX}$$

$$\mathbf{B}_{74} = (-x_{21} - y_{21})\,\mathbf{a}_1 + (-x_{21} + y_{21})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{21}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{21}a - z_{21}c\cos\beta\right)\hat{\mathbf{x}} + y_{21}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{21}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O IX}$$

$$\mathbf{B}_{75} = (-x_{21} + y_{21})\,\mathbf{a}_1 + (-x_{21} - y_{21})\,\mathbf{a}_2 - z_{21}\,\mathbf{a}_3 = (-x_{21}a - z_{21}c\cos\beta)\,\hat{\mathbf{x}} - y_{21}b\,\hat{\mathbf{y}} - z_{21}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O IX}$$

$$\mathbf{B}_{76} = (x_{21} + y_{21})\,\mathbf{a}_1 + (x_{21} - y_{21})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{21}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{21}a + z_{21}c\cos\beta\right)\hat{\mathbf{x}} - y_{21}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{21}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O IX}$$

$$\mathbf{B}_{77} = (x_{22} - y_{22})\,\mathbf{a}_1 + (x_{22} + y_{22})\,\mathbf{a}_2 + z_{22}\,\mathbf{a}_3 = (x_{22}a + z_{22}c\cos\beta)\,\hat{\mathbf{x}} + y_{22}b\,\hat{\mathbf{y}} + z_{22}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O X}$$

$$\mathbf{B}_{78} = (-x_{22} - y_{22})\,\mathbf{a}_1 + (-x_{22} + y_{22})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{22}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{22}a - z_{22}c\cos\beta\right)\hat{\mathbf{x}} + y_{22}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{22}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O X}$$

$$\mathbf{B}_{79} = (-x_{22} + y_{22})\,\mathbf{a}_1 + (-x_{22} - y_{22})\,\mathbf{a}_2 - z_{22}\,\mathbf{a}_3 = (-x_{22}a - z_{22}c\cos\beta)\,\hat{\mathbf{x}} - y_{22}b\,\hat{\mathbf{y}} - z_{22}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O X}$$

$$\mathbf{B}_{80} = (x_{22} + y_{22})\,\mathbf{a}_1 + (x_{22} - y_{22})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{22}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{22}a + z_{22}c\cos\beta\right)\hat{\mathbf{x}} - y_{22}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{22}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O X}$$

$$\mathbf{B}_{81} = (x_{23} - y_{23})\,\mathbf{a}_1 + (x_{23} + y_{23})\,\mathbf{a}_2 + z_{23}\,\mathbf{a}_3 = (x_{23}a + z_{23}c\cos\beta)\,\hat{\mathbf{x}} + y_{23}b\,\hat{\mathbf{y}} + z_{23}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O XI}$$

$$\mathbf{B}_{82} = (-x_{23} - y_{23})\,\mathbf{a}_1 + (-x_{23} + y_{23})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{23}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{23}a - z_{23}c\cos\beta\right)\hat{\mathbf{x}} + y_{23}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{23}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O XI}$$

$$\mathbf{B}_{83} = (-x_{23} + y_{23})\,\mathbf{a}_1 + (-x_{23} - y_{23})\,\mathbf{a}_2 - z_{23}\,\mathbf{a}_3 = (-x_{23}a - z_{23}c\cos\beta)\,\hat{\mathbf{x}} - y_{23}b\,\hat{\mathbf{y}} - z_{23}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O XI}$$

$$\mathbf{B}_{84} = (x_{23} + y_{23})\,\mathbf{a}_1 + (x_{23} - y_{23})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{23}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta + x_{23}a + z_{23}c\cos\beta\right)\hat{\mathbf{x}} - y_{23}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{23}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O XI}$$

$$\mathbf{B}_{85} = (x_{24} - y_{24})\,\mathbf{a}_1 + (x_{24} + y_{24})\,\mathbf{a}_2 + z_{24}\,\mathbf{a}_3 = (x_{24}a + z_{24}c\cos\beta)\,\hat{\mathbf{x}} + y_{24}b\,\hat{\mathbf{y}} + z_{24}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O XII}$$

$$\mathbf{B}_{86} = (-x_{24} - y_{24})\,\mathbf{a}_1 + (-x_{24} + y_{24})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{24}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}c\cos\beta - x_{24}a - z_{24}c\cos\beta\right)\hat{\mathbf{x}} + y_{24}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{24}\right)c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O XII}$$

$$\mathbf{B}_{87} = (-x_{24} + y_{24})\,\mathbf{a}_1 + (-x_{24} - y_{24})\,\mathbf{a}_2 - z_{24}\,\mathbf{a}_3 = (-x_{24}a - z_{24}c\cos\beta)\,\hat{\mathbf{x}} - y_{24}b\,\hat{\mathbf{y}} - z_{24}c\sin\beta\,\hat{\mathbf{z}} \quad (8f) \quad \text{O XII}$$

$\mathbf{B}_{88}$	=	$(x_{24}+y_{24})\,\mathbf{a}_1+(x_{24}-y_{24})\,\mathbf{a}_2+\left(\frac{1}{2}+z_{24}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta+x_{24}a+z_{24}c\cos\beta\right)\hat{\mathbf{x}}-y_{24}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{24}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O XII
$\mathbf{B}_{89}$	=	$(x_{25}-y_{25})\,\mathbf{a}_1+(x_{25}+y_{25})\,\mathbf{a}_2+z_{25}\,\mathbf{a}_3$	=	$(x_{25}a+z_{25}c\cos\beta)\,\hat{\mathbf{x}}+y_{25}b\,\hat{\mathbf{y}}+z_{25}c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O XIII
$\mathbf{B}_{90}$	=	$(-x_{25}-y_{25})\,\mathbf{a}_1+(-x_{25}+y_{25})\,\mathbf{a}_2+\left(\frac{1}{2}-z_{25}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta-x_{25}a-z_{25}c\cos\beta\right)\hat{\mathbf{x}}+y_{25}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{25}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O XIII
$\mathbf{B}_{91}$	=	$(-x_{25}+y_{25})\,\mathbf{a}_1+(-x_{25}-y_{25})\,\mathbf{a}_2-z_{25}\,\mathbf{a}_3$	=	$(-x_{25}a-z_{25}c\cos\beta)\,\hat{\mathbf{x}}-y_{25}b\,\hat{\mathbf{y}}-z_{25}c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O XIII
$\mathbf{B}_{92}$	=	$(x_{25}+y_{25})\,\mathbf{a}_1+(x_{25}-y_{25})\,\mathbf{a}_2+\left(\frac{1}{2}+z_{25}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta+x_{25}a+z_{25}c\cos\beta\right)\hat{\mathbf{x}}-y_{25}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{25}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O XIII
$\mathbf{B}_{93}$	=	$(x_{26}-y_{26})\,\mathbf{a}_1+(x_{26}+y_{26})\,\mathbf{a}_2+z_{26}\,\mathbf{a}_3$	=	$(x_{26}a+z_{26}c\cos\beta)\,\hat{\mathbf{x}}+y_{26}b\,\hat{\mathbf{y}}+z_{26}c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	S I
$\mathbf{B}_{94}$	=	$(-x_{26}-y_{26})\,\mathbf{a}_1+(-x_{26}+y_{26})\,\mathbf{a}_2+\left(\frac{1}{2}-z_{26}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta-x_{26}a-z_{26}c\cos\beta\right)\hat{\mathbf{x}}+y_{26}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{26}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	S I
$\mathbf{B}_{95}$	=	$(-x_{26}+y_{26})\,\mathbf{a}_1+(-x_{26}-y_{26})\,\mathbf{a}_2-z_{26}\,\mathbf{a}_3$	=	$(-x_{26}a-z_{26}c\cos\beta)\,\hat{\mathbf{x}}-y_{26}b\,\hat{\mathbf{y}}-z_{26}c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	S I
$\mathbf{B}_{96}$	=	$(x_{26}+y_{26})\,\mathbf{a}_1+(x_{26}-y_{26})\,\mathbf{a}_2+\left(\frac{1}{2}+z_{26}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta+x_{26}a+z_{26}c\cos\beta\right)\hat{\mathbf{x}}-y_{26}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{26}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	S I
$\mathbf{B}_{97}$	=	$(x_{27}-y_{27})\,\mathbf{a}_1+(x_{27}+y_{27})\,\mathbf{a}_2+z_{27}\,\mathbf{a}_3$	=	$(x_{27}a+z_{27}c\cos\beta)\,\hat{\mathbf{x}}+y_{27}b\,\hat{\mathbf{y}}+z_{27}c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	S II
$\mathbf{B}_{98}$	=	$(-x_{27}-y_{27})\,\mathbf{a}_1+(-x_{27}+y_{27})\,\mathbf{a}_2+\left(\frac{1}{2}-z_{27}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta-x_{27}a-z_{27}c\cos\beta\right)\hat{\mathbf{x}}+y_{27}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{27}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	S II
$\mathbf{B}_{99}$	=	$(-x_{27}+y_{27})\,\mathbf{a}_1+(-x_{27}-y_{27})\,\mathbf{a}_2-z_{27}\,\mathbf{a}_3$	=	$(-x_{27}a-z_{27}c\cos\beta)\,\hat{\mathbf{x}}-y_{27}b\,\hat{\mathbf{y}}-z_{27}c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	S II
$\mathbf{B}_{100}$	=	$(x_{27}+y_{27})\,\mathbf{a}_1+(x_{27}-y_{27})\,\mathbf{a}_2+\left(\frac{1}{2}+z_{27}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta+x_{27}a+z_{27}c\cos\beta\right)\hat{\mathbf{x}}-y_{27}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{27}\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	S II

References:

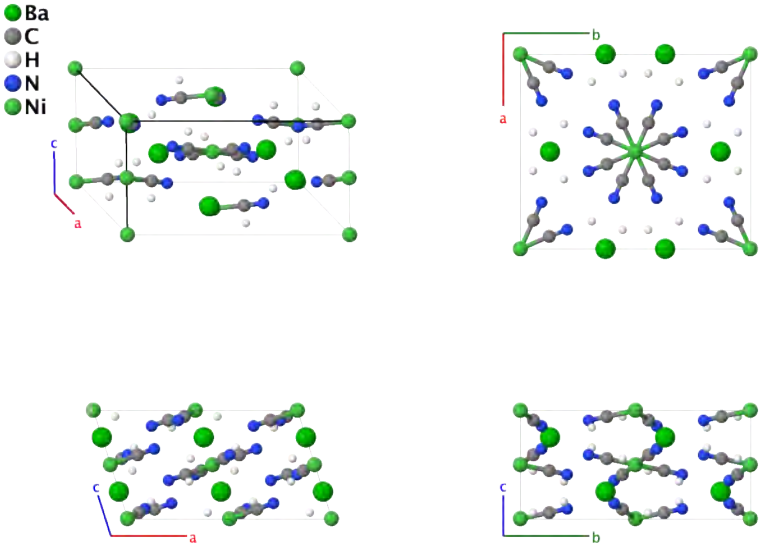
- B. Hertweck, G. Giester, and E. Libowitzky, *The crystal structures of the low-temperature phases of leonite-type compounds, $K_2Me(SO_4)_2{\cdot}4H_2O$ (Me^{2+} = Mg, Mn, Fe)*, Am. Mineral. **86**, 1282–1292 (2001), doi:10.2138/am-2001-1016.
- *$K_2Mn(SO_4)_2{\cdot}4H_2O$ ($K_2Mn[SO_4]_2[H_2O]_4$ mon1, T=185 K) Crystal Structure: Datasheet from "PAULING FILE Multinaries Edition – 2012" in SpringerMaterials (http://materials.springer.com/isp/crystallographic/docs/sd_1811721).* Copyright 2016 Springer-Verlag Berlin Heidelberg & Material Phases Data System (MPDS), Switzerland & National Institute for Materials Science (NIMS), Japan.

Geometry files:

- CIF: pp. 1568
- POSCAR: pp. 1568

$BaNi(CN)_4{\cdot}4H_2O$ ($H4_{22}$) Structure: AB4C4D4E_mC56_15_e_2f_2f_2f_a

http://aflow.org/prototype-encyclopedia/AB4C4D4E_mC56_15_e_2f_2f_2f_a



Prototype	:	$BaC_4(H_2O)_4N_4Ni$
AFLOW prototype label	:	AB4C4D4E_mC56_15_e_2f_2f_2f_a
***Strukturbericht* designation**	:	$H4_{22}$
Pearson symbol	:	mC56
Space group number	:	15
Space group symbol	:	$C2/c$
AFLOW prototype command	:	`aflow --proto=AB4C4D4E_mC56_15_e_2f_2f_2f_a --params=`$a, b/a, c/a, \beta, y_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8$

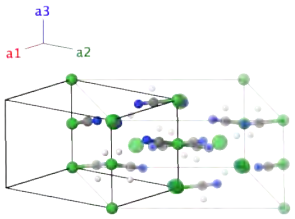
Other compounds with this structure

- $BaC_4N_4Pd{\cdot}4H_2O$ and $BaC_4N_4Pt{\cdot}4H_2O$

- We use the structure found by (Larsen, 1969). It is very similar to the structure found by (Brasseur, 1938) which was designated $H4_{22}$ by (Herrmann, 1941). The newer structure was able to separate the carbon and nitrogen atoms in the cyanide (CN) radical, and also move the barium atom from the (4*b*) to the (4*e*) Wyckoff position.

Base-centered Monoclinic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2}\,b\,\hat{\mathbf{y}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,b\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(4a)$	Ni
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\cos\beta\,\hat{\mathbf{x}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	$(4a)$	Ni
$\mathbf{B}_3$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Ba
$\mathbf{B}_4$	=	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Ba
$\mathbf{B}_5$	=	$(x_3 - y_3)\,\mathbf{a}_1 + (x_3 + y_3)\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	C I
$\mathbf{B}_6$	=	$(-x_3 - y_3)\,\mathbf{a}_1 + (-x_3 + y_3)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_3 a - z_3 c\cos\beta\right)\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	C I
$\mathbf{B}_7$	=	$(-x_3 + y_3)\,\mathbf{a}_1 + (-x_3 - y_3)\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	C I
$\mathbf{B}_8$	=	$(x_3 + y_3)\,\mathbf{a}_1 + (x_3 - y_3)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	C I
$\mathbf{B}_9$	=	$(x_4 - y_4)\,\mathbf{a}_1 + (x_4 + y_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	C II
$\mathbf{B}_{10}$	=	$(-x_4 - y_4)\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_4 a - z_4 c\cos\beta\right)\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	C II
$\mathbf{B}_{11}$	=	$(-x_4 + y_4)\,\mathbf{a}_1 + (-x_4 - y_4)\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	C II
$\mathbf{B}_{12}$	=	$(x_4 + y_4)\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	C II
$\mathbf{B}_{13}$	=	$(x_5 - y_5)\,\mathbf{a}_1 + (x_5 + y_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H_2O I
$\mathbf{B}_{14}$	=	$(-x_5 - y_5)\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_5 a - z_5 c\cos\beta\right)\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H_2O I
$\mathbf{B}_{15}$	=	$(-x_5 + y_5)\,\mathbf{a}_1 + (-x_5 - y_5)\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H_2O I
$\mathbf{B}_{16}$	=	$(x_5 + y_5)\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H_2O I
$\mathbf{B}_{17}$	=	$(x_6 - y_6)\,\mathbf{a}_1 + (x_6 + y_6)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H_2O II
$\mathbf{B}_{18}$	=	$(-x_6 - y_6)\,\mathbf{a}_1 + (-x_6 + y_6)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_6 a - z_6 c\cos\beta\right)\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H_2O II

$\mathbf{B}_{19}$	=	$(-x_6 + y_6)\,\mathbf{a}_1 + (-x_6 - y_6)\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H_2O II
$\mathbf{B}_{20}$	=	$(x_6 + y_6)\,\mathbf{a}_1 + (x_6 - y_6)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H_2O II
$\mathbf{B}_{21}$	=	$(x_7 - y_7)\,\mathbf{a}_1 + (x_7 + y_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	N I
$\mathbf{B}_{22}$	=	$(-x_7 - y_7)\,\mathbf{a}_1 + (-x_7 + y_7)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_7 a - z_7 c\cos\beta\right)\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	N I
$\mathbf{B}_{23}$	=	$(-x_7 + y_7)\,\mathbf{a}_1 + (-x_7 - y_7)\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	N I
$\mathbf{B}_{24}$	=	$(x_7 + y_7)\,\mathbf{a}_1 + (x_7 - y_7)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_7 a + z_7 c\cos\beta\right)\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	N I
$\mathbf{B}_{25}$	=	$(x_8 - y_8)\,\mathbf{a}_1 + (x_8 + y_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	N II
$\mathbf{B}_{26}$	=	$(-x_8 - y_8)\,\mathbf{a}_1 + (-x_8 + y_8)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_8 a - z_8 c\cos\beta\right)\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	N II
$\mathbf{B}_{27}$	=	$(-x_8 + y_8)\,\mathbf{a}_1 + (-x_8 - y_8)\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	N II
$\mathbf{B}_{28}$	=	$(x_8 + y_8)\,\mathbf{a}_1 + (x_8 - y_8)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_8 a + z_8 c\cos\beta\right)\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	N II

References:

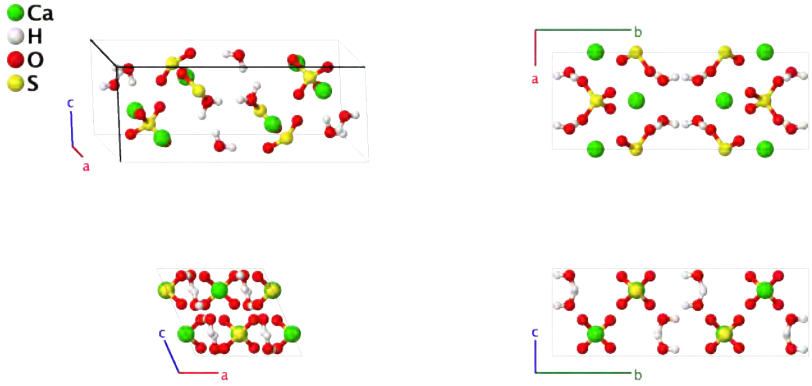
- F. K. Larsen, R. G. Hazell, and S. E. Rasmussen, *The Crystal Structure of Barium Tetracyanonickelate(II) Tetrahydrate*, Acta Chem. Scand. **23**, 61–69 (1969), doi:10.3891/acta.chem.scand.23-0061.
- M. H. Brasseur and M. A. de Rassenfosse, *Structure cristalline des cyanures doubles de baryum à base de platine, de palladium et de nickel*, Bull. Soc. Fr. Mineral. Cristallogr. **61**, 129–136 (1938), doi:10.3406/bulmi.1938.4433.
- K. Herrmann, ed., *Strukturbericht Band VI 1938* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1941).

Geometry files:

- CIF: pp. 1569
- POSCAR: pp. 1569

Gypsum ($CaSO_4 \cdot 2H_2O$, $H4_6$) Structure: AB4C6D_mC48_15_e_2f_3f_e

http://aflow.org/prototype-encyclopedia/AB4C6D_mC48_15_e_2f_3f_e

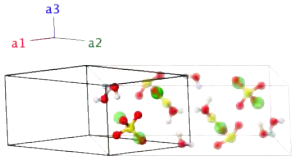


Prototype	:	CaH_4O_6S
AFLOW prototype label	:	AB4C6D_mC48_15_e_2f_3f_e
***Strukturbericht* designation**	:	$H4_6$
Pearson symbol	:	mC48
Space group number	:	15
Space group symbol	:	$C2/c$
AFLOW prototype command	:	`aflow --proto=AB4C6D_mC48_15_e_2f_3f_e --params=`$a, b/a, c/a, \beta, y_1, y_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7$

- (Onorato, 1929) determined that gypsum was in space group $C2/m$ #12, and (Hermann, 1937) assigned this to *Strukturbericht* $H4_6$. However, (Wooster, 1936) and others determined that the correct space group was $C2/c$ #15, and (Gottfried, 1938) assigned the $H4_6$ symbol to this structure. We use the 1 atmosphere data of (Comodi, 2008), who also determine the positions of the hydrogen atoms.

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Ca
$\mathbf{B}_2$	=	$y_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Ca
$\mathbf{B}_3$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S
$\mathbf{B}_4$	=	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	S

$\mathbf{B}_5$	=	$(x_3 - y_3)\,\mathbf{a}_1 + (x_3 + y_3)\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H I
$\mathbf{B}_6$	=	$(-x_3 - y_3)\,\mathbf{a}_1 + (-x_3 + y_3)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_3 a - z_3 c\cos\beta\right)\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H I
$\mathbf{B}_7$	=	$(-x_3 + y_3)\,\mathbf{a}_1 + (-x_3 - y_3)\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H I
$\mathbf{B}_8$	=	$(x_3 + y_3)\,\mathbf{a}_1 + (x_3 - y_3)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H I
$\mathbf{B}_9$	=	$(x_4 - y_4)\,\mathbf{a}_1 + (x_4 + y_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H II
$\mathbf{B}_{10}$	=	$(-x_4 - y_4)\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_4 a - z_4 c\cos\beta\right)\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H II
$\mathbf{B}_{11}$	=	$(-x_4 + y_4)\,\mathbf{a}_1 + (-x_4 - y_4)\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H II
$\mathbf{B}_{12}$	=	$(x_4 + y_4)\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	H II
$\mathbf{B}_{13}$	=	$(x_5 - y_5)\,\mathbf{a}_1 + (x_5 + y_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O I
$\mathbf{B}_{14}$	=	$(-x_5 - y_5)\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_5 a - z_5 c\cos\beta\right)\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O I
$\mathbf{B}_{15}$	=	$(-x_5 + y_5)\,\mathbf{a}_1 + (-x_5 - y_5)\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O I
$\mathbf{B}_{16}$	=	$(x_5 + y_5)\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O I
$\mathbf{B}_{17}$	=	$(x_6 - y_6)\,\mathbf{a}_1 + (x_6 + y_6)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{18}$	=	$(-x_6 - y_6)\,\mathbf{a}_1 + (-x_6 + y_6)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_6 a - z_6 c\cos\beta\right)\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{19}$	=	$(-x_6 + y_6)\,\mathbf{a}_1 + (-x_6 - y_6)\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{20}$	=	$(x_6 + y_6)\,\mathbf{a}_1 + (x_6 - y_6)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{21}$	=	$(x_7 - y_7)\,\mathbf{a}_1 + (x_7 + y_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{22}$	=	$(-x_7 - y_7)\,\mathbf{a}_1 + (-x_7 + y_7)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_7 a - z_7 c\cos\beta\right)\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{23}$	=	$(-x_7 + y_7)\,\mathbf{a}_1 + (-x_7 - y_7)\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{24}$	=	$(x_7 + y_7)\,\mathbf{a}_1 + (x_7 - y_7)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_7 a + z_7 c\cos\beta\right)\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III

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- W. A. Wooster, *On the Crystal Structure of Gypsum, $CaSO_4{\cdot}2H_2O$*, Zeitschrift für Kristallographie - Crystalline Materials **94**, 375–396 (1936), doi:10.1524/zkri.1936.94.1.375.
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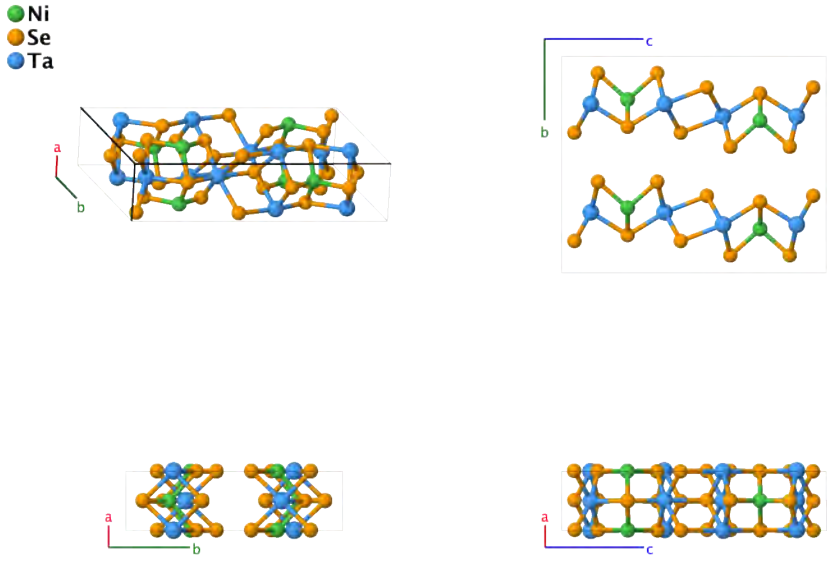
Geometry files:

- CIF: pp. 1570
- POSCAR: pp. 1570

Ta_2NiSe_5 Structure: AB5C2_mC32_15_e_e2f_f

http://aflow.org/prototype-encyclopedia/AB5C2_mC32_15_e_e2f_f

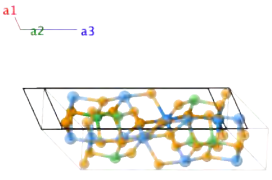


Prototype	:	$NiSe_5Ta_2$
AFLOW prototype label	:	AB5C2_mC32_15_e_e2f_f
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC32
Space group number	:	15
Space group symbol	:	$C2/c$
AFLOW prototype command	:	`aflow --proto=AB5C2_mC32_15_e_e2f_f --params=`$a, b/a, c/a, \beta, y_1, y_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5$

- Above 328 K this transforms into the Ta_2NiS_5 structure.
- When $\beta = 90°$, $x_3 = 1/2$, and $x_4 = x_5 = 0$, this becomes the Ta_2NiS_5 structure.

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$\mathbf{B}_1$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Ni
$\mathbf{B}_2$	=	$y_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Ni
$\mathbf{B}_3$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Se I
$\mathbf{B}_4$	=	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Se I
$\mathbf{B}_5$	=	$(x_3 - y_3)\,\mathbf{a}_1 + (x_3 + y_3)\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Se II
$\mathbf{B}_6$	=	$(-x_3 - y_3)\,\mathbf{a}_1 + (-x_3 + y_3)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_3 a - z_3 c\cos\beta\right)\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Se II
$\mathbf{B}_7$	=	$(-x_3 + y_3)\,\mathbf{a}_1 + (-x_3 - y_3)\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Se II
$\mathbf{B}_8$	=	$(x_3 + y_3)\,\mathbf{a}_1 + (x_3 - y_3)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Se II
$\mathbf{B}_9$	=	$(x_4 - y_4)\,\mathbf{a}_1 + (x_4 + y_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Se III
$\mathbf{B}_{10}$	=	$(-x_4 - y_4)\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_4 a - z_4 c\cos\beta\right)\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Se III
$\mathbf{B}_{11}$	=	$(-x_4 + y_4)\,\mathbf{a}_1 + (-x_4 - y_4)\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Se III
$\mathbf{B}_{12}$	=	$(x_4 + y_4)\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Se III
$\mathbf{B}_{13}$	=	$(x_5 - y_5)\,\mathbf{a}_1 + (x_5 + y_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Ta
$\mathbf{B}_{14}$	=	$(-x_5 - y_5)\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_5 a - z_5 c\cos\beta\right)\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Ta
$\mathbf{B}_{15}$	=	$(-x_5 + y_5)\,\mathbf{a}_1 + (-x_5 - y_5)\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Ta
$\mathbf{B}_{16}$	=	$(x_5 + y_5)\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Ta

References:

- S. A. Sunshine and J. A. Ibers, *Structure and physical properties of the new layered ternary chalcogenides tantalum nickel sulfide (Ta_2NiS_5) and tantalum nickel selenide (Ta_2NiSe_5)*, Inorg. Chem. **24**, 3611–3614 (1985), doi:10.1021/ic00216a027.

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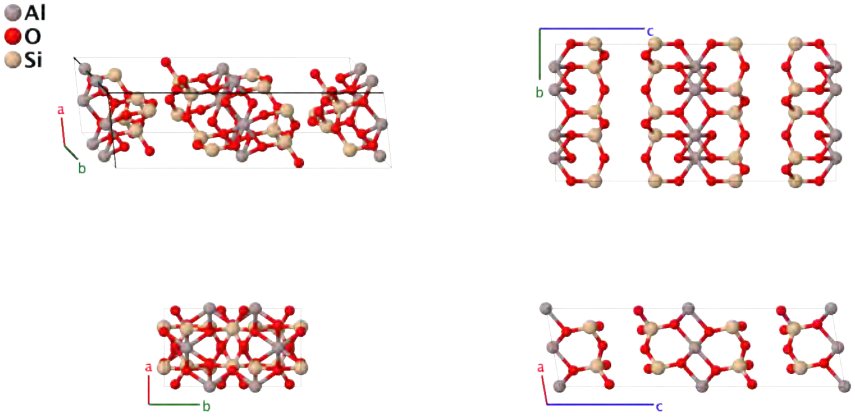
Geometry files:

- CIF: pp. 1570
- POSCAR: pp. 1570

Pyrophyllite [$AlSi_2O_5(OH)$, $S5_6$] Structure: AB5CD2_mC72_15_f_5f_f_2f

http://aflow.org/prototype-encyclopedia/AB5CD2_mC72_15_f_5f_f_2f



Prototype	:	$AlO_5(OH)Si_2$
AFLOW prototype label	:	AB5CD2_mC72_15_f_5f_f_2f
***Strukturbericht* designation**	:	$S5_6$
Pearson symbol	:	mC72
Space group number	:	15
Space group symbol	:	$C2/c$
AFLOW prototype command	:	`aflow --proto=AB5CD2_mC72_15_f_5f_f_2f --params=`$a, b/a, c/a, \beta, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9$

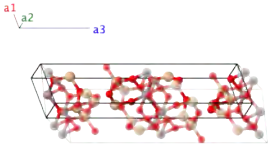
Other compounds with this structure

- $Mg_3Si_4O_{10}(OH)_2$ (talc)

- This a double-layer version of the pyrophyllite structure. (Brindley, 1259) has shown that some samples have only one layer, with a corresponding reduction in the c lattice vector. Both Brindley and (Lee, 1981) found pyrophyllite samples with space group $P\bar{1}$ #2.
- Talc ($Mg_3Si_4O_{10}(OH)_2$) is isostructural with pyrophyllite, with magnesium replacing aluminum on the ($8f$) site and the addition of another magnesium atom on the ($4a$) Wyckoff site.

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$= (x_1 - y_1)\,\mathbf{a}_1 + (x_1 + y_1)\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	$=$	$(x_1 a + z_1 c\cos\beta)\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Al
$\mathbf{B}_2$	$= (-x_1 - y_1)\,\mathbf{a}_1 + (-x_1 + y_1)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}c\cos\beta - x_1 a - z_1 c\cos\beta\right)\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Al
$\mathbf{B}_3$	$= (-x_1 + y_1)\,\mathbf{a}_1 + (-x_1 - y_1)\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	$=$	$(-x_1 a - z_1 c\cos\beta)\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} - z_1 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Al
$\mathbf{B}_4$	$= (x_1 + y_1)\,\mathbf{a}_1 + (x_1 - y_1)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}c\cos\beta + x_1 a + z_1 c\cos\beta\right)\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Al
$\mathbf{B}_5$	$= (x_2 - y_2)\,\mathbf{a}_1 + (x_2 + y_2)\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	$=$	$(x_2 a + z_2 c\cos\beta)\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O I
$\mathbf{B}_6$	$= (-x_2 - y_2)\,\mathbf{a}_1 + (-x_2 + y_2)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}c\cos\beta - x_2 a - z_2 c\cos\beta\right)\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O I
$\mathbf{B}_7$	$= (-x_2 + y_2)\,\mathbf{a}_1 + (-x_2 - y_2)\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	$=$	$(-x_2 a - z_2 c\cos\beta)\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} - z_2 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O I
$\mathbf{B}_8$	$= (x_2 + y_2)\,\mathbf{a}_1 + (x_2 - y_2)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}c\cos\beta + x_2 a + z_2 c\cos\beta\right)\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O I
$\mathbf{B}_9$	$= (x_3 - y_3)\,\mathbf{a}_1 + (x_3 + y_3)\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$=$	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{10}$	$= (-x_3 - y_3)\,\mathbf{a}_1 + (-x_3 + y_3)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}c\cos\beta - x_3 a - z_3 c\cos\beta\right)\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{11}$	$= (-x_3 + y_3)\,\mathbf{a}_1 + (-x_3 - y_3)\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	$=$	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{12}$	$= (x_3 + y_3)\,\mathbf{a}_1 + (x_3 - y_3)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{13}$	$= (x_4 - y_4)\,\mathbf{a}_1 + (x_4 + y_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$=$	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{14}$	$= (-x_4 - y_4)\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}c\cos\beta - x_4 a - z_4 c\cos\beta\right)\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{15}$	$= (-x_4 + y_4)\,\mathbf{a}_1 + (-x_4 - y_4)\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	$=$	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{16}$	$= (x_4 + y_4)\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{17}$	$= (x_5 - y_5)\,\mathbf{a}_1 + (x_5 + y_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	$=$	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O IV
$\mathbf{B}_{18}$	$= (-x_5 - y_5)\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}c\cos\beta - x_5 a - z_5 c\cos\beta\right)\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O IV
$\mathbf{B}_{19}$	$= (-x_5 + y_5)\,\mathbf{a}_1 + (-x_5 - y_5)\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	$=$	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O IV
$\mathbf{B}_{20}$	$= (x_5 + y_5)\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O IV
$\mathbf{B}_{21}$	$= (x_6 - y_6)\,\mathbf{a}_1 + (x_6 + y_6)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	$=$	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O V

$\mathbf{B}_{22}$	=	$(-x_6 - y_6)\,\mathbf{a}_1 + (-x_6 + y_6)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_6 a - z_6 c\cos\beta\right)\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O V
$\mathbf{B}_{23}$	=	$(-x_6 + y_6)\,\mathbf{a}_1 + (-x_6 - y_6)\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O V
$\mathbf{B}_{24}$	=	$(x_6 + y_6)\,\mathbf{a}_1 + (x_6 - y_6)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O V
$\mathbf{B}_{25}$	=	$(x_7 - y_7)\,\mathbf{a}_1 + (x_7 + y_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$(x_7 a + z_7 c\cos\beta)\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	OH
$\mathbf{B}_{26}$	=	$(-x_7 - y_7)\,\mathbf{a}_1 + (-x_7 + y_7)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_7 a - z_7 c\cos\beta\right)\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	OH
$\mathbf{B}_{27}$	=	$(-x_7 + y_7)\,\mathbf{a}_1 + (-x_7 - y_7)\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$(-x_7 a - z_7 c\cos\beta)\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	OH
$\mathbf{B}_{28}$	=	$(x_7 + y_7)\,\mathbf{a}_1 + (x_7 - y_7)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_7 a + z_7 c\cos\beta\right)\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	OH
$\mathbf{B}_{29}$	=	$(x_8 - y_8)\,\mathbf{a}_1 + (x_8 + y_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$(x_8 a + z_8 c\cos\beta)\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Si I
$\mathbf{B}_{30}$	=	$(-x_8 - y_8)\,\mathbf{a}_1 + (-x_8 + y_8)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_8 a - z_8 c\cos\beta\right)\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Si I
$\mathbf{B}_{31}$	=	$(-x_8 + y_8)\,\mathbf{a}_1 + (-x_8 - y_8)\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$(-x_8 a - z_8 c\cos\beta)\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Si I
$\mathbf{B}_{32}$	=	$(x_8 + y_8)\,\mathbf{a}_1 + (x_8 - y_8)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_8 a + z_8 c\cos\beta\right)\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Si I
$\mathbf{B}_{33}$	=	$(x_9 - y_9)\,\mathbf{a}_1 + (x_9 + y_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$(x_9 a + z_9 c\cos\beta)\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Si II
$\mathbf{B}_{34}$	=	$(-x_9 - y_9)\,\mathbf{a}_1 + (-x_9 + y_9)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_9 a - z_9 c\cos\beta\right)\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_9\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Si II
$\mathbf{B}_{35}$	=	$(-x_9 + y_9)\,\mathbf{a}_1 + (-x_9 - y_9)\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$(-x_9 a - z_9 c\cos\beta)\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} - z_9 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Si II
$\mathbf{B}_{36}$	=	$(x_9 + y_9)\,\mathbf{a}_1 + (x_9 - y_9)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_9 a + z_9 c\cos\beta\right)\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Si II

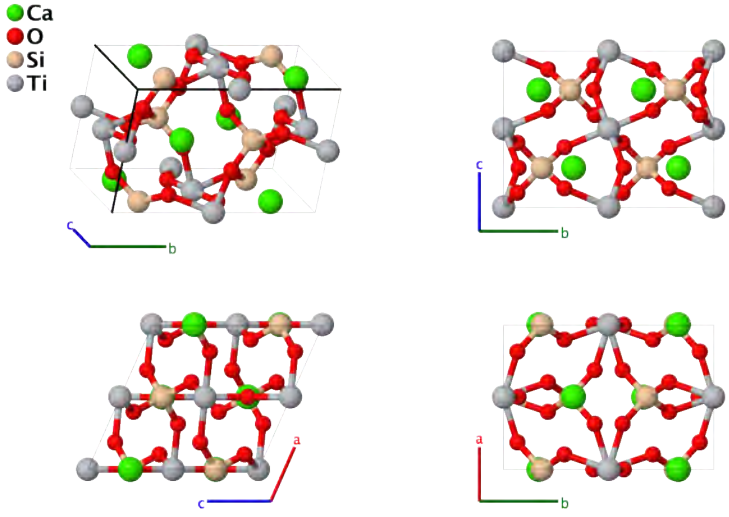
References:
- J. W. Gruner, *The Crystal Structures of Talc and Pyrophyllite*, Zeitschrift für Kristallographie - Crystalline Materials **88**, 412–419 (1934), doi:10.1524/zkri.1934.88.1.412.
- G. W. Brindley and R. Wardle, *Monoclinic and triclinic forms of pyrophyllite and pyrophyllite anhydride*, Am. Mineral. **55**, 1259–1272 (1970).
- J. H. Lee and S. Guggenheim, *Single crystal X-ray refinement of pyrophyllite-1Tc*, Am. Mineral. **66**, 350–357 (1981).

Found in:
- C. Gottfried and F. Schossberger, eds., *Strukturbericht Band III 1933-1935* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:
- CIF: pp. 1571
- POSCAR: pp. 1571

Titanite ($CaTiSiO_5$, $S0_6$) Structure: AB5CD_mC32_15_e_e2f_e_b

http://aflow.org/prototype-encyclopedia/AB5CD_mC32_15_e_e2f_e_b

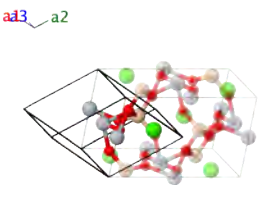


Prototype	:	CaO_5SiTi
AFLOW prototype label	:	AB5CD_mC32_15_e_e2f_e_b
***Strukturbericht* designation**	:	$S0_6$
Pearson symbol	:	mC32
Space group number	:	15
Space group symbol	:	$C2/c$
AFLOW prototype command	:	`aflow --proto=AB5CD_mC32_15_e_e2f_e_b --params=`$a, b/a, c/a, \beta, y_2, y_3, y_4, x_5, y_5, z_5, x_6, y_6, z_6$

- We use the data from sample M28658 to construct our titanite cell.
- (Hermann, 1937) places the titanium atoms at the (4*c*) Wyckoff position rather than the (4*b*) position favored by (Hawthorne, 1991). The biggest difference between the two structures is the distance between the silicon and titanium atoms, which is 3.81 Å in Hawthorne and a rather short 2 Å in Hermann. We therefore favor the Hawthorne structure, and, since it does not otherwise substantially change the results, we continue to use $S0_6$ to describe it.
- (Hermann, 1937) also listed this as the $H5_6$ structure in the index.

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	(4*b*)	Ti
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(a + c\cos\beta)\,\hat{\mathbf{x}} + \frac{1}{2}c\sin\beta\,\hat{\mathbf{z}}$	(4*b*)	Ti
$\mathbf{B}_3$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Ca
$\mathbf{B}_4$	=	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Ca
$\mathbf{B}_5$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	O I
$\mathbf{B}_6$	=	$y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	O I
$\mathbf{B}_7$	=	$-y_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Si
$\mathbf{B}_8$	=	$y_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Si
$\mathbf{B}_9$	=	$(x_5 - y_5)\,\mathbf{a}_1 + (x_5 + y_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	O II
$\mathbf{B}_{10}$	=	$(-x_5 - y_5)\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_5 a - z_5 c\cos\beta\right)\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	O II
$\mathbf{B}_{11}$	=	$(-x_5 + y_5)\,\mathbf{a}_1 + (-x_5 - y_5)\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	O II
$\mathbf{B}_{12}$	=	$(x_5 + y_5)\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	O II
$\mathbf{B}_{13}$	=	$(x_6 - y_6)\,\mathbf{a}_1 + (x_6 + y_6)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	O III
$\mathbf{B}_{14}$	=	$(-x_6 - y_6)\,\mathbf{a}_1 + (-x_6 + y_6)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_6 a - z_6 c\cos\beta\right)\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	O III
$\mathbf{B}_{15}$	=	$(-x_6 + y_6)\,\mathbf{a}_1 + (-x_6 - y_6)\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	O III
$\mathbf{B}_{16}$	=	$(x_6 + y_6)\,\mathbf{a}_1 + (x_6 - y_6)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	(8*f*)	O III

References:

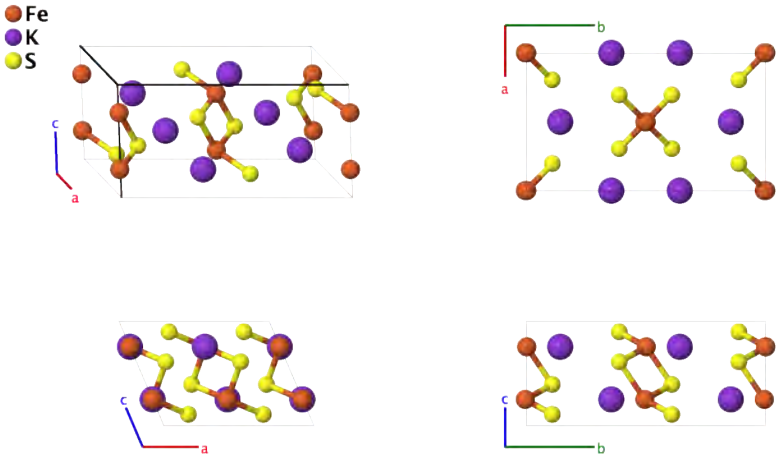
- F. C. Hawthorne, L. A. Groat, M. Raudsepp, N. A. Ball, M. Kimata, F. D. Spike, R. Gaba, N. M. Halden, G. R. Lumpkin, R. C. Ewing, R. B. Greegor, F. W. Lytle, T. Scott Ercit, G. R. Rossman, F. J. Wicks, R. A. Ramik, B. L. Sherriff, M. E. Fleet, and C. McCammon, *Alpha-decay damage in titanite*, Am. Mineral. **76**, 370–396 (1991).
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1571
- POSCAR: pp. 1572

$KFeS_2$ ($F5_a$) Structure: ABC2_mC16_15_e_e_f

http://aflow.org/prototype-encyclopedia/ABC2_mC16_15_e_e_f



Prototype	:	$FeKS_2$
AFLOW prototype label	:	ABC2_mC16_15_e_e_f
***Strukturbericht* designation**	:	$F5_a$
Pearson symbol	:	mC16
Space group number	:	15
Space group symbol	:	$C2/c$
AFLOW prototype command	:	`aflow --proto=ABC2_mC16_15_e_e_f --params=`$a, b/a, c/a, \beta, y_1, y_2, x_3, y_3, z_3$

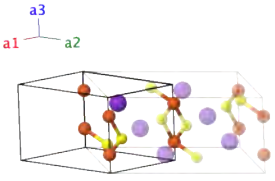
Other compounds with this structure

- $RbFeS_2$, $KFeSe_2$, $RbFeSe_2$, $CsGaS_2$, and $CsGaSe_2$

- This structure forms cages of Fe atoms surrounded by tetrahedra of S atoms.

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Fe
$\mathbf{B}_2$	=	$y_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Fe
$\mathbf{B}_3$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	K

$\mathbf{B}_4$	=	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	K
$\mathbf{B}_5$	=	$(x_3 - y_3)\,\mathbf{a}_1 + (x_3 + y_3)\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	S
$\mathbf{B}_6$	=	$(-x_3 - y_3)\,\mathbf{a}_1 + (-x_3 + y_3)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_3 a - z_3 c\cos\beta\right)\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	S
$\mathbf{B}_7$	=	$(-x_3 + y_3)\,\mathbf{a}_1 + (-x_3 - y_3)\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	S
$\mathbf{B}_8$	=	$(x_3 + y_3)\,\mathbf{a}_1 + (x_3 - y_3)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	S

References:

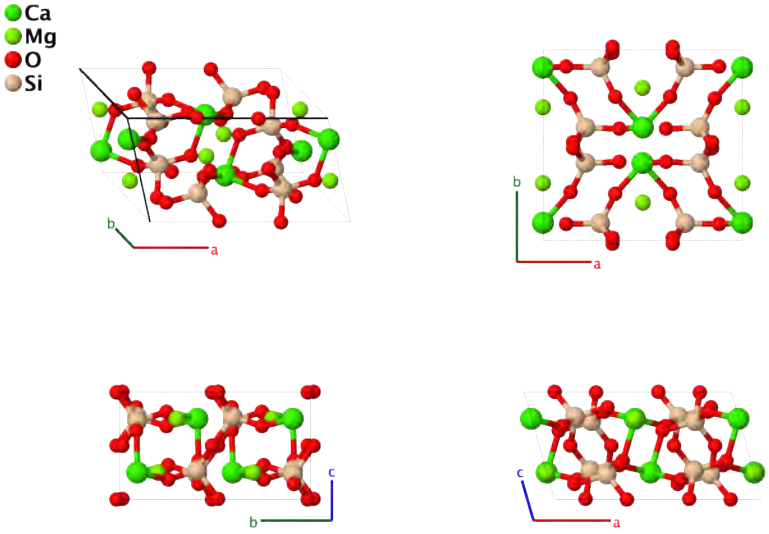
- W. Bronger, A. Kyas, and P. Müller, *The antiferromagnetic structures of $KFeS_2$, $RbFeS_2$, $KFeSe_2$, and $RbFeSe_2$ and the correlation between magnetic moments and crystal field calculations*, J. Solid State Chem. **70**, 262–270 (1987), doi:10.1016/0022-4596(87)90065-X.

Geometry files:

- CIF: pp. 1572
- POSCAR: pp. 1572

Diopside [$CaMg(SiO_3)_2$, $S4_1$] Structure: ABC6D2_mC40_15_e_e_3f_f

http://aflow.org/prototype-encyclopedia/ABC6D2_mC40_15_e_e_3f_f.S4_1



Prototype	:	$CaMgO_6Si_2$
AFLOW prototype label	:	ABC6D2_mC40_15_e_e_3f_f
***Strukturbericht* designation**	:	$S4_1$
Pearson symbol	:	mC40
Space group number	:	15
Space group symbol	:	$C2/c$
AFLOW prototype command	:	`aflow --proto=ABC6D2_mC40_15_e_e_3f_f --params=`$a, b/a, c/a, \beta, y_1, y_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6$

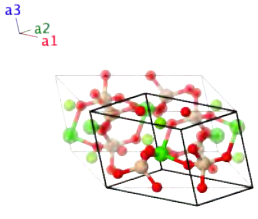
Other compounds with this structure

- $(Ca,Na)(Mg,Fe,Al,Ti)(Si,Al)_2O_6$ (augite)

- (Finger, 1976) list the two ($4e$) positions as 'M1' and 'M2', and both are mixtures of calcium and magnesium. For visual clarity we desigated the first ($4e$) position as calcium and the second as magnesium.
- This structure has the same AFLOW label as esseneite. The structures are generated by the same symmetry operations with different sets of parameters (`--params`) specified in their corresponding CIF files.

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Ca
$\mathbf{B}_2$	=	$y_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Ca
$\mathbf{B}_3$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Mg
$\mathbf{B}_4$	=	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	$(4e)$	Mg
$\mathbf{B}_5$	=	$(x_3 - y_3)\,\mathbf{a}_1 + (x_3 + y_3)\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$(x_3 a + z_3 c\cos\beta)\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O I
$\mathbf{B}_6$	=	$(-x_3 - y_3)\,\mathbf{a}_1 + (-x_3 + y_3)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_3 a - z_3 c\cos\beta\right)\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O I
$\mathbf{B}_7$	=	$(-x_3 + y_3)\,\mathbf{a}_1 + (-x_3 - y_3)\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$(-x_3 a - z_3 c\cos\beta)\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O I
$\mathbf{B}_8$	=	$(x_3 + y_3)\,\mathbf{a}_1 + (x_3 - y_3)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_3 a + z_3 c\cos\beta\right)\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O I
$\mathbf{B}_9$	=	$(x_4 - y_4)\,\mathbf{a}_1 + (x_4 + y_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$(x_4 a + z_4 c\cos\beta)\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{10}$	=	$(-x_4 - y_4)\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_4 a - z_4 c\cos\beta\right)\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{11}$	=	$(-x_4 + y_4)\,\mathbf{a}_1 + (-x_4 - y_4)\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$(-x_4 a - z_4 c\cos\beta)\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{12}$	=	$(x_4 + y_4)\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_4 a + z_4 c\cos\beta\right)\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{13}$	=	$(x_5 - y_5)\,\mathbf{a}_1 + (x_5 + y_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$(x_5 a + z_5 c\cos\beta)\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{14}$	=	$(-x_5 - y_5)\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_5 a - z_5 c\cos\beta\right)\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{15}$	=	$(-x_5 + y_5)\,\mathbf{a}_1 + (-x_5 - y_5)\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$(-x_5 a - z_5 c\cos\beta)\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{16}$	=	$(x_5 + y_5)\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_5 a + z_5 c\cos\beta\right)\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{17}$	=	$(x_6 - y_6)\,\mathbf{a}_1 + (x_6 + y_6)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$(x_6 a + z_6 c\cos\beta)\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Si
$\mathbf{B}_{18}$	=	$(-x_6 - y_6)\,\mathbf{a}_1 + (-x_6 + y_6)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta - x_6 a - z_6 c\cos\beta\right)\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Si
$\mathbf{B}_{19}$	=	$(-x_6 + y_6)\,\mathbf{a}_1 + (-x_6 - y_6)\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$(-x_6 a - z_6 c\cos\beta)\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Si
$\mathbf{B}_{20}$	=	$(x_6 + y_6)\,\mathbf{a}_1 + (x_6 - y_6)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}c\cos\beta + x_6 a + z_6 c\cos\beta\right)\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\sin\beta\,\hat{\mathbf{z}}$	$(8f)$	Si

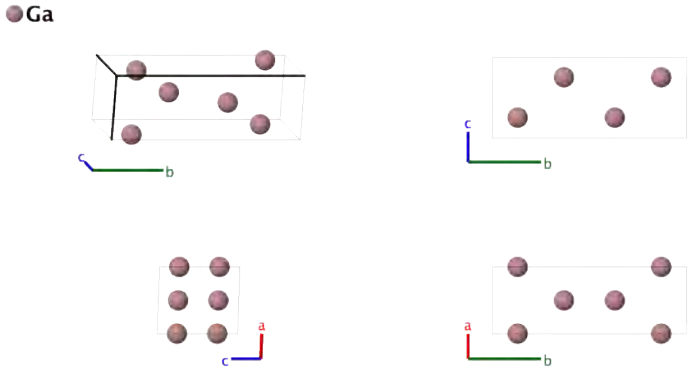
References:

- L. W. Finger and Y. Ohashi, *The thermal expansion of diopside to 800°C and a refinement of the crystal structure at 700°C*, Am. Mineral. **61**, 303–310 (1976).

Geometry files:
- CIF: pp. 1572
- POSCAR: pp. 1573

β-Ga (*obsolete*) Structure: A_mC4_15_e

http://aflow.org/prototype-encyclopedia/A_mC4_15_e

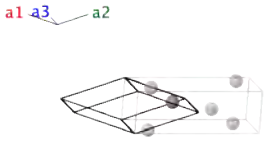


Prototype	:	Ga
AFLOW prototype label	:	A_mC4_15_e
***Strukturbericht* designation**	:	None
Pearson symbol	:	mC4
Space group number	:	15
Space group symbol	:	$C2/c$
AFLOW prototype command	:	`aflow --proto=A_mC4_15_e --params=`$a, b/a, c/a, \beta, y_1$

- β-Ga was proposed as a metastable structure of gallium, visible for short times at atomospheric pressure. The true high-pressure structure of gallium is isostructural with indium (*A*6).

Base-centered Monoclinic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\cos\beta\,\hat{\mathbf{x}} + c\sin\beta\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\cos\beta\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + \frac{1}{4}c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Ga
$\mathbf{B}_2$	=	$y_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\cos\beta\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \frac{3}{4}c\sin\beta\,\hat{\mathbf{z}}$	(4*e*)	Ga

References:

- L. Bosio, A. Defrain, H. Curien, and A. Rimsky, *Structure cristalline du gallium β*, Acta Crystallogr. Sect. B Struct. Sci. **25**, 995 (1969), doi:10.1107/S0567740869003360.

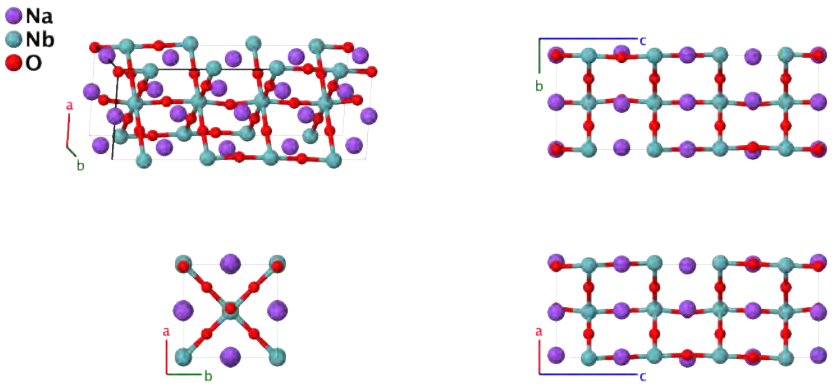
Found in:

- J. Donohue, *The Structures of the Elements* (Robert E. Krieger Publishing Company, Malabar, Florida, 1982). Reprint of the 1974 John Wilely & Sons edition.

Geometry files:
- CIF: pp. 1573
- POSCAR: pp. 1573

$NaNbO_3$ Structure: ABC3_oP40_17_abcd_2e_abcd4e

http://aflow.org/prototype-encyclopedia/ABC3_oP40_17_abcd_2e_abcd4e

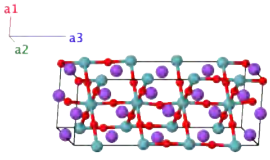


Prototype	:	$NaNbO_3$
AFLOW prototype label	:	ABC3_oP40_17_abcd_2e_abcd4e
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP40
Space group number	:	17
Space group symbol	:	$P222_1$
AFLOW prototype command	:	`aflow --proto=ABC3_oP40_17_abcd_2e_abcd4e --params=`$a, b/a, c/a, x_1, x_2, x_3, x_4, y_5, y_6, y_7, y_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}$

- (Downs, 2003) identifies this as a "possible polymorph of Lueshite."
- If the AFLOW parameters are set to `--params=`$a, 1, \sqrt{8}, 1/2, 0, 0, 1/2, 1/2, 0, 0, 1/2, 0, 0, 3/8, 1/2, 1/2, 3/8, 1/4, 1/4, 3/8, 3/4, 1/4, 3/8, 1/4, 3/4, 3/8, 3/4, 3/4, 3/8$ then the structure is equivalent to Cubic Perovskite $E2_1$.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1$	=	$x_1 a\,\hat{\mathbf{x}}$	$(2a)$	Na I
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(2a)$	Na I
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1$	=	$x_2 a\,\hat{\mathbf{x}}$	$(2a)$	O I
$\mathbf{B}_4$	=	$-x_2\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(2a)$	O I
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$x_3 a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$	$(2b)$	Na II
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(2b)$	Na II

$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$x_4 a\,\hat{\mathbf{x}} + \frac{1}{2} b\,\hat{\mathbf{y}}$	(2*b*)	O II
$\mathbf{B}_8$	=	$-x_4\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + \frac{1}{2} b\,\hat{\mathbf{y}} + \frac{1}{2} c\,\hat{\mathbf{z}}$	(2*b*)	O II
$\mathbf{B}_9$	=	$y_5\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_5 b\,\hat{\mathbf{y}} + \frac{1}{4} c\,\hat{\mathbf{z}}$	(2*c*)	Na III
$\mathbf{B}_{10}$	=	$-y_5\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_5 b\,\hat{\mathbf{y}} + \frac{3}{4} c\,\hat{\mathbf{z}}$	(2*c*)	Na III
$\mathbf{B}_{11}$	=	$y_6\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_6 b\,\hat{\mathbf{y}} + \frac{1}{4} c\,\hat{\mathbf{z}}$	(2*c*)	O III
$\mathbf{B}_{12}$	=	$-y_6\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_6 b\,\hat{\mathbf{y}} + \frac{3}{4} c\,\hat{\mathbf{z}}$	(2*c*)	O III
$\mathbf{B}_{13}$	=	$\frac{1}{2}\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2} a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + \frac{1}{4} c\,\hat{\mathbf{z}}$	(2*d*)	Na IV
$\mathbf{B}_{14}$	=	$\frac{1}{2}\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2} a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \frac{3}{4} c\,\hat{\mathbf{z}}$	(2*d*)	Na IV
$\mathbf{B}_{15}$	=	$\frac{1}{2}\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2} a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + \frac{1}{4} c\,\hat{\mathbf{z}}$	(2*d*)	O IV
$\mathbf{B}_{16}$	=	$\frac{1}{2}\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2} a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \frac{3}{4} c\,\hat{\mathbf{z}}$	(2*d*)	O IV
$\mathbf{B}_{17}$	=	$x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(4*e*)	Nb I
$\mathbf{B}_{18}$	=	$-x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right) c\,\hat{\mathbf{z}}$	(4*e*)	Nb I
$\mathbf{B}_{19}$	=	$-x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + \left(\frac{1}{2} - z_9\right)\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_9\right) c\,\hat{\mathbf{z}}$	(4*e*)	Nb I
$\mathbf{B}_{20}$	=	$x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}}$	(4*e*)	Nb I
$\mathbf{B}_{21}$	=	$x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$x_{10} a\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}}$	(4*e*)	Nb II
$\mathbf{B}_{22}$	=	$-x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{10}\right)\mathbf{a}_3$	=	$-x_{10} a\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{10}\right) c\,\hat{\mathbf{z}}$	(4*e*)	Nb II
$\mathbf{B}_{23}$	=	$-x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{10}\right)\mathbf{a}_3$	=	$-x_{10} a\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{10}\right) c\,\hat{\mathbf{z}}$	(4*e*)	Nb II
$\mathbf{B}_{24}$	=	$x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3$	=	$x_{10} a\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} - z_{10} c\,\hat{\mathbf{z}}$	(4*e*)	Nb II
$\mathbf{B}_{25}$	=	$x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$x_{11} a\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + z_{11} c\,\hat{\mathbf{z}}$	(4*e*)	O V
$\mathbf{B}_{26}$	=	$-x_{11}\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{11}\right)\mathbf{a}_3$	=	$-x_{11} a\,\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{11}\right) c\,\hat{\mathbf{z}}$	(4*e*)	O V
$\mathbf{B}_{27}$	=	$-x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{11}\right)\mathbf{a}_3$	=	$-x_{11} a\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{11}\right) c\,\hat{\mathbf{z}}$	(4*e*)	O V
$\mathbf{B}_{28}$	=	$x_{11}\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3$	=	$x_{11} a\,\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} - z_{11} c\,\hat{\mathbf{z}}$	(4*e*)	O V
$\mathbf{B}_{29}$	=	$x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3$	=	$x_{12} a\,\hat{\mathbf{x}} + y_{12} b\,\hat{\mathbf{y}} + z_{12} c\,\hat{\mathbf{z}}$	(4*e*)	O VI
$\mathbf{B}_{30}$	=	$-x_{12}\,\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{12}\right)\mathbf{a}_3$	=	$-x_{12} a\,\hat{\mathbf{x}} - y_{12} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{12}\right) c\,\hat{\mathbf{z}}$	(4*e*)	O VI
$\mathbf{B}_{31}$	=	$-x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{12}\right)\mathbf{a}_3$	=	$-x_{12} a\,\hat{\mathbf{x}} + y_{12} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{12}\right) c\,\hat{\mathbf{z}}$	(4*e*)	O VI
$\mathbf{B}_{32}$	=	$x_{12}\,\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3$	=	$x_{12} a\,\hat{\mathbf{x}} - y_{12} b\,\hat{\mathbf{y}} - z_{12} c\,\hat{\mathbf{z}}$	(4*e*)	O VI
$\mathbf{B}_{33}$	=	$x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3$	=	$x_{13} a\,\hat{\mathbf{x}} + y_{13} b\,\hat{\mathbf{y}} + z_{13} c\,\hat{\mathbf{z}}$	(4*e*)	O VII
$\mathbf{B}_{34}$	=	$-x_{13}\,\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{13}\right)\mathbf{a}_3$	=	$-x_{13} a\,\hat{\mathbf{x}} - y_{13} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{13}\right) c\,\hat{\mathbf{z}}$	(4*e*)	O VII
$\mathbf{B}_{35}$	=	$-x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{13}\right)\mathbf{a}_3$	=	$-x_{13} a\,\hat{\mathbf{x}} + y_{13} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{13}\right) c\,\hat{\mathbf{z}}$	(4*e*)	O VII
$\mathbf{B}_{36}$	=	$x_{13}\,\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 - z_{13}\,\mathbf{a}_3$	=	$x_{13} a\,\hat{\mathbf{x}} - y_{13} b\,\hat{\mathbf{y}} - z_{13} c\,\hat{\mathbf{z}}$	(4*e*)	O VII
$\mathbf{B}_{37}$	=	$x_{14}\,\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3$	=	$x_{14} a\,\hat{\mathbf{x}} + y_{14} b\,\hat{\mathbf{y}} + z_{14} c\,\hat{\mathbf{z}}$	(4*e*)	O VIII
$\mathbf{B}_{38}$	=	$-x_{14}\,\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{14}\right)\mathbf{a}_3$	=	$-x_{14} a\,\hat{\mathbf{x}} - y_{14} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{14}\right) c\,\hat{\mathbf{z}}$	(4*e*)	O VIII
$\mathbf{B}_{39}$	=	$-x_{14}\,\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{14}\right)\mathbf{a}_3$	=	$-x_{14} a\,\hat{\mathbf{x}} + y_{14} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{14}\right) c\,\hat{\mathbf{z}}$	(4*e*)	O VIII
$\mathbf{B}_{40}$	=	$x_{14}\,\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 - z_{14}\,\mathbf{a}_3$	=	$x_{14} a\,\hat{\mathbf{x}} - y_{14} b\,\hat{\mathbf{y}} - z_{14} c\,\hat{\mathbf{z}}$	(4*e*)	O VIII

References:

- P. Vousden, *The Structure of Ferroelectric Sodium Niobate at Room Temperature*, Acta Cryst. **4**, 545–551 (1951),

doi:10.1107/S0365110X51001768.

Found in:

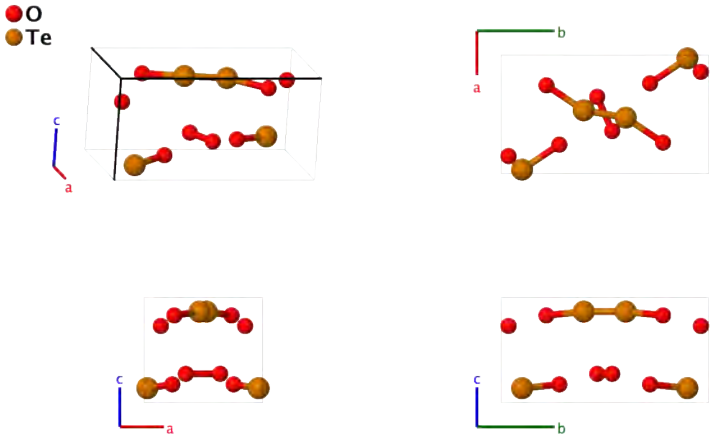
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1573
- POSCAR: pp. 1574

γ-TeO_2 Structure: A2B_oP12_18_2c_c

http://aflow.org/prototype-encyclopedia/A2B_oP12_18_2c_c

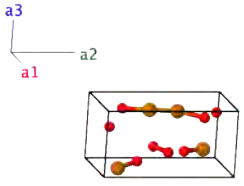


Prototype	:	O_2Te
AFLOW prototype label	:	A2B_oP12_18_2c_c
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP12
Space group number	:	18
Space group symbol	:	$P2_12_12$
AFLOW prototype command	:	`aflow --proto=A2B_oP12_18_2c_c --params=`$a, b/a, c/a, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3$

- This is a metastable state of TeO_2, similar to α-TeO_2 paratelluride. (Champarnaud-Mesjard, 2000) labels the atomic positions as being on the '(4*a*)' Wyckoff position, but this does not exist for space group $P2_12_12$ #18. The atoms are actually on the general site for this space group, (4*c*).

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_3$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_1\right)\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_1\right)b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_4$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_1\right)\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_1\right)b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*c*)	O II

$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_7$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right)b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right)b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*c*)	Te
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*c*)	Te
$\mathbf{B}_{11}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(4*c*)	Te
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(4*c*)	Te

References:

- J. C. Champarnaud-Mesjard, S. Blanchandin, P. Thomas, A. Mirgorodsky, T. Merle-Méjean, and B. Frit, *Crystal structure, Raman spectrum and lattice dynamics of a new metastable form of tellurium dioxide: γ-TeO_2*, J. Phys. Chem. Solids **61**, 1499–1507 (2000), doi:10.1016/S0022-3697(00)00012-3.

Found in:

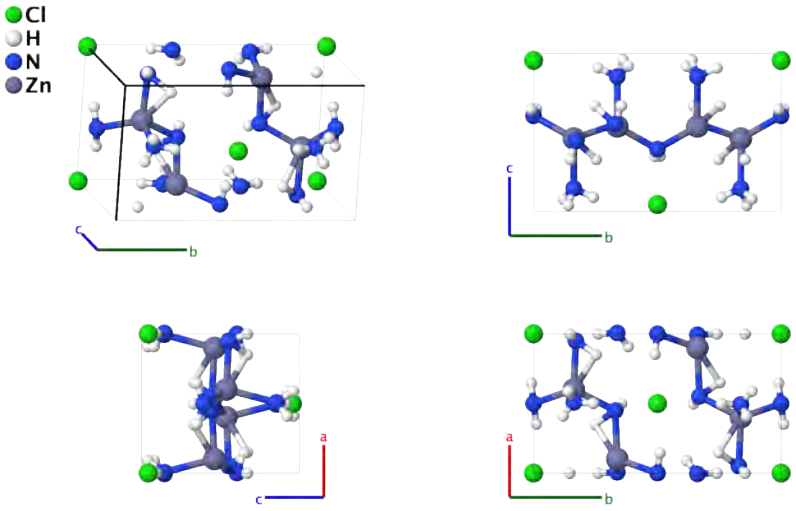
- M. Ceriotti, F. Pietrucci, and M. Bernasconi, Ab initio *study of the vibrational properties of crystalline TeO_2: The α, β, and γ phases*, Phys. Rev. B **73**, 104304 (2006), doi:10.1103/PhysRevB.73.104304.

Geometry files:

- CIF: pp. 1574
- POSCAR: pp. 1574

Diamminetriamidodizinc Chloride ($[Zn_2(NH_3)_2(NH_2)_3]Cl$) Structure: AB12C5D2_oP40_18_a_6c_b2c_c

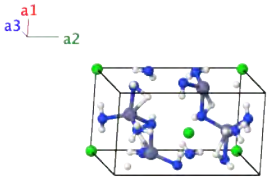
http://aflow.org/prototype-encyclopedia/AB12C5D2_oP40_18_a_6c_b2c_c



Prototype	:	$ClH_{12}N_5Zn_2$
AFLOW prototype label	:	AB12C5D2_oP40_18_a_6c_b2c_c
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP40
Space group number	:	18
Space group symbol	:	$P2_12_12$
AFLOW prototype command	:	`aflow --proto=AB12C5D2_oP40_18_a_6c_b2c_c --params=`$a, b/a, c/a, z_1, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}$

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$z_1\,\mathbf{a}_3$	=	$z_1 c\,\hat{\mathbf{z}}$	$(2a)$	Cl
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	$(2a)$	Cl
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(2b)$	N I

$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 - z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - z_2 c\,\hat{\mathbf{z}}$	(2*b*)	N I
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*c*)	H I
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*c*)	H I
$\mathbf{B}_7$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(4*c*)	H I
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(4*c*)	H I
$\mathbf{B}_9$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(4*c*)	H II
$\mathbf{B}_{10}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(4*c*)	H II
$\mathbf{B}_{11}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(4*c*)	H II
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(4*c*)	H II
$\mathbf{B}_{13}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(4*c*)	H III
$\mathbf{B}_{14}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(4*c*)	H III
$\mathbf{B}_{15}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right)b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(4*c*)	H III
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right)b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(4*c*)	H III
$\mathbf{B}_{17}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(4*c*)	H IV
$\mathbf{B}_{18}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(4*c*)	H IV
$\mathbf{B}_{19}$	=	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right)b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(4*c*)	H IV
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right)b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(4*c*)	H IV
$\mathbf{B}_{21}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(4*c*)	H V
$\mathbf{B}_{22}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(4*c*)	H V
$\mathbf{B}_{23}$	=	$\left(\frac{1}{2} - x_7\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right)b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	(4*c*)	H V
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} + x_7\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_7\right)\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_7\right)b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	(4*c*)	H V
$\mathbf{B}_{25}$	=	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(4*c*)	H VI
$\mathbf{B}_{26}$	=	$-x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(4*c*)	H VI
$\mathbf{B}_{27}$	=	$\left(\frac{1}{2} - x_8\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_8\right)\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_8\right)b\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(4*c*)	H VI
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} + x_8\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_8\right)\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_8\right)b\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(4*c*)	H VI
$\mathbf{B}_{29}$	=	$x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(4*c*)	N II
$\mathbf{B}_{30}$	=	$-x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(4*c*)	N II
$\mathbf{B}_{31}$	=	$\left(\frac{1}{2} - x_9\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_9\right)\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_9\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_9\right)b\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}}$	(4*c*)	N II
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2} + x_9\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_9\right)\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_9\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_9\right)b\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}}$	(4*c*)	N II
$\mathbf{B}_{33}$	=	$x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$x_{10} a\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}}$	(4*c*)	N III
$\mathbf{B}_{34}$	=	$-x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$-x_{10} a\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}}$	(4*c*)	N III
$\mathbf{B}_{35}$	=	$\left(\frac{1}{2} - x_{10}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{10}\right)\mathbf{a}_2 - z_{10}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{10}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{10}\right)b\,\hat{\mathbf{y}} - z_{10} c\,\hat{\mathbf{z}}$	(4*c*)	N III
$\mathbf{B}_{36}$	=	$\left(\frac{1}{2} + x_{10}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{10}\right)\mathbf{a}_2 - z_{10}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{10}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{10}\right)b\,\hat{\mathbf{y}} - z_{10} c\,\hat{\mathbf{z}}$	(4*c*)	N III
$\mathbf{B}_{37}$	=	$x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$x_{11} a\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + z_{11} c\,\hat{\mathbf{z}}$	(4*c*)	Zn
$\mathbf{B}_{38}$	=	$-x_{11}\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$-x_{11} a\,\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} + z_{11} c\,\hat{\mathbf{z}}$	(4*c*)	Zn
$\mathbf{B}_{39}$	=	$\left(\frac{1}{2} - x_{11}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{11}\right)\mathbf{a}_2 - z_{11}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{11}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{11}\right)b\,\hat{\mathbf{y}} - z_{11} c\,\hat{\mathbf{z}}$	(4*c*)	Zn

$$\mathbf{B}_{40} = \left(\tfrac{1}{2} + x_{11}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{11}\right)\mathbf{a}_2 - z_{11}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{11}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{11}\right)b\,\hat{\mathbf{y}} - z_{11}c\,\hat{\mathbf{z}} \quad (4c) \quad \text{Zn}$$

References:

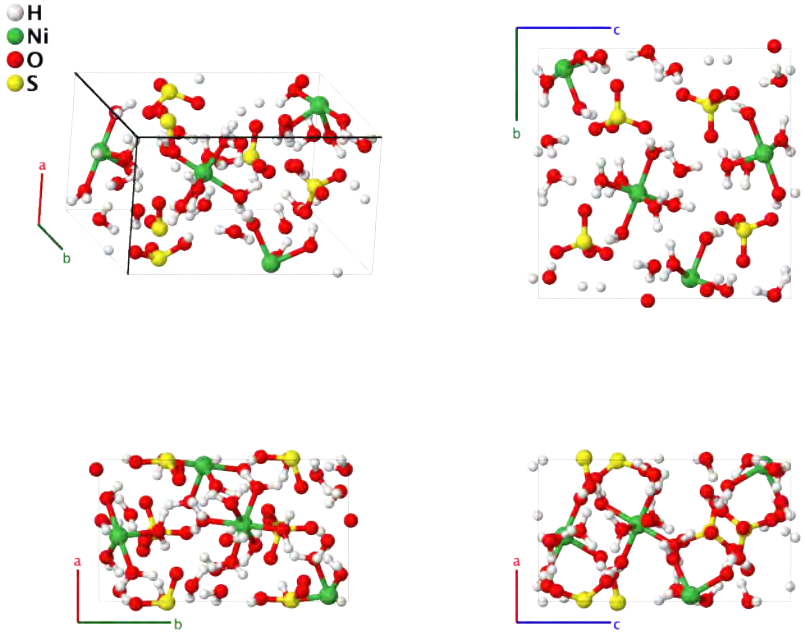
- T. M. M. Richter, S. Strobel, N. S. A. Alt, E. Schlücker, and R. Niewa, *Ammonothermal Synthesis and Crystal Structures of Diamminetriamidodizinc Chloride [$Zn_2(NH_3)_2(NH_2)_3$]Cl and Diamminemonoamidozinc Bromide [$Zn(NH_3)_2(NH_2)$]Br*, Inorganics **4**, 41 (2016), doi:10.3390/inorganics4040041.

Geometry files:

- CIF: pp. 1574
- POSCAR: pp. 1575

Morenosite ($NiSO_4 \cdot 7H_2O$, $H4_{12}$) Structure: A14BC11D_oP108_19_14a_a_11a_a

http://aflow.org/prototype-encyclopedia/A14BC11D_oP108_19_14a_a_11a_a



Prototype	:	$H_{14}NiO_{11}S$
AFLOW prototype label	:	A14BC11D_oP108_19_14a_a_11a_a
***Strukturbericht* designation**	:	$H4_{12}$
Pearson symbol	:	oP108
Space group number	:	19
Space group symbol	:	$P2_12_12_1$
AFLOW prototype command	:	`aflow --proto=A14BC11D_oP108_19_14a_a_11a_a --params=`$a, b/a, c/a, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}, x_{17}, y_{17}, z_{17}, x_{18}, y_{18}, z_{18}, x_{19}, y_{19}, z_{19}, x_{20}, y_{20}, z_{20}, x_{21}, y_{21}, z_{21}, x_{22}, y_{22}, z_{22}, x_{23}, y_{23}, z_{23}, x_{24}, y_{24}, z_{24}, x_{25}, y_{25}, z_{25}, x_{26}, y_{26}, z_{26}, x_{27}, y_{27}, z_{27}$

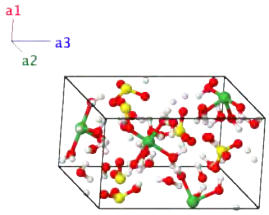
Other compounds with this structure

- $MgSO_4 \cdot 7H_2O$ (epsomite) and $ZnSO_4 \cdot 7H_2O$ (gosalrite)

- We use the structure from the 25 K neutron data taken by (Ptasiewicz-Bak, 1997).

Simple Orthorhombic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= a\,\hat{\mathbf{x}}\\ \mathbf{a}_2 &= b\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	$(4a)$	H I
$\mathbf{B}_2$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right)a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\,\hat{\mathbf{z}}$	$(4a)$	H I
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2} + y_1\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right)c\,\hat{\mathbf{z}}$	$(4a)$	H I
$\mathbf{B}_4$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_1\right)\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_1\right)b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	$(4a)$	H I
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(4a)$	H II
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	$(4a)$	H II
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	$(4a)$	H II
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right)b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	$(4a)$	H II
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(4a)$	H III
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	$(4a)$	H III
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	$(4a)$	H III
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(4a)$	H III
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(4a)$	H IV
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	$(4a)$	H IV
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	$(4a)$	H IV
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	$(4a)$	H IV
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(4a)$	H V
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	$(4a)$	H V
$\mathbf{B}_{19}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	$(4a)$	H V
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right)b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	$(4a)$	H V
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(4a)$	H VI
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\,\hat{\mathbf{z}}$	$(4a)$	H VI
$\mathbf{B}_{23}$	=	$-x_6\,\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\,\hat{\mathbf{z}}$	$(4a)$	H VI
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right)b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	$(4a)$	H VI
$\mathbf{B}_{25}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	$(4a)$	H VII
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} - x_7\right)\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_7\right)a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right)c\,\hat{\mathbf{z}}$	$(4a)$	H VII
$\mathbf{B}_{27}$	=	$-x_7\,\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right)c\,\hat{\mathbf{z}}$	$(4a)$	H VII

$\mathbf{B}_{28}$	$\left(\frac{1}{2}+x_7\right)\mathbf{a}_1+\left(\frac{1}{2}-y_7\right)\mathbf{a}_2-z_7\,\mathbf{a}_3$	$=$	$\left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_7\right)b\,\hat{\mathbf{y}}-z_7c\,\hat{\mathbf{z}}$	(4*a*)	H VII
$\mathbf{B}_{29}$	$x_8\,\mathbf{a}_1+y_8\,\mathbf{a}_2+z_8\,\mathbf{a}_3$	$=$	$x_8a\,\hat{\mathbf{x}}+y_8b\,\hat{\mathbf{y}}+z_8c\,\hat{\mathbf{z}}$	(4*a*)	H VIII
$\mathbf{B}_{30}$	$\left(\frac{1}{2}-x_8\right)\mathbf{a}_1-y_8\,\mathbf{a}_2+\left(\frac{1}{2}+z_8\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}-x_8\right)a\,\hat{\mathbf{x}}-y_8b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_8\right)c\,\hat{\mathbf{z}}$	(4*a*)	H VIII
$\mathbf{B}_{31}$	$-x_8\,\mathbf{a}_1+\left(\frac{1}{2}+y_8\right)\mathbf{a}_2+\left(\frac{1}{2}-z_8\right)\mathbf{a}_3$	$=$	$-x_8a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_8\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_8\right)c\,\hat{\mathbf{z}}$	(4*a*)	H VIII
$\mathbf{B}_{32}$	$\left(\frac{1}{2}+x_8\right)\mathbf{a}_1+\left(\frac{1}{2}-y_8\right)\mathbf{a}_2-z_8\,\mathbf{a}_3$	$=$	$\left(\frac{1}{2}+x_8\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_8\right)b\,\hat{\mathbf{y}}-z_8c\,\hat{\mathbf{z}}$	(4*a*)	H VIII
$\mathbf{B}_{33}$	$x_9\,\mathbf{a}_1+y_9\,\mathbf{a}_2+z_9\,\mathbf{a}_3$	$=$	$x_9a\,\hat{\mathbf{x}}+y_9b\,\hat{\mathbf{y}}+z_9c\,\hat{\mathbf{z}}$	(4*a*)	H IX
$\mathbf{B}_{34}$	$\left(\frac{1}{2}-x_9\right)\mathbf{a}_1-y_9\,\mathbf{a}_2+\left(\frac{1}{2}+z_9\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}-x_9\right)a\,\hat{\mathbf{x}}-y_9b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_9\right)c\,\hat{\mathbf{z}}$	(4*a*)	H IX
$\mathbf{B}_{35}$	$-x_9\,\mathbf{a}_1+\left(\frac{1}{2}+y_9\right)\mathbf{a}_2+\left(\frac{1}{2}-z_9\right)\mathbf{a}_3$	$=$	$-x_9a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_9\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_9\right)c\,\hat{\mathbf{z}}$	(4*a*)	H IX
$\mathbf{B}_{36}$	$\left(\frac{1}{2}+x_9\right)\mathbf{a}_1+\left(\frac{1}{2}-y_9\right)\mathbf{a}_2-z_9\,\mathbf{a}_3$	$=$	$\left(\frac{1}{2}+x_9\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_9\right)b\,\hat{\mathbf{y}}-z_9c\,\hat{\mathbf{z}}$	(4*a*)	H IX
$\mathbf{B}_{37}$	$x_{10}\,\mathbf{a}_1+y_{10}\,\mathbf{a}_2+z_{10}\,\mathbf{a}_3$	$=$	$x_{10}a\,\hat{\mathbf{x}}+y_{10}b\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}}$	(4*a*)	H X
$\mathbf{B}_{38}$	$\left(\frac{1}{2}-x_{10}\right)\mathbf{a}_1-y_{10}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{10}\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{x}}-y_{10}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{10}\right)c\,\hat{\mathbf{z}}$	(4*a*)	H X
$\mathbf{B}_{39}$	$-x_{10}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{10}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{10}\right)\mathbf{a}_3$	$=$	$-x_{10}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{10}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{10}\right)c\,\hat{\mathbf{z}}$	(4*a*)	H X
$\mathbf{B}_{40}$	$\left(\frac{1}{2}+x_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{10}\right)\mathbf{a}_2-z_{10}\,\mathbf{a}_3$	$=$	$\left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{10}\right)b\,\hat{\mathbf{y}}-z_{10}c\,\hat{\mathbf{z}}$	(4*a*)	H X
$\mathbf{B}_{41}$	$x_{11}\,\mathbf{a}_1+y_{11}\,\mathbf{a}_2+z_{11}\,\mathbf{a}_3$	$=$	$x_{11}a\,\hat{\mathbf{x}}+y_{11}b\,\hat{\mathbf{y}}+z_{11}c\,\hat{\mathbf{z}}$	(4*a*)	H XI
$\mathbf{B}_{42}$	$\left(\frac{1}{2}-x_{11}\right)\mathbf{a}_1-y_{11}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{11}\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}-x_{11}\right)a\,\hat{\mathbf{x}}-y_{11}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{11}\right)c\,\hat{\mathbf{z}}$	(4*a*)	H XI
$\mathbf{B}_{43}$	$-x_{11}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{11}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{11}\right)\mathbf{a}_3$	$=$	$-x_{11}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{11}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{11}\right)c\,\hat{\mathbf{z}}$	(4*a*)	H XI
$\mathbf{B}_{44}$	$\left(\frac{1}{2}+x_{11}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{11}\right)\mathbf{a}_2-z_{11}\,\mathbf{a}_3$	$=$	$\left(\frac{1}{2}+x_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{11}\right)b\,\hat{\mathbf{y}}-z_{11}c\,\hat{\mathbf{z}}$	(4*a*)	H XI
$\mathbf{B}_{45}$	$x_{12}\,\mathbf{a}_1+y_{12}\,\mathbf{a}_2+z_{12}\,\mathbf{a}_3$	$=$	$x_{12}a\,\hat{\mathbf{x}}+y_{12}b\,\hat{\mathbf{y}}+z_{12}c\,\hat{\mathbf{z}}$	(4*a*)	H XII
$\mathbf{B}_{46}$	$\left(\frac{1}{2}-x_{12}\right)\mathbf{a}_1-y_{12}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{12}\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}-x_{12}\right)a\,\hat{\mathbf{x}}-y_{12}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{12}\right)c\,\hat{\mathbf{z}}$	(4*a*)	H XII
$\mathbf{B}_{47}$	$-x_{12}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{12}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{12}\right)\mathbf{a}_3$	$=$	$-x_{12}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{12}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{12}\right)c\,\hat{\mathbf{z}}$	(4*a*)	H XII
$\mathbf{B}_{48}$	$\left(\frac{1}{2}+x_{12}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{12}\right)\mathbf{a}_2-z_{12}\,\mathbf{a}_3$	$=$	$\left(\frac{1}{2}+x_{12}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{12}\right)b\,\hat{\mathbf{y}}-z_{12}c\,\hat{\mathbf{z}}$	(4*a*)	H XII
$\mathbf{B}_{49}$	$x_{13}\,\mathbf{a}_1+y_{13}\,\mathbf{a}_2+z_{13}\,\mathbf{a}_3$	$=$	$x_{13}a\,\hat{\mathbf{x}}+y_{13}b\,\hat{\mathbf{y}}+z_{13}c\,\hat{\mathbf{z}}$	(4*a*)	H XIII
$\mathbf{B}_{50}$	$\left(\frac{1}{2}-x_{13}\right)\mathbf{a}_1-y_{13}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{13}\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}-x_{13}\right)a\,\hat{\mathbf{x}}-y_{13}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{13}\right)c\,\hat{\mathbf{z}}$	(4*a*)	H XIII
$\mathbf{B}_{51}$	$-x_{13}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{13}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{13}\right)\mathbf{a}_3$	$=$	$-x_{13}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{13}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{13}\right)c\,\hat{\mathbf{z}}$	(4*a*)	H XIII
$\mathbf{B}_{52}$	$\left(\frac{1}{2}+x_{13}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{13}\right)\mathbf{a}_2-z_{13}\,\mathbf{a}_3$	$=$	$\left(\frac{1}{2}+x_{13}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{13}\right)b\,\hat{\mathbf{y}}-z_{13}c\,\hat{\mathbf{z}}$	(4*a*)	H XIII
$\mathbf{B}_{53}$	$x_{14}\,\mathbf{a}_1+y_{14}\,\mathbf{a}_2+z_{14}\,\mathbf{a}_3$	$=$	$x_{14}a\,\hat{\mathbf{x}}+y_{14}b\,\hat{\mathbf{y}}+z_{14}c\,\hat{\mathbf{z}}$	(4*a*)	H XIV
$\mathbf{B}_{54}$	$\left(\frac{1}{2}-x_{14}\right)\mathbf{a}_1-y_{14}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{14}\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}-x_{14}\right)a\,\hat{\mathbf{x}}-y_{14}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{14}\right)c\,\hat{\mathbf{z}}$	(4*a*)	H XIV
$\mathbf{B}_{55}$	$-x_{14}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{14}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{14}\right)\mathbf{a}_3$	$=$	$-x_{14}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{14}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{14}\right)c\,\hat{\mathbf{z}}$	(4*a*)	H XIV
$\mathbf{B}_{56}$	$\left(\frac{1}{2}+x_{14}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{14}\right)\mathbf{a}_2-z_{14}\,\mathbf{a}_3$	$=$	$\left(\frac{1}{2}+x_{14}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{14}\right)b\,\hat{\mathbf{y}}-z_{14}c\,\hat{\mathbf{z}}$	(4*a*)	H XIV
$\mathbf{B}_{57}$	$x_{15}\,\mathbf{a}_1+y_{15}\,\mathbf{a}_2+z_{15}\,\mathbf{a}_3$	$=$	$x_{15}a\,\hat{\mathbf{x}}+y_{15}b\,\hat{\mathbf{y}}+z_{15}c\,\hat{\mathbf{z}}$	(4*a*)	Ni
$\mathbf{B}_{58}$	$\left(\frac{1}{2}-x_{15}\right)\mathbf{a}_1-y_{15}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{15}\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}-x_{15}\right)a\,\hat{\mathbf{x}}-y_{15}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{15}\right)c\,\hat{\mathbf{z}}$	(4*a*)	Ni
$\mathbf{B}_{59}$	$-x_{15}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{15}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{15}\right)\mathbf{a}_3$	$=$	$-x_{15}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{15}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{15}\right)c\,\hat{\mathbf{z}}$	(4*a*)	Ni

$$
\begin{aligned}
\mathbf{B}_{60} &= \left(\tfrac{1}{2}+x_{15}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{15}\right)\mathbf{a}_2-z_{15}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_{15}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{15}\right)b\,\hat{\mathbf{y}}-z_{15}c\,\hat{\mathbf{z}} && (4a) && \text{Ni}\\
\mathbf{B}_{61} &= x_{16}\,\mathbf{a}_1+y_{16}\,\mathbf{a}_2+z_{16}\,\mathbf{a}_3 &&= x_{16}a\,\hat{\mathbf{x}}+y_{16}b\,\hat{\mathbf{y}}+z_{16}c\,\hat{\mathbf{z}} && (4a) && \text{O I}\\
\mathbf{B}_{62} &= \left(\tfrac{1}{2}-x_{16}\right)\mathbf{a}_1-y_{16}\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_{16}\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_{16}\right)a\,\hat{\mathbf{x}}-y_{16}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{16}\right)c\,\hat{\mathbf{z}} && (4a) && \text{O I}\\
\mathbf{B}_{63} &= -x_{16}\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_{16}\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_{16}\right)\mathbf{a}_3 &&= -x_{16}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{16}\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{16}\right)c\,\hat{\mathbf{z}} && (4a) && \text{O I}\\
\mathbf{B}_{64} &= \left(\tfrac{1}{2}+x_{16}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{16}\right)\mathbf{a}_2-z_{16}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_{16}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{16}\right)b\,\hat{\mathbf{y}}-z_{16}c\,\hat{\mathbf{z}} && (4a) && \text{O I}\\
\mathbf{B}_{65} &= x_{17}\,\mathbf{a}_1+y_{17}\,\mathbf{a}_2+z_{17}\,\mathbf{a}_3 &&= x_{17}a\,\hat{\mathbf{x}}+y_{17}b\,\hat{\mathbf{y}}+z_{17}c\,\hat{\mathbf{z}} && (4a) && \text{O II}\\
\mathbf{B}_{66} &= \left(\tfrac{1}{2}-x_{17}\right)\mathbf{a}_1-y_{17}\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_{17}\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_{17}\right)a\,\hat{\mathbf{x}}-y_{17}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{17}\right)c\,\hat{\mathbf{z}} && (4a) && \text{O II}\\
\mathbf{B}_{67} &= -x_{17}\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_{17}\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_{17}\right)\mathbf{a}_3 &&= -x_{17}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{17}\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{17}\right)c\,\hat{\mathbf{z}} && (4a) && \text{O II}\\
\mathbf{B}_{68} &= \left(\tfrac{1}{2}+x_{17}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{17}\right)\mathbf{a}_2-z_{17}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_{17}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{17}\right)b\,\hat{\mathbf{y}}-z_{17}c\,\hat{\mathbf{z}} && (4a) && \text{O II}\\
\mathbf{B}_{69} &= x_{18}\,\mathbf{a}_1+y_{18}\,\mathbf{a}_2+z_{18}\,\mathbf{a}_3 &&= x_{18}a\,\hat{\mathbf{x}}+y_{18}b\,\hat{\mathbf{y}}+z_{18}c\,\hat{\mathbf{z}} && (4a) && \text{O III}\\
\mathbf{B}_{70} &= \left(\tfrac{1}{2}-x_{18}\right)\mathbf{a}_1-y_{18}\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_{18}\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_{18}\right)a\,\hat{\mathbf{x}}-y_{18}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{18}\right)c\,\hat{\mathbf{z}} && (4a) && \text{O III}\\
\mathbf{B}_{71} &= -x_{18}\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_{18}\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_{18}\right)\mathbf{a}_3 &&= -x_{18}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{18}\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{18}\right)c\,\hat{\mathbf{z}} && (4a) && \text{O III}\\
\mathbf{B}_{72} &= \left(\tfrac{1}{2}+x_{18}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{18}\right)\mathbf{a}_2-z_{18}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_{18}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{18}\right)b\,\hat{\mathbf{y}}-z_{18}c\,\hat{\mathbf{z}} && (4a) && \text{O III}\\
\mathbf{B}_{73} &= x_{19}\,\mathbf{a}_1+y_{19}\,\mathbf{a}_2+z_{19}\,\mathbf{a}_3 &&= x_{19}a\,\hat{\mathbf{x}}+y_{19}b\,\hat{\mathbf{y}}+z_{19}c\,\hat{\mathbf{z}} && (4a) && \text{O IV}\\
\mathbf{B}_{74} &= \left(\tfrac{1}{2}-x_{19}\right)\mathbf{a}_1-y_{19}\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_{19}\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_{19}\right)a\,\hat{\mathbf{x}}-y_{19}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{19}\right)c\,\hat{\mathbf{z}} && (4a) && \text{O IV}\\
\mathbf{B}_{75} &= -x_{19}\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_{19}\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_{19}\right)\mathbf{a}_3 &&= -x_{19}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{19}\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{19}\right)c\,\hat{\mathbf{z}} && (4a) && \text{O IV}\\
\mathbf{B}_{76} &= \left(\tfrac{1}{2}+x_{19}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{19}\right)\mathbf{a}_2-z_{19}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_{19}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{19}\right)b\,\hat{\mathbf{y}}-z_{19}c\,\hat{\mathbf{z}} && (4a) && \text{O IV}\\
\mathbf{B}_{77} &= x_{20}\,\mathbf{a}_1+y_{20}\,\mathbf{a}_2+z_{20}\,\mathbf{a}_3 &&= x_{20}a\,\hat{\mathbf{x}}+y_{20}b\,\hat{\mathbf{y}}+z_{20}c\,\hat{\mathbf{z}} && (4a) && \text{O V}\\
\mathbf{B}_{78} &= \left(\tfrac{1}{2}-x_{20}\right)\mathbf{a}_1-y_{20}\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_{20}\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_{20}\right)a\,\hat{\mathbf{x}}-y_{20}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{20}\right)c\,\hat{\mathbf{z}} && (4a) && \text{O V}\\
\mathbf{B}_{79} &= -x_{20}\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_{20}\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_{20}\right)\mathbf{a}_3 &&= -x_{20}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{20}\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{20}\right)c\,\hat{\mathbf{z}} && (4a) && \text{O V}\\
\mathbf{B}_{80} &= \left(\tfrac{1}{2}+x_{20}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{20}\right)\mathbf{a}_2-z_{20}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_{20}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{20}\right)b\,\hat{\mathbf{y}}-z_{20}c\,\hat{\mathbf{z}} && (4a) && \text{O V}\\
\mathbf{B}_{81} &= x_{21}\,\mathbf{a}_1+y_{21}\,\mathbf{a}_2+z_{21}\,\mathbf{a}_3 &&= x_{21}a\,\hat{\mathbf{x}}+y_{21}b\,\hat{\mathbf{y}}+z_{21}c\,\hat{\mathbf{z}} && (4a) && \text{O VI}\\
\mathbf{B}_{82} &= \left(\tfrac{1}{2}-x_{21}\right)\mathbf{a}_1-y_{21}\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_{21}\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_{21}\right)a\,\hat{\mathbf{x}}-y_{21}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{21}\right)c\,\hat{\mathbf{z}} && (4a) && \text{O VI}\\
\mathbf{B}_{83} &= -x_{21}\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_{21}\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_{21}\right)\mathbf{a}_3 &&= -x_{21}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{21}\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{21}\right)c\,\hat{\mathbf{z}} && (4a) && \text{O VI}\\
\mathbf{B}_{84} &= \left(\tfrac{1}{2}+x_{21}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{21}\right)\mathbf{a}_2-z_{21}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_{21}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{21}\right)b\,\hat{\mathbf{y}}-z_{21}c\,\hat{\mathbf{z}} && (4a) && \text{O VI}\\
\mathbf{B}_{85} &= x_{22}\,\mathbf{a}_1+y_{22}\,\mathbf{a}_2+z_{22}\,\mathbf{a}_3 &&= x_{22}a\,\hat{\mathbf{x}}+y_{22}b\,\hat{\mathbf{y}}+z_{22}c\,\hat{\mathbf{z}} && (4a) && \text{O VII}\\
\mathbf{B}_{86} &= \left(\tfrac{1}{2}-x_{22}\right)\mathbf{a}_1-y_{22}\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_{22}\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_{22}\right)a\,\hat{\mathbf{x}}-y_{22}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{22}\right)c\,\hat{\mathbf{z}} && (4a) && \text{O VII}\\
\mathbf{B}_{87} &= -x_{22}\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_{22}\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_{22}\right)\mathbf{a}_3 &&= -x_{22}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{22}\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{22}\right)c\,\hat{\mathbf{z}} && (4a) && \text{O VII}\\
\mathbf{B}_{88} &= \left(\tfrac{1}{2}+x_{22}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{22}\right)\mathbf{a}_2-z_{22}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_{22}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{22}\right)b\,\hat{\mathbf{y}}-z_{22}c\,\hat{\mathbf{z}} && (4a) && \text{O VII}\\
\mathbf{B}_{89} &= x_{23}\,\mathbf{a}_1+y_{23}\,\mathbf{a}_2+z_{23}\,\mathbf{a}_3 &&= x_{23}a\,\hat{\mathbf{x}}+y_{23}b\,\hat{\mathbf{y}}+z_{23}c\,\hat{\mathbf{z}} && (4a) && \text{O VIII}\\
\mathbf{B}_{90} &= \left(\tfrac{1}{2}-x_{23}\right)\mathbf{a}_1-y_{23}\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_{23}\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_{23}\right)a\,\hat{\mathbf{x}}-y_{23}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{23}\right)c\,\hat{\mathbf{z}} && (4a) && \text{O VIII}
\end{aligned}
$$

$\mathbf{B}_{91}$	=	$-x_{23}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{23}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{23}\right)\mathbf{a}_3$	=	$-x_{23}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{23}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{23}\right)c\,\hat{\mathbf{z}}$	(4*a*)	O VIII
$\mathbf{B}_{92}$	=	$\left(\frac{1}{2}+x_{23}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{23}\right)\mathbf{a}_2-z_{23}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{23}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{23}\right)b\,\hat{\mathbf{y}}-z_{23}c\,\hat{\mathbf{z}}$	(4*a*)	O VIII
$\mathbf{B}_{93}$	=	$x_{24}\,\mathbf{a}_1+y_{24}\,\mathbf{a}_2+z_{24}\,\mathbf{a}_3$	=	$x_{24}a\,\hat{\mathbf{x}}+y_{24}b\,\hat{\mathbf{y}}+z_{24}c\,\hat{\mathbf{z}}$	(4*a*)	O IX
$\mathbf{B}_{94}$	=	$\left(\frac{1}{2}-x_{24}\right)\mathbf{a}_1-y_{24}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{24}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{24}\right)a\,\hat{\mathbf{x}}-y_{24}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{24}\right)c\,\hat{\mathbf{z}}$	(4*a*)	O IX
$\mathbf{B}_{95}$	=	$-x_{24}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{24}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{24}\right)\mathbf{a}_3$	=	$-x_{24}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{24}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{24}\right)c\,\hat{\mathbf{z}}$	(4*a*)	O IX
$\mathbf{B}_{96}$	=	$\left(\frac{1}{2}+x_{24}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{24}\right)\mathbf{a}_2-z_{24}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{24}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{24}\right)b\,\hat{\mathbf{y}}-z_{24}c\,\hat{\mathbf{z}}$	(4*a*)	O IX
$\mathbf{B}_{97}$	=	$x_{25}\,\mathbf{a}_1+y_{25}\,\mathbf{a}_2+z_{25}\,\mathbf{a}_3$	=	$x_{25}a\,\hat{\mathbf{x}}+y_{25}b\,\hat{\mathbf{y}}+z_{25}c\,\hat{\mathbf{z}}$	(4*a*)	O X
$\mathbf{B}_{98}$	=	$\left(\frac{1}{2}-x_{25}\right)\mathbf{a}_1-y_{25}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{25}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{25}\right)a\,\hat{\mathbf{x}}-y_{25}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{25}\right)c\,\hat{\mathbf{z}}$	(4*a*)	O X
$\mathbf{B}_{99}$	=	$-x_{25}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{25}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{25}\right)\mathbf{a}_3$	=	$-x_{25}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{25}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{25}\right)c\,\hat{\mathbf{z}}$	(4*a*)	O X
$\mathbf{B}_{100}$	=	$\left(\frac{1}{2}+x_{25}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{25}\right)\mathbf{a}_2-z_{25}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{25}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{25}\right)b\,\hat{\mathbf{y}}-z_{25}c\,\hat{\mathbf{z}}$	(4*a*)	O X
$\mathbf{B}_{101}$	=	$x_{26}\,\mathbf{a}_1+y_{26}\,\mathbf{a}_2+z_{26}\,\mathbf{a}_3$	=	$x_{26}a\,\hat{\mathbf{x}}+y_{26}b\,\hat{\mathbf{y}}+z_{26}c\,\hat{\mathbf{z}}$	(4*a*)	O XI
$\mathbf{B}_{102}$	=	$\left(\frac{1}{2}-x_{26}\right)\mathbf{a}_1-y_{26}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{26}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{26}\right)a\,\hat{\mathbf{x}}-y_{26}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{26}\right)c\,\hat{\mathbf{z}}$	(4*a*)	O XI
$\mathbf{B}_{103}$	=	$-x_{26}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{26}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{26}\right)\mathbf{a}_3$	=	$-x_{26}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{26}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{26}\right)c\,\hat{\mathbf{z}}$	(4*a*)	O XI
$\mathbf{B}_{104}$	=	$\left(\frac{1}{2}+x_{26}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{26}\right)\mathbf{a}_2-z_{26}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{26}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{26}\right)b\,\hat{\mathbf{y}}-z_{26}c\,\hat{\mathbf{z}}$	(4*a*)	O XI
$\mathbf{B}_{105}$	=	$x_{27}\,\mathbf{a}_1+y_{27}\,\mathbf{a}_2+z_{27}\,\mathbf{a}_3$	=	$x_{27}a\,\hat{\mathbf{x}}+y_{27}b\,\hat{\mathbf{y}}+z_{27}c\,\hat{\mathbf{z}}$	(4*a*)	S
$\mathbf{B}_{106}$	=	$\left(\frac{1}{2}-x_{27}\right)\mathbf{a}_1-y_{27}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{27}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{27}\right)a\,\hat{\mathbf{x}}-y_{27}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{27}\right)c\,\hat{\mathbf{z}}$	(4*a*)	S
$\mathbf{B}_{107}$	=	$-x_{27}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{27}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{27}\right)\mathbf{a}_3$	=	$-x_{27}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{27}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{27}\right)c\,\hat{\mathbf{z}}$	(4*a*)	S
$\mathbf{B}_{108}$	=	$\left(\frac{1}{2}+x_{27}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{27}\right)\mathbf{a}_2-z_{27}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{27}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{27}\right)b\,\hat{\mathbf{y}}-z_{27}c\,\hat{\mathbf{z}}$	(4*a*)	S

References:

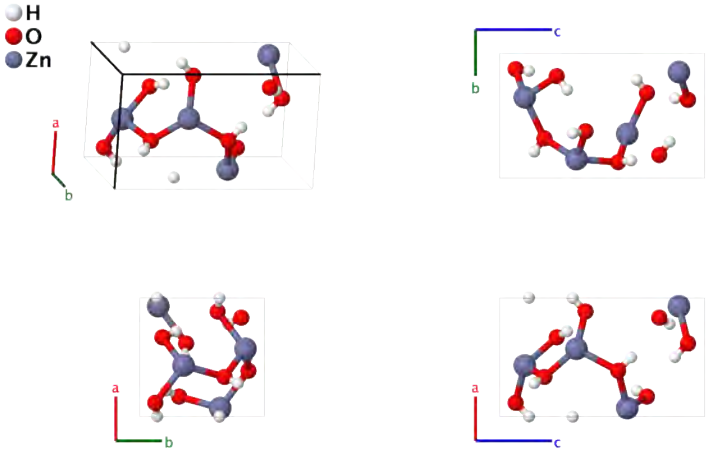
- H. Ptasiewicz-Bak, I. Olovsson, and G. J. McIntyre, *Charge Density in Orthorhombic $NiSO_4{\cdot}7H_2O$ at Room Temperature and 25 K*, Acta Crystallogr. Sect. B Struct. Sci. **53**, 325–336 (1997), doi:10.1107/S0108768196014061.

Geometry files:

- CIF: pp. 1575
- POSCAR: pp. 1576

Wülfingite (ϵ-$Zn(OH)_2$, $C31$) Structure: A2B2C_oP20_19_2a_2a_a

http://aflow.org/prototype-encyclopedia/A2B2C_oP20_19_2a_2a_a



Prototype	:	H_2O_2Zn
AFLOW prototype label	:	A2B2C_oP20_19_2a_2a_a
***Strukturbericht* designation**	:	$C31$
Pearson symbol	:	oP20
Space group number	:	19
Space group symbol	:	$P2_12_12_1$
AFLOW prototype command	:	`aflow --proto=A2B2C_oP20_19_2a_2a_a --params=`$a, b/a, c/a, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5$

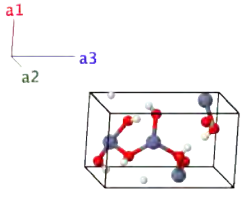
Other compounds with this structure

- β-$Be(OH)_2$

- (Corey, 1933) originally determined the structure of ϵ-$Zn(OH)_2$, but they were unable to locate the hydrogen atoms. This structure was given the *Strukturbericht* designation $C31$ by (Gottfried, 1937). (Stahl, 1998) located the hydrogen atoms, and found that they did not change the space group, so we use the updated structure as our $C31$ prototype.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	($4a$)	H I

$\mathbf{B}_2$	$=$	$\left(\frac{1}{2}-x_1\right)\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \left(\frac{1}{2}+z_1\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}-x_1\right)a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_1\right)c\,\hat{\mathbf{z}}$	(4*a*)	H I
$\mathbf{B}_3$	$=$	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2}+y_1\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_1\right)\mathbf{a}_3$	$=$	$-x_1 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_1\right)c\,\hat{\mathbf{z}}$	(4*a*)	H I
$\mathbf{B}_4$	$=$	$\left(\frac{1}{2}+x_1\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_1\right)\mathbf{a}_2 - z_1\,\mathbf{a}_3$	$=$	$\left(\frac{1}{2}+x_1\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_1\right)b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	(4*a*)	H I
$\mathbf{B}_5$	$=$	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	$=$	$x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*a*)	H II
$\mathbf{B}_6$	$=$	$\left(\frac{1}{2}-x_2\right)\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}-x_2\right)a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_2\right)c\,\hat{\mathbf{z}}$	(4*a*)	H II
$\mathbf{B}_7$	$=$	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2}+y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_2\right)\mathbf{a}_3$	$=$	$-x_2 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_2\right)c\,\hat{\mathbf{z}}$	(4*a*)	H II
$\mathbf{B}_8$	$=$	$\left(\frac{1}{2}+x_2\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_2\right)\mathbf{a}_2 - z_2\,\mathbf{a}_3$	$=$	$\left(\frac{1}{2}+x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_2\right)b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(4*a*)	H II
$\mathbf{B}_9$	$=$	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$=$	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_{10}$	$=$	$\left(\frac{1}{2}-x_3\right)\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_3\right)c\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_{11}$	$=$	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2}+y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_3\right)\mathbf{a}_3$	$=$	$-x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_3\right)c\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_{12}$	$=$	$\left(\frac{1}{2}+x_3\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_3\right)\mathbf{a}_2 - z_3\,\mathbf{a}_3$	$=$	$\left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_3\right)b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_{13}$	$=$	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$=$	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(4*a*)	O II
$\mathbf{B}_{14}$	$=$	$\left(\frac{1}{2}-x_4\right)\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_4\right)c\,\hat{\mathbf{z}}$	(4*a*)	O II
$\mathbf{B}_{15}$	$=$	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2}+y_4\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_4\right)\mathbf{a}_3$	$=$	$-x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_4\right)c\,\hat{\mathbf{z}}$	(4*a*)	O II
$\mathbf{B}_{16}$	$=$	$\left(\frac{1}{2}+x_4\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3$	$=$	$\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_4\right)b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(4*a*)	O II
$\mathbf{B}_{17}$	$=$	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	$=$	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(4*a*)	Zn
$\mathbf{B}_{18}$	$=$	$\left(\frac{1}{2}-x_5\right)\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	(4*a*)	Zn
$\mathbf{B}_{19}$	$=$	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2}+y_5\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_5\right)\mathbf{a}_3$	$=$	$-x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_5\right)c\,\hat{\mathbf{z}}$	(4*a*)	Zn
$\mathbf{B}_{20}$	$=$	$\left(\frac{1}{2}+x_5\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3$	$=$	$\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_5\right)b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(4*a*)	Zn

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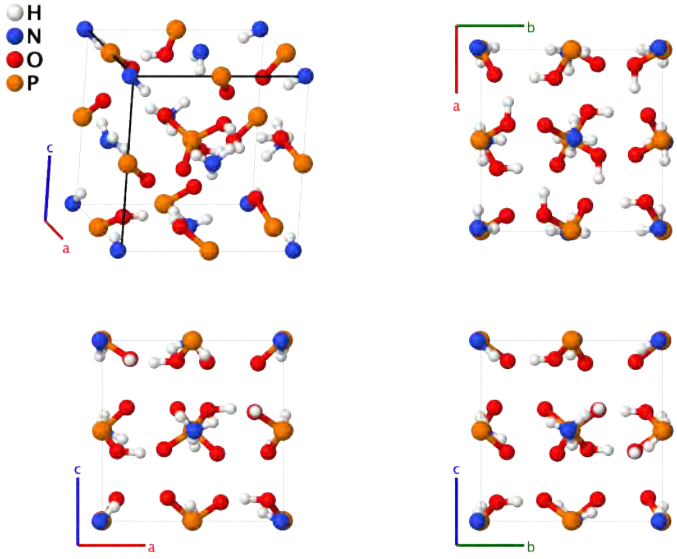
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1576
- POSCAR: pp. 1577

Ferroelectric $NH_4H_2PO_4$ Structure: A6BC4D_oP48_19_6a_a_4a_a

http://aflow.org/prototype-encyclopedia/A6BC4D_oP48_19_6a_a_4a_a



Prototype	:	H_6NO_4P
AFLOW prototype label	:	A6BC4D_oP48_19_6a_a_4a_a
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP48
Space group number	:	19
Space group symbol	:	$P2_12_12_1$
AFLOW prototype command	:	`aflow --proto=A6BC4D_oP48_19_6a_a_4a_a --params=`$a, b/a, c/a, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}$

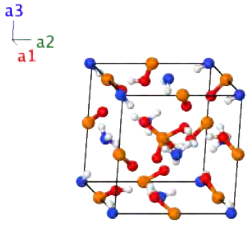
Other compounds with this structure

- $NH_4H_2AsO_4$

- Below 148 K, $NH_4H_2PO_4$ undergoes a transition from a tetragonal paraelectric phase to this orthorhombic ferroelectric phase. This seems to occur because the hydrogen atoms associated with the PO_4 ions become frozen into place.
- The data for this structure was taken at 143 K.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(4*a*)	H I
$\mathbf{B}_2$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right)a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\,\hat{\mathbf{z}}$	(4*a*)	H I
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2} + y_1\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right)c\,\hat{\mathbf{z}}$	(4*a*)	H I
$\mathbf{B}_4$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_1\right)\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_1\right)b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	(4*a*)	H I
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*a*)	H II
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	(4*a*)	H II
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	(4*a*)	H II
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right)b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(4*a*)	H II
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*a*)	H III
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(4*a*)	H III
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(4*a*)	H III
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(4*a*)	H III
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(4*a*)	H IV
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(4*a*)	H IV
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(4*a*)	H IV
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(4*a*)	H IV
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(4*a*)	H V
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	(4*a*)	H V
$\mathbf{B}_{19}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	(4*a*)	H V
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right)b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(4*a*)	H V
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(4*a*)	H VI
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\,\hat{\mathbf{z}}$	(4*a*)	H VI
$\mathbf{B}_{23}$	=	$-x_6\,\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\,\hat{\mathbf{z}}$	(4*a*)	H VI
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right)b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(4*a*)	H VI
$\mathbf{B}_{25}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(4*a*)	N
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} - x_7\right)\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_7\right)a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right)c\,\hat{\mathbf{z}}$	(4*a*)	N
$\mathbf{B}_{27}$	=	$-x_7\,\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right)c\,\hat{\mathbf{z}}$	(4*a*)	N
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} + x_7\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_7\right)\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_7\right)b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	(4*a*)	N
$\mathbf{B}_{29}$	=	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2} - x_8\right)\mathbf{a}_1 - y_8\,\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_8\right)a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right)c\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_{31}$	=	$-x_8\,\mathbf{a}_1 + \left(\frac{1}{2} + y_8\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_8\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right)c\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2} + x_8\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_8\right)\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_8\right)b\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_{33}$	=	$x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(4*a*)	O II
$\mathbf{B}_{34}$	=	$\left(\frac{1}{2} - x_9\right)\mathbf{a}_1 - y_9\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_9\right)a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right)c\,\hat{\mathbf{z}}$	(4*a*)	O II

$\mathbf{B}_{35}$	=	$-x_9\,\mathbf{a}_1+\left(\frac{1}{2}+y_9\right)\mathbf{a}_2+\left(\frac{1}{2}-z_9\right)\mathbf{a}_3$	=	$-x_9a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_9\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_9\right)c\,\hat{\mathbf{z}}$	(4*a*)	O II
$\mathbf{B}_{36}$	=	$\left(\frac{1}{2}+x_9\right)\mathbf{a}_1+\left(\frac{1}{2}-y_9\right)\mathbf{a}_2-z_9\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_9\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_9\right)b\,\hat{\mathbf{y}}-z_9c\,\hat{\mathbf{z}}$	(4*a*)	O II
$\mathbf{B}_{37}$	=	$x_{10}\,\mathbf{a}_1+y_{10}\,\mathbf{a}_2+z_{10}\,\mathbf{a}_3$	=	$x_{10}a\,\hat{\mathbf{x}}+y_{10}b\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}}$	(4*a*)	O III
$\mathbf{B}_{38}$	=	$\left(\frac{1}{2}-x_{10}\right)\mathbf{a}_1-y_{10}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{10}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{x}}-y_{10}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{10}\right)c\,\hat{\mathbf{z}}$	(4*a*)	O III
$\mathbf{B}_{39}$	=	$-x_{10}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{10}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{10}\right)\mathbf{a}_3$	=	$-x_{10}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{10}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{10}\right)c\,\hat{\mathbf{z}}$	(4*a*)	O III
$\mathbf{B}_{40}$	=	$\left(\frac{1}{2}+x_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{10}\right)\mathbf{a}_2-z_{10}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{10}\right)b\,\hat{\mathbf{y}}-z_{10}c\,\hat{\mathbf{z}}$	(4*a*)	O III
$\mathbf{B}_{41}$	=	$x_{11}\,\mathbf{a}_1+y_{11}\,\mathbf{a}_2+z_{11}\,\mathbf{a}_3$	=	$x_{11}a\,\hat{\mathbf{x}}+y_{11}b\,\hat{\mathbf{y}}+z_{11}c\,\hat{\mathbf{z}}$	(4*a*)	O IV
$\mathbf{B}_{42}$	=	$\left(\frac{1}{2}-x_{11}\right)\mathbf{a}_1-y_{11}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{11}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{11}\right)a\,\hat{\mathbf{x}}-y_{11}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{11}\right)c\,\hat{\mathbf{z}}$	(4*a*)	O IV
$\mathbf{B}_{43}$	=	$-x_{11}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{11}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{11}\right)\mathbf{a}_3$	=	$-x_{11}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{11}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{11}\right)c\,\hat{\mathbf{z}}$	(4*a*)	O IV
$\mathbf{B}_{44}$	=	$\left(\frac{1}{2}+x_{11}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{11}\right)\mathbf{a}_2-z_{11}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{11}\right)b\,\hat{\mathbf{y}}-z_{11}c\,\hat{\mathbf{z}}$	(4*a*)	O IV
$\mathbf{B}_{45}$	=	$x_{12}\,\mathbf{a}_1+y_{12}\,\mathbf{a}_2+z_{12}\,\mathbf{a}_3$	=	$x_{12}a\,\hat{\mathbf{x}}+y_{12}b\,\hat{\mathbf{y}}+z_{12}c\,\hat{\mathbf{z}}$	(4*a*)	P
$\mathbf{B}_{46}$	=	$\left(\frac{1}{2}-x_{12}\right)\mathbf{a}_1-y_{12}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{12}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{12}\right)a\,\hat{\mathbf{x}}-y_{12}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{12}\right)c\,\hat{\mathbf{z}}$	(4*a*)	P
$\mathbf{B}_{47}$	=	$-x_{12}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{12}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{12}\right)\mathbf{a}_3$	=	$-x_{12}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{12}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{12}\right)c\,\hat{\mathbf{z}}$	(4*a*)	P
$\mathbf{B}_{48}$	=	$\left(\frac{1}{2}+x_{12}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{12}\right)\mathbf{a}_2-z_{12}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{12}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{12}\right)b\,\hat{\mathbf{y}}-z_{12}c\,\hat{\mathbf{z}}$	(4*a*)	P

References:

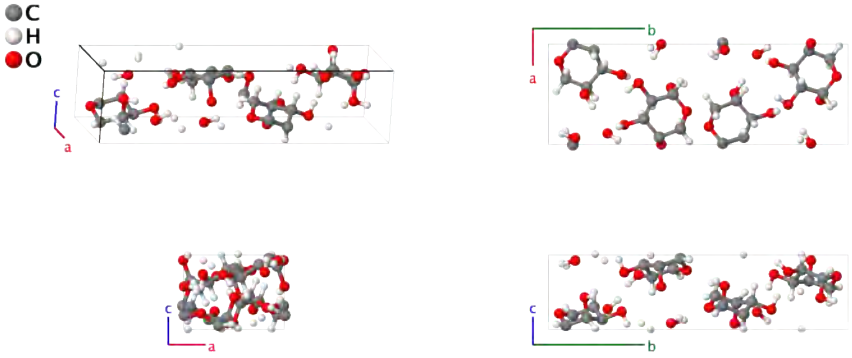
- T. Fukami, S. Akahoshi, K. Hukuda, and T. Yagi, *Refinement of the Crystal Structure of $NH_4H_2PO_4$ above and below Antiferroelectric Phase Transition Temperature*, J. Phys. Soc. Jpn. **56**, 2223–2224 (1987), doi:10.1143/JPSJ.56.2223.

Geometry files:

- CIF: pp. 1577
- POSCAR: pp. 1577

β-Arabinose [$(CH_2O)_{20}$] Structure: AB2C_oP80_19_5a_10a_5a

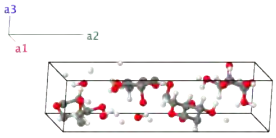
http://aflow.org/prototype-encyclopedia/AB2C_oP80_19_5a_10a_5a



Prototype	:	CH_2O
AFLOW prototype label	:	AB2C_oP80_19_5a_10a_5a
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP80
Space group number	:	19
Space group symbol	:	$P2_12_12_1$
AFLOW prototype command	:	`aflow --proto=AB2C_oP80_19_5a_10a_5a --params=`$a, b/a, c/a, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}, x_{17}, y_{17}, z_{17}, x_{18}, y_{18}, z_{18}, x_{19}, y_{19}, z_{19}, x_{20}, y_{20}, z_{20}$

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	$(4a)$	C I
$\mathbf{B}_2$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right)a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\,\hat{\mathbf{z}}$	$(4a)$	C I
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2} + y_1\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right)c\,\hat{\mathbf{z}}$	$(4a)$	C I
$\mathbf{B}_4$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_1\right)\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_1\right)b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	$(4a)$	C I
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(4a)$	C II
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	$(4a)$	C II
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	$(4a)$	C II

$\mathbf{B}_8$	=	$\left(\frac{1}{2}+x_2\right)\mathbf{a}_1+\left(\frac{1}{2}-y_2\right)\mathbf{a}_2-z_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_2\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_2\right)b\,\hat{\mathbf{y}}-z_2c\,\hat{\mathbf{z}}$	(4*a*)	C II
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1+y_3\,\mathbf{a}_2+z_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}}+y_3b\,\hat{\mathbf{y}}+z_3c\,\hat{\mathbf{z}}$	(4*a*)	C III
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2}-x_3\right)\mathbf{a}_1-y_3\,\mathbf{a}_2+\left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{x}}-y_3b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_3\right)c\,\hat{\mathbf{z}}$	(4*a*)	C III
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1+\left(\frac{1}{2}+y_3\right)\mathbf{a}_2+\left(\frac{1}{2}-z_3\right)\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_3\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_3\right)c\,\hat{\mathbf{z}}$	(4*a*)	C III
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2}+x_3\right)\mathbf{a}_1+\left(\frac{1}{2}-y_3\right)\mathbf{a}_2-z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_3\right)b\,\hat{\mathbf{y}}-z_3c\,\hat{\mathbf{z}}$	(4*a*)	C III
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1+y_4\,\mathbf{a}_2+z_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}}+y_4b\,\hat{\mathbf{y}}+z_4c\,\hat{\mathbf{z}}$	(4*a*)	C IV
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2}-x_4\right)\mathbf{a}_1-y_4\,\mathbf{a}_2+\left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{x}}-y_4b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_4\right)c\,\hat{\mathbf{z}}$	(4*a*)	C IV
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1+\left(\frac{1}{2}+y_4\right)\mathbf{a}_2+\left(\frac{1}{2}-z_4\right)\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_4\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_4\right)c\,\hat{\mathbf{z}}$	(4*a*)	C IV
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2}+x_4\right)\mathbf{a}_1+\left(\frac{1}{2}-y_4\right)\mathbf{a}_2-z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_4\right)b\,\hat{\mathbf{y}}-z_4c\,\hat{\mathbf{z}}$	(4*a*)	C IV
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1+y_5\,\mathbf{a}_2+z_5\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}}+y_5b\,\hat{\mathbf{y}}+z_5c\,\hat{\mathbf{z}}$	(4*a*)	C V
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2}-x_5\right)\mathbf{a}_1-y_5\,\mathbf{a}_2+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{x}}-y_5b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	(4*a*)	C V
$\mathbf{B}_{19}$	=	$-x_5\,\mathbf{a}_1+\left(\frac{1}{2}+y_5\right)\mathbf{a}_2+\left(\frac{1}{2}-z_5\right)\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_5\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_5\right)c\,\hat{\mathbf{z}}$	(4*a*)	C V
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2}+x_5\right)\mathbf{a}_1+\left(\frac{1}{2}-y_5\right)\mathbf{a}_2-z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_5\right)b\,\hat{\mathbf{y}}-z_5c\,\hat{\mathbf{z}}$	(4*a*)	C V
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1+y_6\,\mathbf{a}_2+z_6\,\mathbf{a}_3$	=	$x_6a\,\hat{\mathbf{x}}+y_6b\,\hat{\mathbf{y}}+z_6c\,\hat{\mathbf{z}}$	(4*a*)	H I
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2}-x_6\right)\mathbf{a}_1-y_6\,\mathbf{a}_2+\left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{x}}-y_6b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	(4*a*)	H I
$\mathbf{B}_{23}$	=	$-x_6\,\mathbf{a}_1+\left(\frac{1}{2}+y_6\right)\mathbf{a}_2+\left(\frac{1}{2}-z_6\right)\mathbf{a}_3$	=	$-x_6a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_6\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_6\right)c\,\hat{\mathbf{z}}$	(4*a*)	H I
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2}+x_6\right)\mathbf{a}_1+\left(\frac{1}{2}-y_6\right)\mathbf{a}_2-z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_6\right)b\,\hat{\mathbf{y}}-z_6c\,\hat{\mathbf{z}}$	(4*a*)	H I
$\mathbf{B}_{25}$	=	$x_7\,\mathbf{a}_1+y_7\,\mathbf{a}_2+z_7\,\mathbf{a}_3$	=	$x_7a\,\hat{\mathbf{x}}+y_7b\,\hat{\mathbf{y}}+z_7c\,\hat{\mathbf{z}}$	(4*a*)	H II
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2}-x_7\right)\mathbf{a}_1-y_7\,\mathbf{a}_2+\left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_7\right)a\,\hat{\mathbf{x}}-y_7b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	(4*a*)	H II
$\mathbf{B}_{27}$	=	$-x_7\,\mathbf{a}_1+\left(\frac{1}{2}+y_7\right)\mathbf{a}_2+\left(\frac{1}{2}-z_7\right)\mathbf{a}_3$	=	$-x_7a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_7\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_7\right)c\,\hat{\mathbf{z}}$	(4*a*)	H II
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2}+x_7\right)\mathbf{a}_1+\left(\frac{1}{2}-y_7\right)\mathbf{a}_2-z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_7\right)b\,\hat{\mathbf{y}}-z_7c\,\hat{\mathbf{z}}$	(4*a*)	H II
$\mathbf{B}_{29}$	=	$x_8\,\mathbf{a}_1+y_8\,\mathbf{a}_2+z_8\,\mathbf{a}_3$	=	$x_8a\,\hat{\mathbf{x}}+y_8b\,\hat{\mathbf{y}}+z_8c\,\hat{\mathbf{z}}$	(4*a*)	H III
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2}-x_8\right)\mathbf{a}_1-y_8\,\mathbf{a}_2+\left(\frac{1}{2}+z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_8\right)a\,\hat{\mathbf{x}}-y_8b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_8\right)c\,\hat{\mathbf{z}}$	(4*a*)	H III
$\mathbf{B}_{31}$	=	$-x_8\,\mathbf{a}_1+\left(\frac{1}{2}+y_8\right)\mathbf{a}_2+\left(\frac{1}{2}-z_8\right)\mathbf{a}_3$	=	$-x_8a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_8\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_8\right)c\,\hat{\mathbf{z}}$	(4*a*)	H III
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2}+x_8\right)\mathbf{a}_1+\left(\frac{1}{2}-y_8\right)\mathbf{a}_2-z_8\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_8\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_8\right)b\,\hat{\mathbf{y}}-z_8c\,\hat{\mathbf{z}}$	(4*a*)	H III
$\mathbf{B}_{33}$	=	$x_9\,\mathbf{a}_1+y_9\,\mathbf{a}_2+z_9\,\mathbf{a}_3$	=	$x_9a\,\hat{\mathbf{x}}+y_9b\,\hat{\mathbf{y}}+z_9c\,\hat{\mathbf{z}}$	(4*a*)	H IV
$\mathbf{B}_{34}$	=	$\left(\frac{1}{2}-x_9\right)\mathbf{a}_1-y_9\,\mathbf{a}_2+\left(\frac{1}{2}+z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_9\right)a\,\hat{\mathbf{x}}-y_9b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_9\right)c\,\hat{\mathbf{z}}$	(4*a*)	H IV
$\mathbf{B}_{35}$	=	$-x_9\,\mathbf{a}_1+\left(\frac{1}{2}+y_9\right)\mathbf{a}_2+\left(\frac{1}{2}-z_9\right)\mathbf{a}_3$	=	$-x_9a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_9\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_9\right)c\,\hat{\mathbf{z}}$	(4*a*)	H IV
$\mathbf{B}_{36}$	=	$\left(\frac{1}{2}+x_9\right)\mathbf{a}_1+\left(\frac{1}{2}-y_9\right)\mathbf{a}_2-z_9\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_9\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_9\right)b\,\hat{\mathbf{y}}-z_9c\,\hat{\mathbf{z}}$	(4*a*)	H IV
$\mathbf{B}_{37}$	=	$x_{10}\,\mathbf{a}_1+y_{10}\,\mathbf{a}_2+z_{10}\,\mathbf{a}_3$	=	$x_{10}a\,\hat{\mathbf{x}}+y_{10}b\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}}$	(4*a*)	H V
$\mathbf{B}_{38}$	=	$\left(\frac{1}{2}-x_{10}\right)\mathbf{a}_1-y_{10}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{10}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{x}}-y_{10}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{10}\right)c\,\hat{\mathbf{z}}$	(4*a*)	H V
$\mathbf{B}_{39}$	=	$-x_{10}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{10}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{10}\right)\mathbf{a}_3$	=	$-x_{10}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{10}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{10}\right)c\,\hat{\mathbf{z}}$	(4*a*)	H V
$\mathbf{B}_{40}$	=	$\left(\frac{1}{2}+x_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{10}\right)\mathbf{a}_2-z_{10}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{10}\right)b\,\hat{\mathbf{y}}-z_{10}c\,\hat{\mathbf{z}}$	(4*a*)	H V
$\mathbf{B}_{41}$	=	$x_{11}\,\mathbf{a}_1+y_{11}\,\mathbf{a}_2+z_{11}\,\mathbf{a}_3$	=	$x_{11}a\,\hat{\mathbf{x}}+y_{11}b\,\hat{\mathbf{y}}+z_{11}c\,\hat{\mathbf{z}}$	(4*a*)	H VI
$\mathbf{B}_{42}$	=	$\left(\frac{1}{2}-x_{11}\right)\mathbf{a}_1-y_{11}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{11}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{11}\right)a\,\hat{\mathbf{x}}-y_{11}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{11}\right)c\,\hat{\mathbf{z}}$	(4*a*)	H VI

$$
\begin{array}{lcccccl}
\mathbf{B}_{43} & = & -x_{11}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{11}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{11}\right)\mathbf{a}_3 & = & -x_{11}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{11}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{11}\right)c\,\hat{\mathbf{z}} & (4a) & \text{H VI} \\
\mathbf{B}_{44} & = & \left(\frac{1}{2} + x_{11}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{11}\right)\mathbf{a}_2 - z_{11}\,\mathbf{a}_3 & = & \left(\frac{1}{2} + x_{11}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{11}\right)b\,\hat{\mathbf{y}} - z_{11}c\,\hat{\mathbf{z}} & (4a) & \text{H VI} \\
\mathbf{B}_{45} & = & x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 & = & x_{12}a\,\hat{\mathbf{x}} + y_{12}b\,\hat{\mathbf{y}} + z_{12}c\,\hat{\mathbf{z}} & (4a) & \text{H VII} \\
\mathbf{B}_{46} & = & \left(\frac{1}{2} - x_{12}\right)\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{12}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_{12}\right)a\,\hat{\mathbf{x}} - y_{12}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{12}\right)c\,\hat{\mathbf{z}} & (4a) & \text{H VII} \\
\mathbf{B}_{47} & = & -x_{12}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{12}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{12}\right)\mathbf{a}_3 & = & -x_{12}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{12}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{12}\right)c\,\hat{\mathbf{z}} & (4a) & \text{H VII} \\
\mathbf{B}_{48} & = & \left(\frac{1}{2} + x_{12}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{12}\right)\mathbf{a}_2 - z_{12}\,\mathbf{a}_3 & = & \left(\frac{1}{2} + x_{12}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{12}\right)b\,\hat{\mathbf{y}} - z_{12}c\,\hat{\mathbf{z}} & (4a) & \text{H VII} \\
\mathbf{B}_{49} & = & x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3 & = & x_{13}a\,\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} + z_{13}c\,\hat{\mathbf{z}} & (4a) & \text{H VIII} \\
\mathbf{B}_{50} & = & \left(\frac{1}{2} - x_{13}\right)\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{13}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_{13}\right)a\,\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{13}\right)c\,\hat{\mathbf{z}} & (4a) & \text{H VIII} \\
\mathbf{B}_{51} & = & -x_{13}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{13}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{13}\right)\mathbf{a}_3 & = & -x_{13}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{13}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{13}\right)c\,\hat{\mathbf{z}} & (4a) & \text{H VIII} \\
\mathbf{B}_{52} & = & \left(\frac{1}{2} + x_{13}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{13}\right)\mathbf{a}_2 - z_{13}\,\mathbf{a}_3 & = & \left(\frac{1}{2} + x_{13}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{13}\right)b\,\hat{\mathbf{y}} - z_{13}c\,\hat{\mathbf{z}} & (4a) & \text{H VIII} \\
\mathbf{B}_{53} & = & x_{14}\,\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3 & = & x_{14}a\,\hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} + z_{14}c\,\hat{\mathbf{z}} & (4a) & \text{H IX} \\
\mathbf{B}_{54} & = & \left(\frac{1}{2} - x_{14}\right)\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{14}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_{14}\right)a\,\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{14}\right)c\,\hat{\mathbf{z}} & (4a) & \text{H IX} \\
\mathbf{B}_{55} & = & -x_{14}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{14}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{14}\right)\mathbf{a}_3 & = & -x_{14}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{14}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{14}\right)c\,\hat{\mathbf{z}} & (4a) & \text{H IX} \\
\mathbf{B}_{56} & = & \left(\frac{1}{2} + x_{14}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{14}\right)\mathbf{a}_2 - z_{14}\,\mathbf{a}_3 & = & \left(\frac{1}{2} + x_{14}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{14}\right)b\,\hat{\mathbf{y}} - z_{14}c\,\hat{\mathbf{z}} & (4a) & \text{H IX} \\
\mathbf{B}_{57} & = & x_{15}\,\mathbf{a}_1 + y_{15}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 & = & x_{15}a\,\hat{\mathbf{x}} + y_{15}b\,\hat{\mathbf{y}} + z_{15}c\,\hat{\mathbf{z}} & (4a) & \text{H X} \\
\mathbf{B}_{58} & = & \left(\frac{1}{2} - x_{15}\right)\mathbf{a}_1 - y_{15}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{15}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_{15}\right)a\,\hat{\mathbf{x}} - y_{15}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{15}\right)c\,\hat{\mathbf{z}} & (4a) & \text{H X} \\
\mathbf{B}_{59} & = & -x_{15}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{15}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{15}\right)\mathbf{a}_3 & = & -x_{15}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{15}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{15}\right)c\,\hat{\mathbf{z}} & (4a) & \text{H X} \\
\mathbf{B}_{60} & = & \left(\frac{1}{2} + x_{15}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{15}\right)\mathbf{a}_2 - z_{15}\,\mathbf{a}_3 & = & \left(\frac{1}{2} + x_{15}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{15}\right)b\,\hat{\mathbf{y}} - z_{15}c\,\hat{\mathbf{z}} & (4a) & \text{H X} \\
\mathbf{B}_{61} & = & x_{16}\,\mathbf{a}_1 + y_{16}\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 & = & x_{16}a\,\hat{\mathbf{x}} + y_{16}b\,\hat{\mathbf{y}} + z_{16}c\,\hat{\mathbf{z}} & (4a) & \text{O I} \\
\mathbf{B}_{62} & = & \left(\frac{1}{2} - x_{16}\right)\mathbf{a}_1 - y_{16}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{16}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_{16}\right)a\,\hat{\mathbf{x}} - y_{16}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{16}\right)c\,\hat{\mathbf{z}} & (4a) & \text{O I} \\
\mathbf{B}_{63} & = & -x_{16}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{16}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{16}\right)\mathbf{a}_3 & = & -x_{16}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{16}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{16}\right)c\,\hat{\mathbf{z}} & (4a) & \text{O I} \\
\mathbf{B}_{64} & = & \left(\frac{1}{2} + x_{16}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{16}\right)\mathbf{a}_2 - z_{16}\,\mathbf{a}_3 & = & \left(\frac{1}{2} + x_{16}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{16}\right)b\,\hat{\mathbf{y}} - z_{16}c\,\hat{\mathbf{z}} & (4a) & \text{O I} \\
\mathbf{B}_{65} & = & x_{17}\,\mathbf{a}_1 + y_{17}\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 & = & x_{17}a\,\hat{\mathbf{x}} + y_{17}b\,\hat{\mathbf{y}} + z_{17}c\,\hat{\mathbf{z}} & (4a) & \text{O II} \\
\mathbf{B}_{66} & = & \left(\frac{1}{2} - x_{17}\right)\mathbf{a}_1 - y_{17}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{17}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_{17}\right)a\,\hat{\mathbf{x}} - y_{17}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{17}\right)c\,\hat{\mathbf{z}} & (4a) & \text{O II} \\
\mathbf{B}_{67} & = & -x_{17}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{17}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{17}\right)\mathbf{a}_3 & = & -x_{17}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{17}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{17}\right)c\,\hat{\mathbf{z}} & (4a) & \text{O II} \\
\mathbf{B}_{68} & = & \left(\frac{1}{2} + x_{17}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{17}\right)\mathbf{a}_2 - z_{17}\,\mathbf{a}_3 & = & \left(\frac{1}{2} + x_{17}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{17}\right)b\,\hat{\mathbf{y}} - z_{17}c\,\hat{\mathbf{z}} & (4a) & \text{O II} \\
\mathbf{B}_{69} & = & x_{18}\,\mathbf{a}_1 + y_{18}\,\mathbf{a}_2 + z_{18}\,\mathbf{a}_3 & = & x_{18}a\,\hat{\mathbf{x}} + y_{18}b\,\hat{\mathbf{y}} + z_{18}c\,\hat{\mathbf{z}} & (4a) & \text{O III} \\
\mathbf{B}_{70} & = & \left(\frac{1}{2} - x_{18}\right)\mathbf{a}_1 - y_{18}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{18}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_{18}\right)a\,\hat{\mathbf{x}} - y_{18}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{18}\right)c\,\hat{\mathbf{z}} & (4a) & \text{O III} \\
\mathbf{B}_{71} & = & -x_{18}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{18}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{18}\right)\mathbf{a}_3 & = & -x_{18}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{18}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{18}\right)c\,\hat{\mathbf{z}} & (4a) & \text{O III} \\
\mathbf{B}_{72} & = & \left(\frac{1}{2} + x_{18}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{18}\right)\mathbf{a}_2 - z_{18}\,\mathbf{a}_3 & = & \left(\frac{1}{2} + x_{18}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{18}\right)b\,\hat{\mathbf{y}} - z_{18}c\,\hat{\mathbf{z}} & (4a) & \text{O III}
\end{array}
$$

$\mathbf{B}_{73}$	=	$x_{19}\,\mathbf{a}_1 + y_{19}\,\mathbf{a}_2 + z_{19}\,\mathbf{a}_3$	=	$x_{19}a\,\hat{\mathbf{x}} + y_{19}b\,\hat{\mathbf{y}} + z_{19}c\,\hat{\mathbf{z}}$	(4*a*)	O IV
$\mathbf{B}_{74}$	=	$\left(\frac{1}{2} - x_{19}\right)\mathbf{a}_1 - y_{19}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{19}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{19}\right)a\,\hat{\mathbf{x}} - y_{19}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{19}\right)c\,\hat{\mathbf{z}}$	(4*a*)	O IV
$\mathbf{B}_{75}$	=	$-x_{19}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{19}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{19}\right)\mathbf{a}_3$	=	$-x_{19}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{19}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{19}\right)c\,\hat{\mathbf{z}}$	(4*a*)	O IV
$\mathbf{B}_{76}$	=	$\left(\frac{1}{2} + x_{19}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{19}\right)\mathbf{a}_2 - z_{19}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{19}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{19}\right)b\,\hat{\mathbf{y}} - z_{19}c\,\hat{\mathbf{z}}$	(4*a*)	O IV
$\mathbf{B}_{77}$	=	$x_{20}\,\mathbf{a}_1 + y_{20}\,\mathbf{a}_2 + z_{20}\,\mathbf{a}_3$	=	$x_{20}a\,\hat{\mathbf{x}} + y_{20}b\,\hat{\mathbf{y}} + z_{20}c\,\hat{\mathbf{z}}$	(4*a*)	O V
$\mathbf{B}_{78}$	=	$\left(\frac{1}{2} - x_{20}\right)\mathbf{a}_1 - y_{20}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{20}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{20}\right)a\,\hat{\mathbf{x}} - y_{20}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{20}\right)c\,\hat{\mathbf{z}}$	(4*a*)	O V
$\mathbf{B}_{79}$	=	$-x_{20}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{20}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{20}\right)\mathbf{a}_3$	=	$-x_{20}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{20}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{20}\right)c\,\hat{\mathbf{z}}$	(4*a*)	O V
$\mathbf{B}_{80}$	=	$\left(\frac{1}{2} + x_{20}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{20}\right)\mathbf{a}_2 - z_{20}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{20}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{20}\right)b\,\hat{\mathbf{y}} - z_{20}c\,\hat{\mathbf{z}}$	(4*a*)	O V

References:

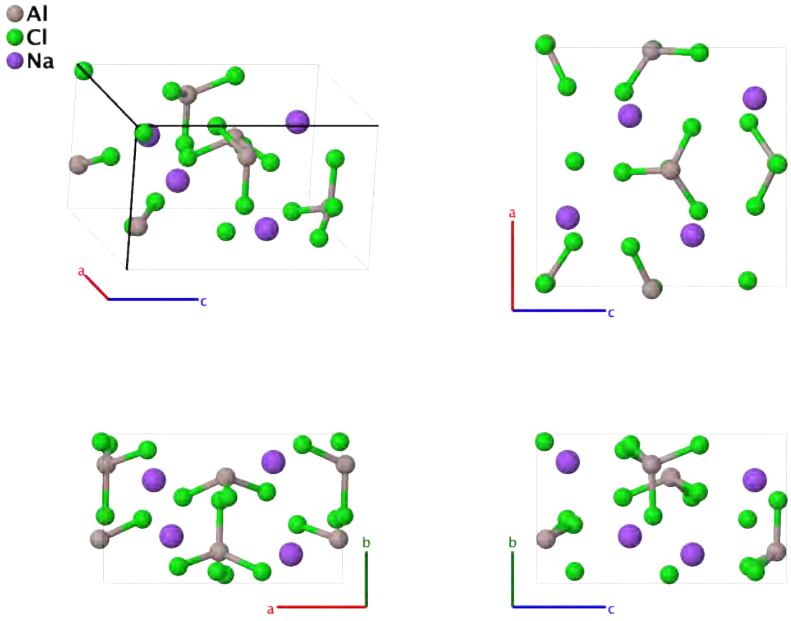
- A. Hordvik, *Refinement of the Crystal Structure of β-Arabinose*, Acta Chem. Scand. **15**, 16–30 (1961), doi:10.3891/acta.chem.scand.15-0016.

Geometry files:

- CIF: pp. 1577
- POSCAR: pp. 1578

$NaAlCl_4$ Structure: AB4C_oP24_19_a_4a_a

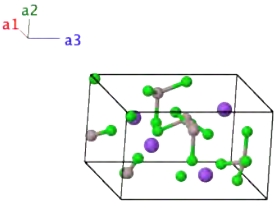
http://aflow.org/prototype-encyclopedia/AB4C_oP24_19_a_4a_a



Prototype	:	$AlCl_4Na$
AFLOW prototype label	:	AB4C_oP24_19_a_4a_a
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP24
Space group number	:	19
Space group symbol	:	$P2_12_12_1$
AFLOW prototype command	:	`aflow --proto=AB4C_oP24_19_a_4a_a --params=`$a, b/a, c/a, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6$

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	$(4a)$	Al
$\mathbf{B}_2$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right)a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\,\hat{\mathbf{z}}$	$(4a)$	Al

$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2}+y_1\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_1\right)\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_1\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_1\right) c\,\hat{\mathbf{z}}$	(4*a*)	Al
$\mathbf{B}_4$	=	$\left(\frac{1}{2}+x_1\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_1\right)\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_1\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_1\right) b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	(4*a*)	Al
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*a*)	Cl I
$\mathbf{B}_6$	=	$\left(\frac{1}{2}-x_2\right)\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_2\right) a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_2\right) c\,\hat{\mathbf{z}}$	(4*a*)	Cl I
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2}+y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_2\right)\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_2\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_2\right) c\,\hat{\mathbf{z}}$	(4*a*)	Cl I
$\mathbf{B}_8$	=	$\left(\frac{1}{2}+x_2\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_2\right)\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_2\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_2\right) b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(4*a*)	Cl I
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*a*)	Cl II
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2}-x_3\right)\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_3\right) a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_3\right) c\,\hat{\mathbf{z}}$	(4*a*)	Cl II
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2}+y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_3\right)\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_3\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_3\right) c\,\hat{\mathbf{z}}$	(4*a*)	Cl II
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2}+x_3\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_3\right)\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_3\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_3\right) b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(4*a*)	Cl II
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(4*a*)	Cl III
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2}-x_4\right)\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_4\right) a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_4\right) c\,\hat{\mathbf{z}}$	(4*a*)	Cl III
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2}+y_4\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_4\right)\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_4\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_4\right) c\,\hat{\mathbf{z}}$	(4*a*)	Cl III
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2}+x_4\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_4\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_4\right) b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(4*a*)	Cl III
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(4*a*)	Cl IV
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2}-x_5\right)\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_5\right) a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_5\right) c\,\hat{\mathbf{z}}$	(4*a*)	Cl IV
$\mathbf{B}_{19}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2}+y_5\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_5\right)\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_5\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_5\right) c\,\hat{\mathbf{z}}$	(4*a*)	Cl IV
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2}+x_5\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_5\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_5\right) b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(4*a*)	Cl IV
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(4*a*)	Na
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2}-x_6\right)\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_6\right) a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_6\right) c\,\hat{\mathbf{z}}$	(4*a*)	Na
$\mathbf{B}_{23}$	=	$-x_6\,\mathbf{a}_1 + \left(\frac{1}{2}+y_6\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_6\right)\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_6\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_6\right) c\,\hat{\mathbf{z}}$	(4*a*)	Na
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2}+x_6\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_6\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_6\right) b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(4*a*)	Na

References:

- G. Mairesse, P. Barbier, and J.-P. Wignacourt, *Comparison of the crystal structures of alkaline (M = Li, Na, K, Rb, Cs) and pseudo-alkaline (M = NO,NH$_4$) tetrachloroaluminates, MAlCl$_4$*, Acta Crystallogr. Sect. B Struct. Sci. **35**, 1573–1580 (1979), doi:10.1107/S0567740879007160.

Found in:

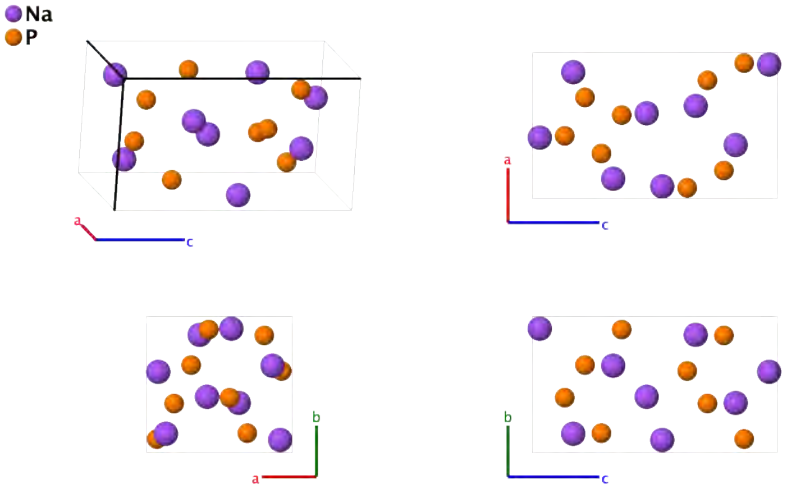
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1578
- POSCAR: pp. 1579

NaP Structure: AB_oP16_19_2a_2a

http://aflow.org/prototype-encyclopedia/AB_oP16_19_2a_2a



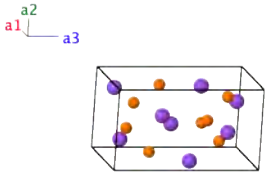
Prototype	:	NaP
AFLOW prototype label	:	AB_oP16_19_2a_2a
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP16
Space group number	:	19
Space group symbol	:	$P2_12_12_1$
AFLOW prototype command	:	`aflow --proto=AB_oP16_19_2a_2a --params=`$a, b/a, c/a, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4$

Other compounds with this structure

- CsSb and KP

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	$(4a)$	Na I
$\mathbf{B}_2$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right)a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\,\hat{\mathbf{z}}$	$(4a)$	Na I
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2} + y_1\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right)c\,\hat{\mathbf{z}}$	$(4a)$	Na I
$\mathbf{B}_4$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_1\right)\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_1\right)b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	$(4a)$	Na I

$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*a*)	Na II
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	(4*a*)	Na II
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right) c\,\hat{\mathbf{z}}$	(4*a*)	Na II
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right) b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(4*a*)	Na II
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*a*)	P I
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right) a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	(4*a*)	P I
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right) c\,\hat{\mathbf{z}}$	(4*a*)	P I
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right) b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(4*a*)	P I
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(4*a*)	P II
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	(4*a*)	P II
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\,\hat{\mathbf{z}}$	(4*a*)	P II
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right) b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(4*a*)	P II

References:

- H. G. von Schnering and W. Hönle, *Zur Chemie und Strukturchemie der Phosphide und Polyphosphide. 20. Darstellung, Struktur und Eigenschaften der Alkalimetallmonophosphide NaP und KP*, Z. Anorg. Allg. Chem. **456**, 194–206 (1979), doi:10.1002/zaac.19794560121.

Found in:

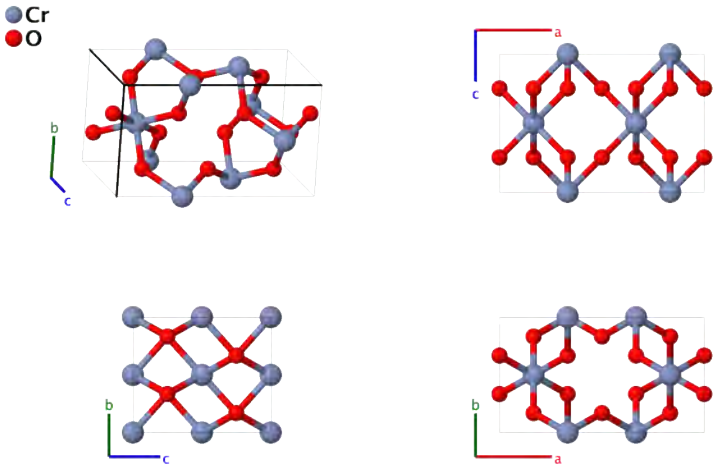
- P. Villars (Chief Editor), *NaP Crystal Structure*, http://materials.springer.com/isp/crystallographic/docs/sd_1250260 (2016). PAULING FILE in: Inorganic Solid Phases, SpringerMaterials (online database).

Geometry files:

- CIF: pp. 1579
- POSCAR: pp. 1579

$D0_7$ (CrO_3) (*obsolete*) Structure: AB3_oC16_20_a_bc

http://aflow.org/prototype-encyclopedia/AB3_oC16_20_a_bc

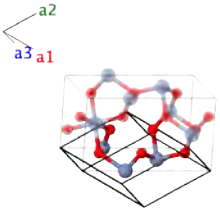


Prototype	:	CrO_3
AFLOW prototype label	:	AB3_oC16_20_a_bc
***Strukturbericht* designation**	:	$D0_7$
Pearson symbol	:	oC16
Space group number	:	20
Space group symbol	:	$C222_1$
AFLOW prototype command	:	`aflow --proto=AB3_oC16_20_a_bc --params=`$a, b/a, c/a, x_1, y_2, x_3, y_3, z_3$

- This is the determination of the orthorhombic structure of CrO_3 given the $D0_7$ designation by (Hermann, 1937). This structure is superceded by the one found by (Byström, 1950), which we call the "orthorhombic CrO_3 phase,". The current structure has the Cr atom at the center of a distorted oxygen octahedron, while the newer structure has the Cr atom at center of a distorted tetrahedron.

Base-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + x_1\,\mathbf{a}_2$	=	$x_1 a\,\hat{\mathbf{x}}$	$(4a)$	Cr
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 - x_1\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(4a)$	Cr

$\mathbf{B}_3$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_2 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*b*)	O I
$\mathbf{B}_4$	=	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_2 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(4*b*)	O I
$\mathbf{B}_5$	=	$(x_3 - y_3)\,\mathbf{a}_1 + (x_3 + y_3)\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(8*c*)	O II
$\mathbf{B}_6$	=	$(-x_3 + y_3)\,\mathbf{a}_1 + (-x_3 - y_3)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(8*c*)	O II
$\mathbf{B}_7$	=	$(-x_3 - y_3)\,\mathbf{a}_1 + (-x_3 + y_3)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(8*c*)	O II
$\mathbf{B}_8$	=	$(x_3 + y_3)\,\mathbf{a}_1 + (x_3 - y_3)\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(8*c*)	O II

References:

- H. Bräkken, *Die Kristallstrukturen der Trioxyde von Chrom, Molybdän und Wolfram*, Zeitschrift für Kristallographie - Crystalline Materials **78**, 484–488 (1931), doi:10.1524/zkri.1931.78.1.484.
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturebericht Band II, 1928-1932* (Akademsiche Verlagsgesellschaft M. B. H, Leipzig, 1937).

Found in:

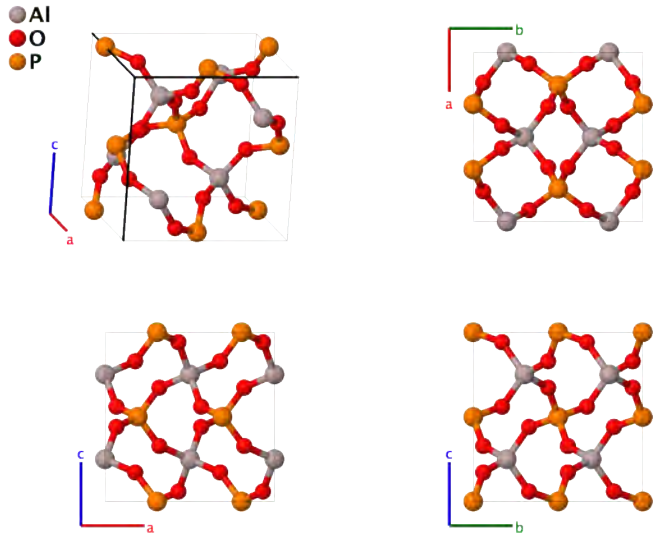
- A. Byström and K.-A. Wilhelmi, *The Crystal Structure of Chromium Trioxide*, Acta Chem. Scand. **4**, 1131–1141 (1950), doi:10.3891/acta.chem.scand.04-1131.

Geometry files:

- CIF: pp. 1579
- POSCAR: pp. 1580

$AlPO_4$ "low cristobalite type" Structure: AB4C_oC24_20_b_2c_a

http://aflow.org/prototype-encyclopedia/AB4C_oC24_20_b_2c_a



Prototype	:	AlO_4P
AFLOW prototype label	:	AB4C_oC24_20_b_2c_a
***Strukturbericht* designation**	:	None
Pearson symbol	:	oC24
Space group number	:	20
Space group symbol	:	$C222_1$
AFLOW prototype command	:	`aflow --proto=AB4C_oC24_20_b_2c_a --params=`$a, b/a, c/a, x_1, y_2, x_3, y_3, z_3, x_4, y_4, z_4$

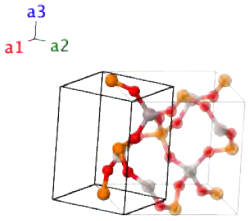
Other compounds with this structure

- $GaPO_4$

- Aluminum and Gallium Phosphate are closely related to the low cristobalite structure of SiO_2.
- The change in symmetry from $Fd\bar{3}m$ in SiO_2 to $C222_1$ in $AlPO_4$ is caused by the replacement of the silicon atoms by two types of atoms.
- The structure is "pseudo-tetragonal", with $a = b$, but the presence of ordered aluminum and phosphorus atoms reduces the symmetry to orthogonal.

Base-centered Orthorhombic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2}\,b\,\hat{\mathbf{y}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,b\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + x_1\,\mathbf{a}_2$	=	$x_1 a\,\hat{\mathbf{x}}$	($4a$)	P
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 - x_1\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4a$)	P
$\mathbf{B}_3$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_2 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($4b$)	Al
$\mathbf{B}_4$	=	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_2 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	($4b$)	Al
$\mathbf{B}_5$	=	$(x_3 - y_3)\,\mathbf{a}_1 + (x_3 + y_3)\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	($8c$)	O I
$\mathbf{B}_6$	=	$(-x_3 + y_3)\,\mathbf{a}_1 + (-x_3 - y_3)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	($8c$)	O I
$\mathbf{B}_7$	=	$(-x_3 - y_3)\,\mathbf{a}_1 + (-x_3 + y_3)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	($8c$)	O I
$\mathbf{B}_8$	=	$(x_3 + y_3)\,\mathbf{a}_1 + (x_3 - y_3)\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	($8c$)	O I
$\mathbf{B}_9$	=	$(x_4 - y_4)\,\mathbf{a}_1 + (x_4 + y_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	($8c$)	O II
$\mathbf{B}_{10}$	=	$(-x_4 + y_4)\,\mathbf{a}_1 + (-x_4 - y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	($8c$)	O II
$\mathbf{B}_{11}$	=	$(-x_4 - y_4)\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	($8c$)	O II
$\mathbf{B}_{12}$	=	$(x_4 + y_4)\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	($8c$)	O II

References:

- R. C. L. Mooney, *The crystal structure of aluminium phosphate and gallium phosphate, low-cristobalite type*, Acta Cryst. **9**, 728–734 (1956), doi:10.1107/S0365110X56001996.

Found in:

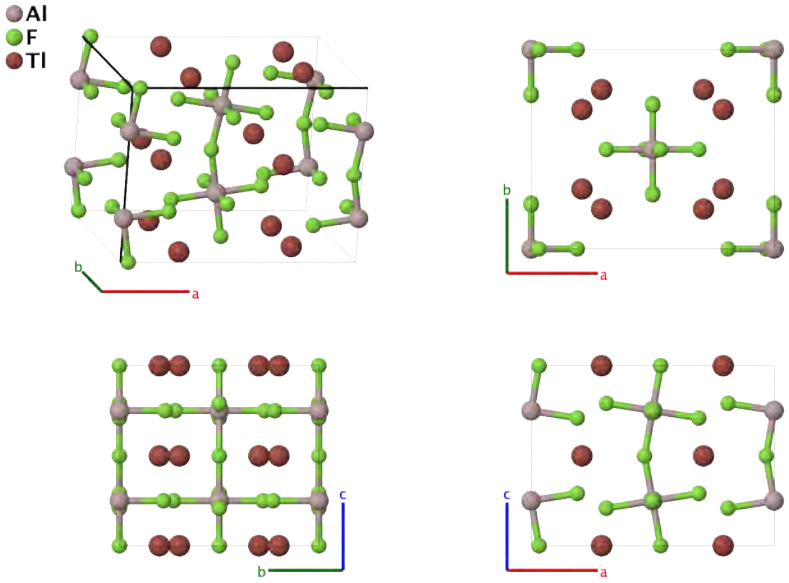
- D. M. Hatch, S. Ghose, and J. L. Bjorkstam, *The $\alpha - \beta$ phase transition in $AlPO_4$ cristobalite: Symmetry analysis, domain structure and transition dynamics*, Phys. Chem. Miner. **21**, 67–77 (1994), doi:10.1007/BF00205217.

Geometry files:

- CIF: pp. 1580
- POSCAR: pp. 1580

Tl_2AlF_5 ($K3_3$) Structure: AB5C2_oC32_20_b_a2bc_c

http://aflow.org/prototype-encyclopedia/AB5C2_oC32_20_b_a2bc_c

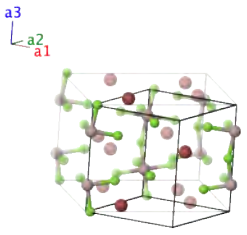


Prototype	:	AlF_5Tl_2
AFLOW prototype label	:	AB5C2_oC32_20_b_a2bc_c
***Strukturbericht* designation**	:	$K3_3$
Pearson symbol	:	oC32
Space group number	:	20
Space group symbol	:	$C222_1$
AFLOW prototype command	:	`aflow --proto=AB5C2_oC32_20_b_a2bc_c --params=`$a, b/a, c/a, x_1, y_2, y_3, y_4, x_5, y_5, z_5, x_6, y_6, z_6$

- There are several problems with this structure:
- First, (Brosset, 1937) gives coordinates $y_3 = -y_2$, which gives the structure an inversion site and makes the space group *Cmcm*. We have adjusted the coordinates of the third atom slightly to avoid this.
- Second, (Molinier, 1993) makes the argument that the structure observed by Brosset is actually $Tl_2AlF_5{\cdot}H_2O$.
- Finally, (Pabst, 1950) indicates that a 1942 work of Brosset refines the structure, but we have not been able to access this paper in any form, nor have we found any reference to this structure in the literature.

Base-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + x_1\,\mathbf{a}_2$	=	$x_1 a\,\hat{\mathbf{x}}$	(4*a*)	F I
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 - x_1\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*a*)	F I
$\mathbf{B}_3$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_2 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*b*)	Al
$\mathbf{B}_4$	=	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_2 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(4*b*)	Al
$\mathbf{B}_5$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_3 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*b*)	F II
$\mathbf{B}_6$	=	$y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_3 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(4*b*)	F II
$\mathbf{B}_7$	=	$-y_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_4 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*b*)	F III
$\mathbf{B}_8$	=	$y_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_4 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(4*b*)	F III
$\mathbf{B}_9$	=	$(x_5 - y_5)\,\mathbf{a}_1 + (x_5 + y_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(8*c*)	F IV
$\mathbf{B}_{10}$	=	$(-x_5 + y_5)\,\mathbf{a}_1 + (-x_5 - y_5)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	(8*c*)	F IV
$\mathbf{B}_{11}$	=	$(-x_5 - y_5)\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	(8*c*)	F IV
$\mathbf{B}_{12}$	=	$(x_5 + y_5)\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(8*c*)	F IV
$\mathbf{B}_{13}$	=	$(x_6 - y_6)\,\mathbf{a}_1 + (x_6 + y_6)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(8*c*)	Tl
$\mathbf{B}_{14}$	=	$(-x_6 + y_6)\,\mathbf{a}_1 + (-x_6 - y_6)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\,\hat{\mathbf{z}}$	(8*c*)	Tl
$\mathbf{B}_{15}$	=	$(-x_6 - y_6)\,\mathbf{a}_1 + (-x_6 + y_6)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\,\hat{\mathbf{z}}$	(8*c*)	Tl
$\mathbf{B}_{16}$	=	$(x_6 + y_6)\,\mathbf{a}_1 + (x_6 - y_6)\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(8*c*)	Tl

References:

- C. Brosset, *Herstellung und Kristallbau der Verbindungen $TlAlF_4$ und Tl_2AlF_5*, Z. Anorg. Allg. Chem. **235**, 139–147 (1937), doi:10.1002/zaac.19372350119.
- M. Molinier and W. Massa, *Refinement of the structure of $Tl_2AlF_5 \cdot H_2O$*, Acta Crystallogr. C **49**, 782–784 (1993), doi:10.1107/S010827019201148X.

Found in:

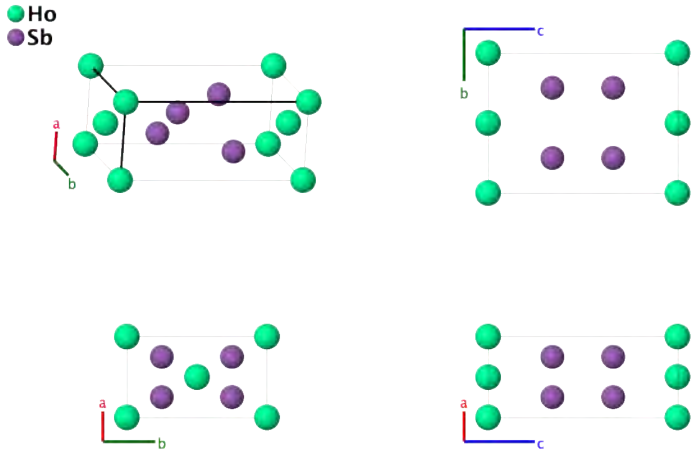
- A. Pabst, *A Structural Classification of Fluoaluminates*, Am. Mineral. **35**, 149–165 (1950).

Geometry files:

- CIF: pp. 1580
- POSCAR: pp. 1581

$HoSb_2$ Structure: AB2_oC6_21_a_k

http://aflow.org/prototype-encyclopedia/AB2_oC6_21_a_k.HoSb2



Prototype	:	$HoSb_2$
AFLOW prototype label	:	AB2_oC6_21_a_k
***Strukturbericht* designation**	:	None
Pearson symbol	:	oC6
Space group number	:	21
Space group symbol	:	$C222$
AFLOW prototype command	:	`aflow --proto=AB2_oC6_21_a_k --params=`$a, b/a, c/a, z_2$

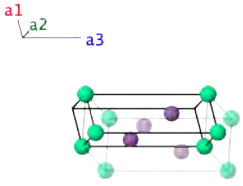
Other compounds with this structure

- $LuSb_2$, YSb_2, $DySb_2$ (HT), $GdSb_2$ (HT), and $TbSb_2$ (HT)

- Measurements were performed at a pressure of 65 kbar.
- The primitive unit cell is nearly hexagonal.
- The author states "We are well aware of the fact that the structure presented here may indeed be only a subcell of the true structure."
- $HoSb_2$ and Ta_2H share the same AFLOW prototype label, AB2_oC6_21_a_k. The structures are generated by the same symmetry operations with different sets of parameters (`--params`) specified in their corresponding CIF files.

Base-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($2a$)	Ho
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	($4k$)	Sb
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_1 - z_2\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} - \frac{1}{4}b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	($4k$)	Sb

References:

- Q. Johnson, *The Crystal Structure of High-Pressure Synthesized Holmium Diantimonde*, Inorg. Chem. **10**, 2089–2090 (1971), doi:10.1021/ic50103a059.

Found in:

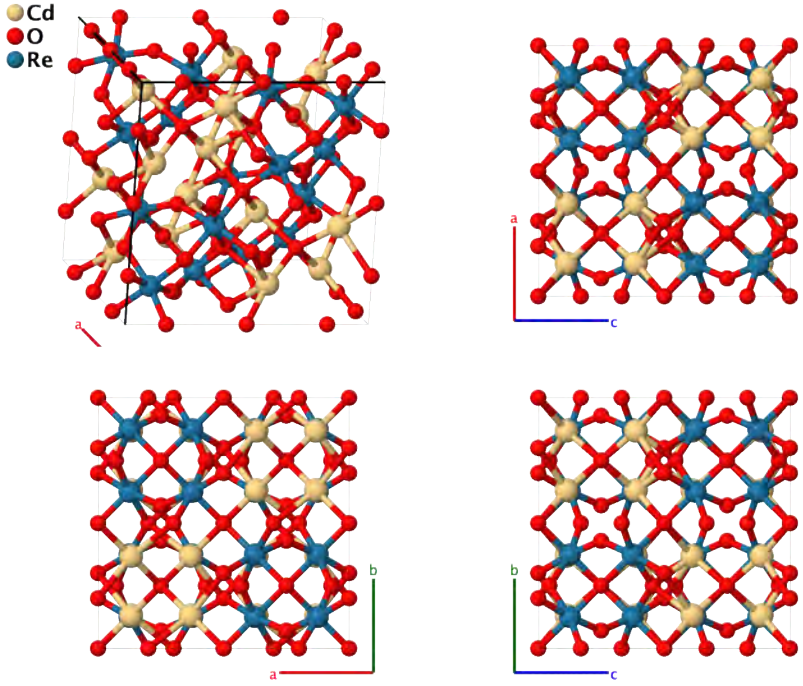
- M. N. Abdusaljamova, O. R. Burnashev, and K. E. Mironov, *The Ho-Sb Alloy System*, J. Less-Common Met. **102**, L19–L22 (1984), doi:10.1016/0022-5088(84)90403-X.

Geometry files:

- CIF: pp. 1581
- POSCAR: pp. 1581

Predicted Phase IV $Cd_2Re_2O_7$ Structure: A2B7C2_oF88_22_k_bdefghij_k

http://aflow.org/prototype-encyclopedia/A2B7C2_oF88_22_k_bdefghij_k

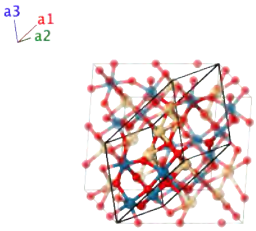


Prototype	:	$Cd_2O_7Re_2$
AFLOW prototype label	:	A2B7C2_oF88_22_k_bdefghij_k
***Strukturbericht* designation**	:	None
Pearson symbol	:	oF88
Space group number	:	22
Space group symbol	:	$F222$
AFLOW prototype command	:	`aflow --proto=A2B7C2_oF88_22_k_bdefghij_k --params=`$a, b/a, c/a, x_3, y_4, z_5, z_6, y_7, x_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}$

- $Cd_2Re_2O_7$ exhibits a number of phases. We will use the notation of (Kapcia, 2019) to describe them:
  - Phase I: above 200 K, the system takes on the cubic pyrochlore ($E8_1$) structure.
  - Phase II: in the range 120-200 K the system is in the tetragonal $I\bar{4}m2$ #119 structure.
  - Phase III: in the range 80-120 K the system is in the tetragonal $I4_122$ #98 structure.
  - Phase IV: (Kapcia, 2019) did a first-principles study of this system and found that below 80 K Phase III develops a soft phonon mode which transforms the system into an orthorhombic $F222$ #22 structure. (This structure)
  - (Norman, 2019) points out that both Phase III and Phase IV structures have issues.
- Phase IV is extremely close to Phase II. If AFLOW-SYM and FINDSYM allow symmetry tolerances of 0.002 Å, the orthorhombic phase becomes tetragonal.

Face-centered Orthorhombic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,b\,\hat{\mathbf{y}} + \tfrac{1}{2}\,c\,\hat{\mathbf{z}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,c\,\hat{\mathbf{z}}\\ \mathbf{a}_3 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,b\,\hat{\mathbf{y}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*b*)	O I
$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(4*d*)	O II
$\mathbf{B}_3$	=	$-x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}}$	(8*e*)	O III
$\mathbf{B}_4$	=	$x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}}$	(8*e*)	O III
$\mathbf{B}_5$	=	$y_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + y_4\,\mathbf{a}_3$	=	$y_4 b\,\hat{\mathbf{y}}$	(8*f*)	O IV
$\mathbf{B}_6$	=	$-y_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 - y_4\,\mathbf{a}_3$	=	$-y_4 b\,\hat{\mathbf{y}}$	(8*f*)	O IV
$\mathbf{B}_7$	=	$z_5\,\mathbf{a}_1 + z_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$z_5 c\,\hat{\mathbf{z}}$	(8*g*)	O V
$\mathbf{B}_8$	=	$-z_5\,\mathbf{a}_1 - z_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$-z_5 c\,\hat{\mathbf{z}}$	(8*g*)	O V
$\mathbf{B}_9$	=	$z_6\,\mathbf{a}_1 + z_6\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(8*h*)	O VI
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - z_6\right)\mathbf{a}_1 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\,\hat{\mathbf{z}}$	(8*h*)	O VI
$\mathbf{B}_{11}$	=	$y_7\,\mathbf{a}_1 + \left(\frac{1}{2} - y_7\right)\mathbf{a}_2 + y_7\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(8*i*)	O VII
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} - y_7\right)\mathbf{a}_1 + y_7\,\mathbf{a}_2 + \left(\frac{1}{2} - y_7\right)\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_7\right)b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(8*i*)	O VII
$\mathbf{B}_{13}$	=	$\left(\frac{1}{2} - x_8\right)\mathbf{a}_1 + x_8\,\mathbf{a}_2 + x_8\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(8*j*)	O VIII
$\mathbf{B}_{14}$	=	$x_8\,\mathbf{a}_1 + \left(\frac{1}{2} - x_8\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_8\right)a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(8*j*)	O VIII
$\mathbf{B}_{15}$	=	$(-x_9 + y_9 + z_9)\,\mathbf{a}_1 + (x_9 - y_9 + z_9)\,\mathbf{a}_2 + (x_9 + y_9 - z_9)\,\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(16*k*)	Cd
$\mathbf{B}_{16}$	=	$(x_9 - y_9 + z_9)\,\mathbf{a}_1 + (-x_9 + y_9 + z_9)\,\mathbf{a}_2 + (-x_9 - y_9 - z_9)\,\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(16*k*)	Cd
$\mathbf{B}_{17}$	=	$(x_9 + y_9 - z_9)\,\mathbf{a}_1 + (-x_9 - y_9 - z_9)\,\mathbf{a}_2 + (-x_9 + y_9 + z_9)\,\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}}$	(16*k*)	Cd
$\mathbf{B}_{18}$	=	$(-x_9 - y_9 - z_9)\,\mathbf{a}_1 + (x_9 + y_9 - z_9)\,\mathbf{a}_2 + (x_9 - y_9 + z_9)\,\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}}$	(16*k*)	Cd
$\mathbf{B}_{19}$	=	$(-x_{10} + y_{10} + z_{10})\,\mathbf{a}_1 + (x_{10} - y_{10} + z_{10})\,\mathbf{a}_2 + (x_{10} + y_{10} - z_{10})\,\mathbf{a}_3$	=	$x_{10} a\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}}$	(16*k*)	Re
$\mathbf{B}_{20}$	=	$(x_{10} - y_{10} + z_{10})\,\mathbf{a}_1 + (-x_{10} + y_{10} + z_{10})\,\mathbf{a}_2 + (-x_{10} - y_{10} - z_{10})\,\mathbf{a}_3$	=	$-x_{10} a\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}}$	(16*k*)	Re
$\mathbf{B}_{21}$	=	$(x_{10} + y_{10} - z_{10})\,\mathbf{a}_1 + (-x_{10} - y_{10} - z_{10})\,\mathbf{a}_2 + (-x_{10} + y_{10} + z_{10})\,\mathbf{a}_3$	=	$-x_{10} a\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} - z_{10} c\,\hat{\mathbf{z}}$	(16*k*)	Re
$\mathbf{B}_{22}$	=	$(-x_{10} - y_{10} - z_{10})\,\mathbf{a}_1 + (x_{10} + y_{10} - z_{10})\,\mathbf{a}_2 + (x_{10} - y_{10} + z_{10})\,\mathbf{a}_3$	=	$x_{10} a\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} - z_{10} c\,\hat{\mathbf{z}}$	(16*k*)	Re

References:

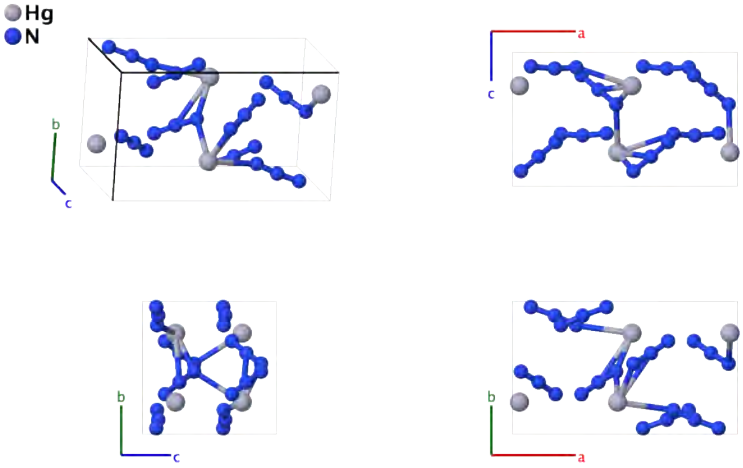
- K. J. Kapcia, M. Reedyk, M. Hajialamdari, A. Ptok, P. Piekarz, F. S. Razavi, A. M. Oleś, and R. K. Kremer, *Discovery of a low-temperature orthorhombic phase of the $Cd_2Re_2O_7$ superconductor*, Phys. Rev. Research **2**, 033108 (2020), doi:10.1103/PhysRevResearch.2.033108.

Geometry files:

- CIF: pp. 1581
- POSCAR: pp. 1582

Mercury (II) Azide [$Hg(N_3)_2$] Structure: AB6_oP28_29_a_6a

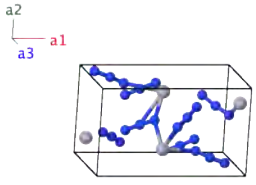
http://aflow.org/prototype-encyclopedia/AB6_oP28_29_a_6a



Prototype	:	HgN_6
AFLOW prototype label	:	AB6_oP28_29_a_6a
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP28
Space group number	:	29
Space group symbol	:	$Pca2_1$
AFLOW prototype command	:	`aflow --proto=AB6_oP28_29_a_6a --params=`$a, b/a, c/a, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7$

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(4*a*)	Hg
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	(4*a*)	Hg
$\mathbf{B}_3$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 - y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right) a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(4*a*)	Hg
$\mathbf{B}_4$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + y_1\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	(4*a*)	Hg
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*a*)	N I

$\mathbf{B}_{6}$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	(4*a*)	N I
$\mathbf{B}_{7}$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 - y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*a*)	N I
$\mathbf{B}_{8}$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	(4*a*)	N I
$\mathbf{B}_{9}$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*a*)	N II
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	(4*a*)	N II
$\mathbf{B}_{11}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 - y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right) a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*a*)	N II
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right) a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	(4*a*)	N II
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(4*a*)	N III
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	(4*a*)	N III
$\mathbf{B}_{15}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 - y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(4*a*)	N III
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	(4*a*)	N III
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(4*a*)	N IV
$\mathbf{B}_{18}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	(4*a*)	N IV
$\mathbf{B}_{19}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 - y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(4*a*)	N IV
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + y_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right) a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	(4*a*)	N IV
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(4*a*)	N V
$\mathbf{B}_{22}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	(4*a*)	N V
$\mathbf{B}_{23}$	=	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 - y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(4*a*)	N V
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 + y_6\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	(4*a*)	N V
$\mathbf{B}_{25}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(4*a*)	N VI
$\mathbf{B}_{26}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right) c\,\hat{\mathbf{z}}$	(4*a*)	N VI
$\mathbf{B}_{27}$	=	$\left(\frac{1}{2} + x_7\right)\mathbf{a}_1 - y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_7\right) a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(4*a*)	N VI
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} - x_7\right)\mathbf{a}_1 + y_7\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_7\right) a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right) c\,\hat{\mathbf{z}}$	(4*a*)	N VI

References:

- U. Müller, *Die Kristallstruktur von α-Quecksilber(II)-Azid*, Z. Anorg. Allg. Chem. **399**, 183–192 (1973), doi:10.1002/zaac.19733990207.

Found in:

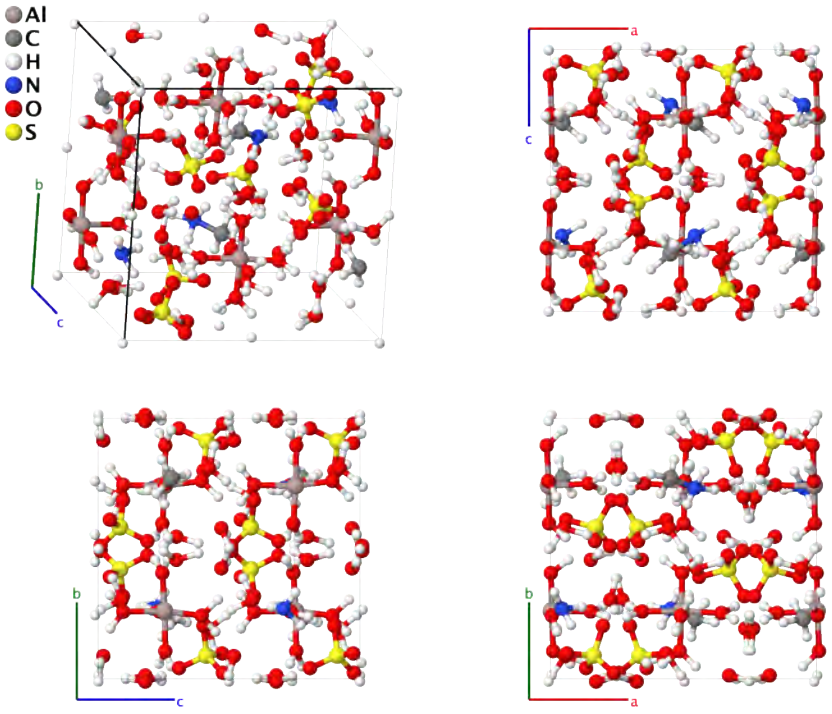
- T. B. Massalski, H. Okamoto, P. R. Subramanian, and L. Kacprzak, eds., *Binary Alloy Phase Diagrams*, vol. 3 (ASM International, Materials Park, Ohio, USA, 1990), 2nd edn. Hf-Re to Zn-Zr.

Geometry files:

- CIF: pp. 1582
- POSCAR: pp. 1582

Low-Temperature $(NH_3CH_3)Al(SO_4)_2{\cdot}12H_2O$ Structure: ABC30DE20F2_oP220_29_a_a_30a_a_20a_2a

http://aflow.org/prototype-encyclopedia/ABC30DE20F2_oP220_29_a_a_30a_a_20a_2a



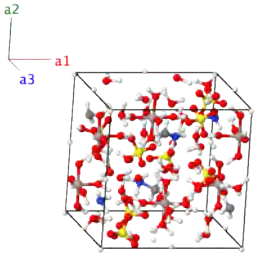
Prototype	:	$AlCH_{30}NO_{20}S_2$
AFLOW prototype label	:	ABC30DE20F2_oP220_29_a_a_30a_a_20a_2a
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP220
Space group number	:	29
Space group symbol	:	$Pca2_1$
AFLOW prototype command	:	`aflow --proto=ABC30DE20F2_oP220_29_a_a_30a_a_20a_2a --params=`$a, b/a, c/a, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}, x_{17}, y_{17}, z_{17}, x_{18}, y_{18}, z_{18}, x_{19}, y_{19}, z_{19}, x_{20}, y_{20}, z_{20}, x_{21}, y_{21}, z_{21}, x_{22}, y_{22}, z_{22}, x_{23}, y_{23}, z_{23}, x_{24}, y_{24}, z_{24}, x_{25}, y_{25}, z_{25}, x_{26}, y_{26}, z_{26}, x_{27}, y_{27}, z_{27}, x_{28}, y_{28}, z_{28}, x_{29}, y_{29}, z_{29}, x_{30}, y_{30}, z_{30}, x_{31}, y_{31}, z_{31}, x_{32}, y_{32}, z_{32}, x_{33}, y_{33}, z_{33}, x_{34}, y_{34}, z_{34}, x_{35}, y_{35}, z_{35}, x_{36}, y_{36}, z_{36}, x_{37}, y_{37}, z_{37}, x_{38}, y_{38}, z_{38}, x_{39}, y_{39}, z_{39}, x_{40}, y_{40}, z_{40}, x_{41}, y_{41}, z_{41}, x_{42}, y_{42}, z_{42}, x_{43}, y_{43}, z_{43}, x_{44}, y_{44}, z_{44}, x_{45}, y_{45}, z_{45}, x_{46}, y_{46}, z_{46}, x_{47}, y_{47}, z_{47}, x_{48}, y_{48}, z_{48}, x_{49}, y_{49}, z_{49}, x_{50}, y_{50}, z_{50}, x_{51}, y_{51}, z_{51}, x_{52}, y_{52}, z_{52}, x_{53}, y_{53}, z_{53}, x_{54}, y_{54}, z_{54}, x_{55}, y_{55}, z_{55}$

- The alums have the general formula $AB(XO_4)_2{\cdot}12H_2O$, where A is a monovalent ion, B is a trivalent ion, and X is a chalcogen. In most cases atom B is aluminum and atom X is sulfur, leading to the name alum.
- All alums have their room-temperature form in space group $Pa\bar{3}$ #205, but the bonding between the A and B ions and the XO_4 complex can be quite different.

- (Lipson, 1935ab) described three general forms of alum based on the sizes of the monovalent ions. Each of these forms was given a *Strukturbericht* designation by (Gottfried, 1937):
  - α-alum, with intermediate sized ions, prototype $KAl(SO_4)_2{\cdot}12H_2O$, $H4_{13}$,
  - β-alum, with large ions, prototype $(NH_3CH_3)Al(SO_4)_2{\cdot}12H_2O$, $H4_{14}$, and
  - γ-alum, with small ions, prototype $NaAl(SO_4)_2{\cdot}12H_2O$, $H4_{15}$.
- This classification scheme is not complete, *e.g.*, (Ledsham, 1968) points out that $NaCr(SO_4)_2{\cdot}12H_2O$ does not fit into any of these categories, and that the actual structure depends on the combination of monovalent and trivalent ions.
- As noted above, the $Pa\overline{3}$ structures of alum are the room temperature form. As the temperature decreases the alum structure may transform. For example, in the temperature range 150-170 K, the β-alum $(NH_3CH_3)Al(SO_4)_2{\cdot}12H_2O$ transforms into this orthorhombic structure with fully ordered NH_3CH_3 ions.
- The data presented here was taken at 113 K.

Simple Orthorhombic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= a\,\hat{\mathbf{x}}\\ \mathbf{a}_2 &= b\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	$(4a)$	Al
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	$(4a)$	Al
$\mathbf{B}_3$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 - y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right) a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	$(4a)$	Al
$\mathbf{B}_4$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + y_1\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	$(4a)$	Al
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(4a)$	C
$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	$(4a)$	C
$\mathbf{B}_7$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 - y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(4a)$	C
$\mathbf{B}_8$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	$(4a)$	C
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(4a)$	H I
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	$(4a)$	H I
$\mathbf{B}_{11}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 - y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right) a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(4a)$	H I
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right) a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	$(4a)$	H I
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(4a)$	H II
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	$(4a)$	H II
$\mathbf{B}_{15}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 - y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(4a)$	H II
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	$(4a)$	H II
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(4a)$	H III

$$
\begin{aligned}
\mathbf{B}_{18} &= -x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_3 &&= -x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_5\right) c\,\hat{\mathbf{z}} && (4a) && \text{H III}\\
\mathbf{B}_{19} &= \left(\tfrac{1}{2} + x_5\right)\mathbf{a}_1 - y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}} && (4a) && \text{H III}\\
\mathbf{B}_{20} &= \left(\tfrac{1}{2} - x_5\right)\mathbf{a}_1 + y_5\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_5\right) a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_5\right) c\,\hat{\mathbf{z}} && (4a) && \text{H III}\\
\mathbf{B}_{21} &= x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 &&= x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}} && (4a) && \text{H IV}\\
\mathbf{B}_{22} &= -x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 &&= -x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right) c\,\hat{\mathbf{z}} && (4a) && \text{H IV}\\
\mathbf{B}_{23} &= \left(\tfrac{1}{2} + x_6\right)\mathbf{a}_1 - y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}} && (4a) && \text{H IV}\\
\mathbf{B}_{24} &= \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_1 + y_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right) c\,\hat{\mathbf{z}} && (4a) && \text{H IV}\\
\mathbf{B}_{25} &= x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 &&= x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}} && (4a) && \text{H V}\\
\mathbf{B}_{26} &= -x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_7\right)\mathbf{a}_3 &&= -x_7 a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_7\right) c\,\hat{\mathbf{z}} && (4a) && \text{H V}\\
\mathbf{B}_{27} &= \left(\tfrac{1}{2} + x_7\right)\mathbf{a}_1 - y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_7\right) a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}} && (4a) && \text{H V}\\
\mathbf{B}_{28} &= \left(\tfrac{1}{2} - x_7\right)\mathbf{a}_1 + y_7\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_7\right) a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_7\right) c\,\hat{\mathbf{z}} && (4a) && \text{H V}\\
\mathbf{B}_{29} &= x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 &&= x_8 a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}} && (4a) && \text{H VI}\\
\mathbf{B}_{30} &= -x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_8\right)\mathbf{a}_3 &&= -x_8 a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_8\right) c\,\hat{\mathbf{z}} && (4a) && \text{H VI}\\
\mathbf{B}_{31} &= \left(\tfrac{1}{2} + x_8\right)\mathbf{a}_1 - y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_8\right) a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}} && (4a) && \text{H VI}\\
\mathbf{B}_{32} &= \left(\tfrac{1}{2} - x_8\right)\mathbf{a}_1 + y_8\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_8\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_8\right) a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_8\right) c\,\hat{\mathbf{z}} && (4a) && \text{H VI}\\
\mathbf{B}_{33} &= x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 &&= x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}} && (4a) && \text{H VII}\\
\mathbf{B}_{34} &= -x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_3 &&= -x_9 a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_9\right) c\,\hat{\mathbf{z}} && (4a) && \text{H VII}\\
\mathbf{B}_{35} &= \left(\tfrac{1}{2} + x_9\right)\mathbf{a}_1 - y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_9\right) a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}} && (4a) && \text{H VII}\\
\mathbf{B}_{36} &= \left(\tfrac{1}{2} - x_9\right)\mathbf{a}_1 + y_9\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_9\right) a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_9\right) c\,\hat{\mathbf{z}} && (4a) && \text{H VII}\\
\mathbf{B}_{37} &= x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 &&= x_{10} a\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}} && (4a) && \text{H VIII}\\
\mathbf{B}_{38} &= -x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{10}\right)\mathbf{a}_3 &&= -x_{10} a\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{10}\right) c\,\hat{\mathbf{z}} && (4a) && \text{H VIII}\\
\mathbf{B}_{39} &= \left(\tfrac{1}{2} + x_{10}\right)\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_{10}\right) a\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}} && (4a) && \text{H VIII}\\
\mathbf{B}_{40} &= \left(\tfrac{1}{2} - x_{10}\right)\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{10}\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_{10}\right) a\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{10}\right) c\,\hat{\mathbf{z}} && (4a) && \text{H VIII}\\
\mathbf{B}_{41} &= x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 &&= x_{11} a\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + z_{11} c\,\hat{\mathbf{z}} && (4a) && \text{H IX}\\
\mathbf{B}_{42} &= -x_{11}\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{11}\right)\mathbf{a}_3 &&= -x_{11} a\,\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{11}\right) c\,\hat{\mathbf{z}} && (4a) && \text{H IX}\\
\mathbf{B}_{43} &= \left(\tfrac{1}{2} + x_{11}\right)\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_{11}\right) a\,\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} + z_{11} c\,\hat{\mathbf{z}} && (4a) && \text{H IX}\\
\mathbf{B}_{44} &= \left(\tfrac{1}{2} - x_{11}\right)\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{11}\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_{11}\right) a\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{11}\right) c\,\hat{\mathbf{z}} && (4a) && \text{H IX}\\
\mathbf{B}_{45} &= x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 &&= x_{12} a\,\hat{\mathbf{x}} + y_{12} b\,\hat{\mathbf{y}} + z_{12} c\,\hat{\mathbf{z}} && (4a) && \text{H X}\\
\mathbf{B}_{46} &= -x_{12}\,\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{12}\right)\mathbf{a}_3 &&= -x_{12} a\,\hat{\mathbf{x}} - y_{12} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{12}\right) c\,\hat{\mathbf{z}} && (4a) && \text{H X}\\
\mathbf{B}_{47} &= \left(\tfrac{1}{2} + x_{12}\right)\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_{12}\right) a\,\hat{\mathbf{x}} - y_{12} b\,\hat{\mathbf{y}} + z_{12} c\,\hat{\mathbf{z}} && (4a) && \text{H X}\\
\mathbf{B}_{48} &= \left(\tfrac{1}{2} - x_{12}\right)\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{12}\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_{12}\right) a\,\hat{\mathbf{x}} + y_{12} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{12}\right) c\,\hat{\mathbf{z}} && (4a) && \text{H X}\\
\mathbf{B}_{49} &= x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3 &&= x_{13} a\,\hat{\mathbf{x}} + y_{13} b\,\hat{\mathbf{y}} + z_{13} c\,\hat{\mathbf{z}} && (4a) && \text{H XI}\\
\mathbf{B}_{50} &= -x_{13}\,\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{13}\right)\mathbf{a}_3 &&= -x_{13} a\,\hat{\mathbf{x}} - y_{13} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{13}\right) c\,\hat{\mathbf{z}} && (4a) && \text{H XI}\\
\mathbf{B}_{51} &= \left(\tfrac{1}{2} + x_{13}\right)\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_{13}\right) a\,\hat{\mathbf{x}} - y_{13} b\,\hat{\mathbf{y}} + z_{13} c\,\hat{\mathbf{z}} && (4a) && \text{H XI}\\
\mathbf{B}_{52} &= \left(\tfrac{1}{2} - x_{13}\right)\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{13}\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_{13}\right) a\,\hat{\mathbf{x}} + y_{13} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{13}\right) c\,\hat{\mathbf{z}} && (4a) && \text{H XI}\\
\mathbf{B}_{53} &= x_{14}\,\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3 &&= x_{14} a\,\hat{\mathbf{x}} + y_{14} b\,\hat{\mathbf{y}} + z_{14} c\,\hat{\mathbf{z}} && (4a) && \text{H XII}
\end{aligned}
$$

$$
\begin{array}{lcccccc}
\mathbf{B}_{54} & = & -x_{14}\,\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{14}\right)\mathbf{a}_3 & = & -x_{14}a\,\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{14}\right)c\,\hat{\mathbf{z}} & (4a) & \text{H XII} \\
\mathbf{B}_{55} & = & \left(\frac{1}{2} + x_{14}\right)\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3 & = & \left(\frac{1}{2} + x_{14}\right)a\,\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} + z_{14}c\,\hat{\mathbf{z}} & (4a) & \text{H XII} \\
\mathbf{B}_{56} & = & \left(\frac{1}{2} - x_{14}\right)\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{14}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_{14}\right)a\,\hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{14}\right)c\,\hat{\mathbf{z}} & (4a) & \text{H XII} \\
\mathbf{B}_{57} & = & x_{15}\,\mathbf{a}_1 + y_{15}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 & = & x_{15}a\,\hat{\mathbf{x}} + y_{15}b\,\hat{\mathbf{y}} + z_{15}c\,\hat{\mathbf{z}} & (4a) & \text{H XIII} \\
\mathbf{B}_{58} & = & -x_{15}\,\mathbf{a}_1 - y_{15}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{15}\right)\mathbf{a}_3 & = & -x_{15}a\,\hat{\mathbf{x}} - y_{15}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{15}\right)c\,\hat{\mathbf{z}} & (4a) & \text{H XIII} \\
\mathbf{B}_{59} & = & \left(\frac{1}{2} + x_{15}\right)\mathbf{a}_1 - y_{15}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 & = & \left(\frac{1}{2} + x_{15}\right)a\,\hat{\mathbf{x}} - y_{15}b\,\hat{\mathbf{y}} + z_{15}c\,\hat{\mathbf{z}} & (4a) & \text{H XIII} \\
\mathbf{B}_{60} & = & \left(\frac{1}{2} - x_{15}\right)\mathbf{a}_1 + y_{15}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{15}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_{15}\right)a\,\hat{\mathbf{x}} + y_{15}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{15}\right)c\,\hat{\mathbf{z}} & (4a) & \text{H XIII} \\
\mathbf{B}_{61} & = & x_{16}\,\mathbf{a}_1 + y_{16}\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 & = & x_{16}a\,\hat{\mathbf{x}} + y_{16}b\,\hat{\mathbf{y}} + z_{16}c\,\hat{\mathbf{z}} & (4a) & \text{H XIV} \\
\mathbf{B}_{62} & = & -x_{16}\,\mathbf{a}_1 - y_{16}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{16}\right)\mathbf{a}_3 & = & -x_{16}a\,\hat{\mathbf{x}} - y_{16}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{16}\right)c\,\hat{\mathbf{z}} & (4a) & \text{H XIV} \\
\mathbf{B}_{63} & = & \left(\frac{1}{2} + x_{16}\right)\mathbf{a}_1 - y_{16}\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 & = & \left(\frac{1}{2} + x_{16}\right)a\,\hat{\mathbf{x}} - y_{16}b\,\hat{\mathbf{y}} + z_{16}c\,\hat{\mathbf{z}} & (4a) & \text{H XIV} \\
\mathbf{B}_{64} & = & \left(\frac{1}{2} - x_{16}\right)\mathbf{a}_1 + y_{16}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{16}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_{16}\right)a\,\hat{\mathbf{x}} + y_{16}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{16}\right)c\,\hat{\mathbf{z}} & (4a) & \text{H XIV} \\
\mathbf{B}_{65} & = & x_{17}\,\mathbf{a}_1 + y_{17}\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 & = & x_{17}a\,\hat{\mathbf{x}} + y_{17}b\,\hat{\mathbf{y}} + z_{17}c\,\hat{\mathbf{z}} & (4a) & \text{H XV} \\
\mathbf{B}_{66} & = & -x_{17}\,\mathbf{a}_1 - y_{17}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{17}\right)\mathbf{a}_3 & = & -x_{17}a\,\hat{\mathbf{x}} - y_{17}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{17}\right)c\,\hat{\mathbf{z}} & (4a) & \text{H XV} \\
\mathbf{B}_{67} & = & \left(\frac{1}{2} + x_{17}\right)\mathbf{a}_1 - y_{17}\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 & = & \left(\frac{1}{2} + x_{17}\right)a\,\hat{\mathbf{x}} - y_{17}b\,\hat{\mathbf{y}} + z_{17}c\,\hat{\mathbf{z}} & (4a) & \text{H XV} \\
\mathbf{B}_{68} & = & \left(\frac{1}{2} - x_{17}\right)\mathbf{a}_1 + y_{17}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{17}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_{17}\right)a\,\hat{\mathbf{x}} + y_{17}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{17}\right)c\,\hat{\mathbf{z}} & (4a) & \text{H XV} \\
\mathbf{B}_{69} & = & x_{18}\,\mathbf{a}_1 + y_{18}\,\mathbf{a}_2 + z_{18}\,\mathbf{a}_3 & = & x_{18}a\,\hat{\mathbf{x}} + y_{18}b\,\hat{\mathbf{y}} + z_{18}c\,\hat{\mathbf{z}} & (4a) & \text{H XVI} \\
\mathbf{B}_{70} & = & -x_{18}\,\mathbf{a}_1 - y_{18}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{18}\right)\mathbf{a}_3 & = & -x_{18}a\,\hat{\mathbf{x}} - y_{18}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{18}\right)c\,\hat{\mathbf{z}} & (4a) & \text{H XVI} \\
\mathbf{B}_{71} & = & \left(\frac{1}{2} + x_{18}\right)\mathbf{a}_1 - y_{18}\,\mathbf{a}_2 + z_{18}\,\mathbf{a}_3 & = & \left(\frac{1}{2} + x_{18}\right)a\,\hat{\mathbf{x}} - y_{18}b\,\hat{\mathbf{y}} + z_{18}c\,\hat{\mathbf{z}} & (4a) & \text{H XVI} \\
\mathbf{B}_{72} & = & \left(\frac{1}{2} - x_{18}\right)\mathbf{a}_1 + y_{18}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{18}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_{18}\right)a\,\hat{\mathbf{x}} + y_{18}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{18}\right)c\,\hat{\mathbf{z}} & (4a) & \text{H XVI} \\
\mathbf{B}_{73} & = & x_{19}\,\mathbf{a}_1 + y_{19}\,\mathbf{a}_2 + z_{19}\,\mathbf{a}_3 & = & x_{19}a\,\hat{\mathbf{x}} + y_{19}b\,\hat{\mathbf{y}} + z_{19}c\,\hat{\mathbf{z}} & (4a) & \text{H XVII} \\
\mathbf{B}_{74} & = & -x_{19}\,\mathbf{a}_1 - y_{19}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{19}\right)\mathbf{a}_3 & = & -x_{19}a\,\hat{\mathbf{x}} - y_{19}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{19}\right)c\,\hat{\mathbf{z}} & (4a) & \text{H XVII} \\
\mathbf{B}_{75} & = & \left(\frac{1}{2} + x_{19}\right)\mathbf{a}_1 - y_{19}\,\mathbf{a}_2 + z_{19}\,\mathbf{a}_3 & = & \left(\frac{1}{2} + x_{19}\right)a\,\hat{\mathbf{x}} - y_{19}b\,\hat{\mathbf{y}} + z_{19}c\,\hat{\mathbf{z}} & (4a) & \text{H XVII} \\
\mathbf{B}_{76} & = & \left(\frac{1}{2} - x_{19}\right)\mathbf{a}_1 + y_{19}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{19}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_{19}\right)a\,\hat{\mathbf{x}} + y_{19}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{19}\right)c\,\hat{\mathbf{z}} & (4a) & \text{H XVII} \\
\mathbf{B}_{77} & = & x_{20}\,\mathbf{a}_1 + y_{20}\,\mathbf{a}_2 + z_{20}\,\mathbf{a}_3 & = & x_{20}a\,\hat{\mathbf{x}} + y_{20}b\,\hat{\mathbf{y}} + z_{20}c\,\hat{\mathbf{z}} & (4a) & \text{H XVIII} \\
\mathbf{B}_{78} & = & -x_{20}\,\mathbf{a}_1 - y_{20}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{20}\right)\mathbf{a}_3 & = & -x_{20}a\,\hat{\mathbf{x}} - y_{20}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{20}\right)c\,\hat{\mathbf{z}} & (4a) & \text{H XVIII} \\
\mathbf{B}_{79} & = & \left(\frac{1}{2} + x_{20}\right)\mathbf{a}_1 - y_{20}\,\mathbf{a}_2 + z_{20}\,\mathbf{a}_3 & = & \left(\frac{1}{2} + x_{20}\right)a\,\hat{\mathbf{x}} - y_{20}b\,\hat{\mathbf{y}} + z_{20}c\,\hat{\mathbf{z}} & (4a) & \text{H XVIII} \\
\mathbf{B}_{80} & = & \left(\frac{1}{2} - x_{20}\right)\mathbf{a}_1 + y_{20}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{20}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_{20}\right)a\,\hat{\mathbf{x}} + y_{20}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{20}\right)c\,\hat{\mathbf{z}} & (4a) & \text{H XVIII} \\
\mathbf{B}_{81} & = & x_{21}\,\mathbf{a}_1 + y_{21}\,\mathbf{a}_2 + z_{21}\,\mathbf{a}_3 & = & x_{21}a\,\hat{\mathbf{x}} + y_{21}b\,\hat{\mathbf{y}} + z_{21}c\,\hat{\mathbf{z}} & (4a) & \text{H XIX} \\
\mathbf{B}_{82} & = & -x_{21}\,\mathbf{a}_1 - y_{21}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{21}\right)\mathbf{a}_3 & = & -x_{21}a\,\hat{\mathbf{x}} - y_{21}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{21}\right)c\,\hat{\mathbf{z}} & (4a) & \text{H XIX} \\
\mathbf{B}_{83} & = & \left(\frac{1}{2} + x_{21}\right)\mathbf{a}_1 - y_{21}\,\mathbf{a}_2 + z_{21}\,\mathbf{a}_3 & = & \left(\frac{1}{2} + x_{21}\right)a\,\hat{\mathbf{x}} - y_{21}b\,\hat{\mathbf{y}} + z_{21}c\,\hat{\mathbf{z}} & (4a) & \text{H XIX} \\
\mathbf{B}_{84} & = & \left(\frac{1}{2} - x_{21}\right)\mathbf{a}_1 + y_{21}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{21}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_{21}\right)a\,\hat{\mathbf{x}} + y_{21}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{21}\right)c\,\hat{\mathbf{z}} & (4a) & \text{H XIX} \\
\mathbf{B}_{85} & = & x_{22}\,\mathbf{a}_1 + y_{22}\,\mathbf{a}_2 + z_{22}\,\mathbf{a}_3 & = & x_{22}a\,\hat{\mathbf{x}} + y_{22}b\,\hat{\mathbf{y}} + z_{22}c\,\hat{\mathbf{z}} & (4a) & \text{H XX} \\
\mathbf{B}_{86} & = & -x_{22}\,\mathbf{a}_1 - y_{22}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{22}\right)\mathbf{a}_3 & = & -x_{22}a\,\hat{\mathbf{x}} - y_{22}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{22}\right)c\,\hat{\mathbf{z}} & (4a) & \text{H XX} \\
\mathbf{B}_{87} & = & \left(\frac{1}{2} + x_{22}\right)\mathbf{a}_1 - y_{22}\,\mathbf{a}_2 + z_{22}\,\mathbf{a}_3 & = & \left(\frac{1}{2} + x_{22}\right)a\,\hat{\mathbf{x}} - y_{22}b\,\hat{\mathbf{y}} + z_{22}c\,\hat{\mathbf{z}} & (4a) & \text{H XX} \\
\mathbf{B}_{88} & = & \left(\frac{1}{2} - x_{22}\right)\mathbf{a}_1 + y_{22}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{22}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_{22}\right)a\,\hat{\mathbf{x}} + y_{22}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{22}\right)c\,\hat{\mathbf{z}} & (4a) & \text{H XX} \\
\mathbf{B}_{89} & = & x_{23}\,\mathbf{a}_1 + y_{23}\,\mathbf{a}_2 + z_{23}\,\mathbf{a}_3 & = & x_{23}a\,\hat{\mathbf{x}} + y_{23}b\,\hat{\mathbf{y}} + z_{23}c\,\hat{\mathbf{z}} & (4a) & \text{H XXI}
\end{array}
$$

$$\mathbf{B}_{90} = -x_{23}\,\mathbf{a}_1 - y_{23}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{23}\right)\mathbf{a}_3 = -x_{23}a\,\hat{\mathbf{x}} - y_{23}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{23}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXI}$$

$$\mathbf{B}_{91} = \left(\tfrac{1}{2} + x_{23}\right)\mathbf{a}_1 - y_{23}\,\mathbf{a}_2 + z_{23}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{23}\right)a\,\hat{\mathbf{x}} - y_{23}b\,\hat{\mathbf{y}} + z_{23}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXI}$$

$$\mathbf{B}_{92} = \left(\tfrac{1}{2} - x_{23}\right)\mathbf{a}_1 + y_{23}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{23}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{23}\right)a\,\hat{\mathbf{x}} + y_{23}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{23}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXI}$$

$$\mathbf{B}_{93} = x_{24}\,\mathbf{a}_1 + y_{24}\,\mathbf{a}_2 + z_{24}\,\mathbf{a}_3 = x_{24}a\,\hat{\mathbf{x}} + y_{24}b\,\hat{\mathbf{y}} + z_{24}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXII}$$

$$\mathbf{B}_{94} = -x_{24}\,\mathbf{a}_1 - y_{24}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{24}\right)\mathbf{a}_3 = -x_{24}a\,\hat{\mathbf{x}} - y_{24}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{24}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXII}$$

$$\mathbf{B}_{95} = \left(\tfrac{1}{2} + x_{24}\right)\mathbf{a}_1 - y_{24}\,\mathbf{a}_2 + z_{24}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{24}\right)a\,\hat{\mathbf{x}} - y_{24}b\,\hat{\mathbf{y}} + z_{24}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXII}$$

$$\mathbf{B}_{96} = \left(\tfrac{1}{2} - x_{24}\right)\mathbf{a}_1 + y_{24}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{24}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{24}\right)a\,\hat{\mathbf{x}} + y_{24}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{24}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXII}$$

$$\mathbf{B}_{97} = x_{25}\,\mathbf{a}_1 + y_{25}\,\mathbf{a}_2 + z_{25}\,\mathbf{a}_3 = x_{25}a\,\hat{\mathbf{x}} + y_{25}b\,\hat{\mathbf{y}} + z_{25}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXIII}$$

$$\mathbf{B}_{98} = -x_{25}\,\mathbf{a}_1 - y_{25}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{25}\right)\mathbf{a}_3 = -x_{25}a\,\hat{\mathbf{x}} - y_{25}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{25}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXIII}$$

$$\mathbf{B}_{99} = \left(\tfrac{1}{2} + x_{25}\right)\mathbf{a}_1 - y_{25}\,\mathbf{a}_2 + z_{25}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{25}\right)a\,\hat{\mathbf{x}} - y_{25}b\,\hat{\mathbf{y}} + z_{25}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXIII}$$

$$\mathbf{B}_{100} = \left(\tfrac{1}{2} - x_{25}\right)\mathbf{a}_1 + y_{25}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{25}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{25}\right)a\,\hat{\mathbf{x}} + y_{25}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{25}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXIII}$$

$$\mathbf{B}_{101} = x_{26}\,\mathbf{a}_1 + y_{26}\,\mathbf{a}_2 + z_{26}\,\mathbf{a}_3 = x_{26}a\,\hat{\mathbf{x}} + y_{26}b\,\hat{\mathbf{y}} + z_{26}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXIV}$$

$$\mathbf{B}_{102} = -x_{26}\,\mathbf{a}_1 - y_{26}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{26}\right)\mathbf{a}_3 = -x_{26}a\,\hat{\mathbf{x}} - y_{26}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{26}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXIV}$$

$$\mathbf{B}_{103} = \left(\tfrac{1}{2} + x_{26}\right)\mathbf{a}_1 - y_{26}\,\mathbf{a}_2 + z_{26}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{26}\right)a\,\hat{\mathbf{x}} - y_{26}b\,\hat{\mathbf{y}} + z_{26}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXIV}$$

$$\mathbf{B}_{104} = \left(\tfrac{1}{2} - x_{26}\right)\mathbf{a}_1 + y_{26}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{26}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{26}\right)a\,\hat{\mathbf{x}} + y_{26}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{26}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXIV}$$

$$\mathbf{B}_{105} = x_{27}\,\mathbf{a}_1 + y_{27}\,\mathbf{a}_2 + z_{27}\,\mathbf{a}_3 = x_{27}a\,\hat{\mathbf{x}} + y_{27}b\,\hat{\mathbf{y}} + z_{27}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXV}$$

$$\mathbf{B}_{106} = -x_{27}\,\mathbf{a}_1 - y_{27}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{27}\right)\mathbf{a}_3 = -x_{27}a\,\hat{\mathbf{x}} - y_{27}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{27}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXV}$$

$$\mathbf{B}_{107} = \left(\tfrac{1}{2} + x_{27}\right)\mathbf{a}_1 - y_{27}\,\mathbf{a}_2 + z_{27}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{27}\right)a\,\hat{\mathbf{x}} - y_{27}b\,\hat{\mathbf{y}} + z_{27}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXV}$$

$$\mathbf{B}_{108} = \left(\tfrac{1}{2} - x_{27}\right)\mathbf{a}_1 + y_{27}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{27}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{27}\right)a\,\hat{\mathbf{x}} + y_{27}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{27}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXV}$$

$$\mathbf{B}_{109} = x_{28}\,\mathbf{a}_1 + y_{28}\,\mathbf{a}_2 + z_{28}\,\mathbf{a}_3 = x_{28}a\,\hat{\mathbf{x}} + y_{28}b\,\hat{\mathbf{y}} + z_{28}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXVI}$$

$$\mathbf{B}_{110} = -x_{28}\,\mathbf{a}_1 - y_{28}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{28}\right)\mathbf{a}_3 = -x_{28}a\,\hat{\mathbf{x}} - y_{28}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{28}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXVI}$$

$$\mathbf{B}_{111} = \left(\tfrac{1}{2} + x_{28}\right)\mathbf{a}_1 - y_{28}\,\mathbf{a}_2 + z_{28}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{28}\right)a\,\hat{\mathbf{x}} - y_{28}b\,\hat{\mathbf{y}} + z_{28}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXVI}$$

$$\mathbf{B}_{112} = \left(\tfrac{1}{2} - x_{28}\right)\mathbf{a}_1 + y_{28}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{28}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{28}\right)a\,\hat{\mathbf{x}} + y_{28}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{28}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXVI}$$

$$\mathbf{B}_{113} = x_{29}\,\mathbf{a}_1 + y_{29}\,\mathbf{a}_2 + z_{29}\,\mathbf{a}_3 = x_{29}a\,\hat{\mathbf{x}} + y_{29}b\,\hat{\mathbf{y}} + z_{29}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXVII}$$

$$\mathbf{B}_{114} = -x_{29}\,\mathbf{a}_1 - y_{29}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{29}\right)\mathbf{a}_3 = -x_{29}a\,\hat{\mathbf{x}} - y_{29}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{29}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXVII}$$

$$\mathbf{B}_{115} = \left(\tfrac{1}{2} + x_{29}\right)\mathbf{a}_1 - y_{29}\,\mathbf{a}_2 + z_{29}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{29}\right)a\,\hat{\mathbf{x}} - y_{29}b\,\hat{\mathbf{y}} + z_{29}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXVII}$$

$$\mathbf{B}_{116} = \left(\tfrac{1}{2} - x_{29}\right)\mathbf{a}_1 + y_{29}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{29}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{29}\right)a\,\hat{\mathbf{x}} + y_{29}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{29}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXVII}$$

$$\mathbf{B}_{117} = x_{30}\,\mathbf{a}_1 + y_{30}\,\mathbf{a}_2 + z_{30}\,\mathbf{a}_3 = x_{30}a\,\hat{\mathbf{x}} + y_{30}b\,\hat{\mathbf{y}} + z_{30}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXVIII}$$

$$\mathbf{B}_{118} = -x_{30}\,\mathbf{a}_1 - y_{30}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{30}\right)\mathbf{a}_3 = -x_{30}a\,\hat{\mathbf{x}} - y_{30}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{30}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXVIII}$$

$$\mathbf{B}_{119} = \left(\tfrac{1}{2} + x_{30}\right)\mathbf{a}_1 - y_{30}\,\mathbf{a}_2 + z_{30}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{30}\right)a\,\hat{\mathbf{x}} - y_{30}b\,\hat{\mathbf{y}} + z_{30}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXVIII}$$

$$\mathbf{B}_{120} = \left(\tfrac{1}{2} - x_{30}\right)\mathbf{a}_1 + y_{30}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{30}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{30}\right)a\,\hat{\mathbf{x}} + y_{30}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{30}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXVIII}$$

$$\mathbf{B}_{121} = x_{31}\,\mathbf{a}_1 + y_{31}\,\mathbf{a}_2 + z_{31}\,\mathbf{a}_3 = x_{31}a\,\hat{\mathbf{x}} + y_{31}b\,\hat{\mathbf{y}} + z_{31}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXIX}$$

$$\mathbf{B}_{122} = -x_{31}\,\mathbf{a}_1 - y_{31}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{31}\right)\mathbf{a}_3 = -x_{31}a\,\hat{\mathbf{x}} - y_{31}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{31}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXIX}$$

$$\mathbf{B}_{123} = \left(\tfrac{1}{2} + x_{31}\right)\mathbf{a}_1 - y_{31}\,\mathbf{a}_2 + z_{31}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{31}\right)a\,\hat{\mathbf{x}} - y_{31}b\,\hat{\mathbf{y}} + z_{31}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXIX}$$

$$\mathbf{B}_{124} = \left(\tfrac{1}{2} - x_{31}\right)\mathbf{a}_1 + y_{31}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{31}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{31}\right)a\,\hat{\mathbf{x}} + y_{31}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{31}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXIX}$$

$$\mathbf{B}_{125} = x_{32}\,\mathbf{a}_1 + y_{32}\,\mathbf{a}_2 + z_{32}\,\mathbf{a}_3 = x_{32}a\,\hat{\mathbf{x}} + y_{32}b\,\hat{\mathbf{y}} + z_{32}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXX}$$

$$\mathbf{B}_{126} = -x_{32}\,\mathbf{a}_1 - y_{32}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{32}\right)\mathbf{a}_3 = -x_{32}a\,\hat{\mathbf{x}} - y_{32}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{32}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXX}$$

$$\mathbf{B}_{127} = \left(\tfrac{1}{2} + x_{32}\right)\mathbf{a}_1 - y_{32}\,\mathbf{a}_2 + z_{32}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{32}\right)a\,\hat{\mathbf{x}} - y_{32}b\,\hat{\mathbf{y}} + z_{32}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXX}$$

$$\mathbf{B}_{128} = \left(\tfrac{1}{2} - x_{32}\right)\mathbf{a}_1 + y_{32}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{32}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{32}\right)a\,\hat{\mathbf{x}} + y_{32}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{32}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{H XXX}$$

$$\mathbf{B}_{129} = x_{33}\,\mathbf{a}_1 + y_{33}\,\mathbf{a}_2 + z_{33}\,\mathbf{a}_3 = x_{33}a\,\hat{\mathbf{x}} + y_{33}b\,\hat{\mathbf{y}} + z_{33}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{N}$$

$$\mathbf{B}_{130} = -x_{33}\,\mathbf{a}_1 - y_{33}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{33}\right)\mathbf{a}_3 = -x_{33}a\,\hat{\mathbf{x}} - y_{33}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{33}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{N}$$

$$\mathbf{B}_{131} = \left(\tfrac{1}{2} + x_{33}\right)\mathbf{a}_1 - y_{33}\,\mathbf{a}_2 + z_{33}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{33}\right)a\,\hat{\mathbf{x}} - y_{33}b\,\hat{\mathbf{y}} + z_{33}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{N}$$

$$\mathbf{B}_{132} = \left(\tfrac{1}{2} - x_{33}\right)\mathbf{a}_1 + y_{33}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{33}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{33}\right)a\,\hat{\mathbf{x}} + y_{33}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{33}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{N}$$

$$\mathbf{B}_{133} = x_{34}\,\mathbf{a}_1 + y_{34}\,\mathbf{a}_2 + z_{34}\,\mathbf{a}_3 = x_{34}a\,\hat{\mathbf{x}} + y_{34}b\,\hat{\mathbf{y}} + z_{34}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O I}$$

$$\mathbf{B}_{134} = -x_{34}\,\mathbf{a}_1 - y_{34}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{34}\right)\mathbf{a}_3 = -x_{34}a\,\hat{\mathbf{x}} - y_{34}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{34}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O I}$$

$$\mathbf{B}_{135} = \left(\tfrac{1}{2} + x_{34}\right)\mathbf{a}_1 - y_{34}\,\mathbf{a}_2 + z_{34}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{34}\right)a\,\hat{\mathbf{x}} - y_{34}b\,\hat{\mathbf{y}} + z_{34}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O I}$$

$$\mathbf{B}_{136} = \left(\tfrac{1}{2} - x_{34}\right)\mathbf{a}_1 + y_{34}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{34}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{34}\right)a\,\hat{\mathbf{x}} + y_{34}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{34}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O I}$$

$$\mathbf{B}_{137} = x_{35}\,\mathbf{a}_1 + y_{35}\,\mathbf{a}_2 + z_{35}\,\mathbf{a}_3 = x_{35}a\,\hat{\mathbf{x}} + y_{35}b\,\hat{\mathbf{y}} + z_{35}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O II}$$

$$\mathbf{B}_{138} = -x_{35}\,\mathbf{a}_1 - y_{35}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{35}\right)\mathbf{a}_3 = -x_{35}a\,\hat{\mathbf{x}} - y_{35}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{35}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O II}$$

$$\mathbf{B}_{139} = \left(\tfrac{1}{2} + x_{35}\right)\mathbf{a}_1 - y_{35}\,\mathbf{a}_2 + z_{35}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{35}\right)a\,\hat{\mathbf{x}} - y_{35}b\,\hat{\mathbf{y}} + z_{35}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O II}$$

$$\mathbf{B}_{140} = \left(\tfrac{1}{2} - x_{35}\right)\mathbf{a}_1 + y_{35}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{35}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{35}\right)a\,\hat{\mathbf{x}} + y_{35}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{35}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O II}$$

$$\mathbf{B}_{141} = x_{36}\,\mathbf{a}_1 + y_{36}\,\mathbf{a}_2 + z_{36}\,\mathbf{a}_3 = x_{36}a\,\hat{\mathbf{x}} + y_{36}b\,\hat{\mathbf{y}} + z_{36}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O III}$$

$$\mathbf{B}_{142} = -x_{36}\,\mathbf{a}_1 - y_{36}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{36}\right)\mathbf{a}_3 = -x_{36}a\,\hat{\mathbf{x}} - y_{36}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{36}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O III}$$

$$\mathbf{B}_{143} = \left(\tfrac{1}{2} + x_{36}\right)\mathbf{a}_1 - y_{36}\,\mathbf{a}_2 + z_{36}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{36}\right)a\,\hat{\mathbf{x}} - y_{36}b\,\hat{\mathbf{y}} + z_{36}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O III}$$

$$\mathbf{B}_{144} = \left(\tfrac{1}{2} - x_{36}\right)\mathbf{a}_1 + y_{36}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{36}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{36}\right)a\,\hat{\mathbf{x}} + y_{36}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{36}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O III}$$

$$\mathbf{B}_{145} = x_{37}\,\mathbf{a}_1 + y_{37}\,\mathbf{a}_2 + z_{37}\,\mathbf{a}_3 = x_{37}a\,\hat{\mathbf{x}} + y_{37}b\,\hat{\mathbf{y}} + z_{37}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O IV}$$

$$\mathbf{B}_{146} = -x_{37}\,\mathbf{a}_1 - y_{37}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{37}\right)\mathbf{a}_3 = -x_{37}a\,\hat{\mathbf{x}} - y_{37}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{37}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O IV}$$

$$\mathbf{B}_{147} = \left(\tfrac{1}{2} + x_{37}\right)\mathbf{a}_1 - y_{37}\,\mathbf{a}_2 + z_{37}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{37}\right)a\,\hat{\mathbf{x}} - y_{37}b\,\hat{\mathbf{y}} + z_{37}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O IV}$$

$$\mathbf{B}_{148} = \left(\tfrac{1}{2} - x_{37}\right)\mathbf{a}_1 + y_{37}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{37}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{37}\right)a\,\hat{\mathbf{x}} + y_{37}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{37}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O IV}$$

$$\mathbf{B}_{149} = x_{38}\,\mathbf{a}_1 + y_{38}\,\mathbf{a}_2 + z_{38}\,\mathbf{a}_3 = x_{38}a\,\hat{\mathbf{x}} + y_{38}b\,\hat{\mathbf{y}} + z_{38}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O V}$$

$$\mathbf{B}_{150} = -x_{38}\,\mathbf{a}_1 - y_{38}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{38}\right)\mathbf{a}_3 = -x_{38}a\,\hat{\mathbf{x}} - y_{38}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{38}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O V}$$

$$\mathbf{B}_{151} = \left(\tfrac{1}{2} + x_{38}\right)\mathbf{a}_1 - y_{38}\,\mathbf{a}_2 + z_{38}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{38}\right)a\,\hat{\mathbf{x}} - y_{38}b\,\hat{\mathbf{y}} + z_{38}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O V}$$

$$\mathbf{B}_{152} = \left(\tfrac{1}{2} - x_{38}\right)\mathbf{a}_1 + y_{38}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{38}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{38}\right)a\,\hat{\mathbf{x}} + y_{38}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{38}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O V}$$

$$\mathbf{B}_{153} = x_{39}\,\mathbf{a}_1 + y_{39}\,\mathbf{a}_2 + z_{39}\,\mathbf{a}_3 = x_{39}a\,\hat{\mathbf{x}} + y_{39}b\,\hat{\mathbf{y}} + z_{39}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O VI}$$

$$\mathbf{B}_{154} = -x_{39}\,\mathbf{a}_1 - y_{39}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{39}\right)\mathbf{a}_3 = -x_{39}a\,\hat{\mathbf{x}} - y_{39}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{39}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O VI}$$

$$\mathbf{B}_{155} = \left(\tfrac{1}{2} + x_{39}\right)\mathbf{a}_1 - y_{39}\,\mathbf{a}_2 + z_{39}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{39}\right)a\,\hat{\mathbf{x}} - y_{39}b\,\hat{\mathbf{y}} + z_{39}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O VI}$$

$$\mathbf{B}_{156} = \left(\tfrac{1}{2} - x_{39}\right)\mathbf{a}_1 + y_{39}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{39}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{39}\right)a\,\hat{\mathbf{x}} + y_{39}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{39}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O VI}$$

$$\mathbf{B}_{157} = x_{40}\,\mathbf{a}_1 + y_{40}\,\mathbf{a}_2 + z_{40}\,\mathbf{a}_3 = x_{40}a\,\hat{\mathbf{x}} + y_{40}b\,\hat{\mathbf{y}} + z_{40}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O VII}$$

$$\mathbf{B}_{158} = -x_{40}\,\mathbf{a}_1 - y_{40}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{40}\right)\mathbf{a}_3 = -x_{40}a\,\hat{\mathbf{x}} - y_{40}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{40}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O VII}$$

$$\mathbf{B}_{159} = \left(\tfrac{1}{2} + x_{40}\right)\mathbf{a}_1 - y_{40}\,\mathbf{a}_2 + z_{40}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{40}\right)a\,\hat{\mathbf{x}} - y_{40}b\,\hat{\mathbf{y}} + z_{40}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O VII}$$

$$\mathbf{B}_{160} = \left(\tfrac{1}{2} - x_{40}\right)\mathbf{a}_1 + y_{40}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{40}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{40}\right)a\,\hat{\mathbf{x}} + y_{40}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{40}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O VII}$$

$$\mathbf{B}_{161} = x_{41}\,\mathbf{a}_1 + y_{41}\,\mathbf{a}_2 + z_{41}\,\mathbf{a}_3 = x_{41}a\,\hat{\mathbf{x}} + y_{41}b\,\hat{\mathbf{y}} + z_{41}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O VIII}$$

$$\mathbf{B}_{162} = -x_{41}\,\mathbf{a}_1 - y_{41}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{41}\right)\mathbf{a}_3 = -x_{41}a\,\hat{\mathbf{x}} - y_{41}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{41}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O VIII}$$

$$\mathbf{B}_{163} = \left(\tfrac{1}{2} + x_{41}\right)\mathbf{a}_1 - y_{41}\,\mathbf{a}_2 + z_{41}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{41}\right)a\,\hat{\mathbf{x}} - y_{41}b\,\hat{\mathbf{y}} + z_{41}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O VIII}$$

$$\mathbf{B}_{164} = \left(\tfrac{1}{2} - x_{41}\right)\mathbf{a}_1 + y_{41}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{41}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{41}\right)a\,\hat{\mathbf{x}} + y_{41}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{41}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O VIII}$$

$$\mathbf{B}_{165} = x_{42}\,\mathbf{a}_1 + y_{42}\,\mathbf{a}_2 + z_{42}\,\mathbf{a}_3 = x_{42}a\,\hat{\mathbf{x}} + y_{42}b\,\hat{\mathbf{y}} + z_{42}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O IX}$$

$$\mathbf{B}_{166} = -x_{42}\,\mathbf{a}_1 - y_{42}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{42}\right)\mathbf{a}_3 = -x_{42}a\,\hat{\mathbf{x}} - y_{42}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{42}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O IX}$$

$$\mathbf{B}_{167} = \left(\tfrac{1}{2} + x_{42}\right)\mathbf{a}_1 - y_{42}\,\mathbf{a}_2 + z_{42}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{42}\right)a\,\hat{\mathbf{x}} - y_{42}b\,\hat{\mathbf{y}} + z_{42}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O IX}$$

$$\mathbf{B}_{168} = \left(\tfrac{1}{2} - x_{42}\right)\mathbf{a}_1 + y_{42}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{42}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{42}\right)a\,\hat{\mathbf{x}} + y_{42}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{42}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O IX}$$

$$\mathbf{B}_{169} = x_{43}\,\mathbf{a}_1 + y_{43}\,\mathbf{a}_2 + z_{43}\,\mathbf{a}_3 = x_{43}a\,\hat{\mathbf{x}} + y_{43}b\,\hat{\mathbf{y}} + z_{43}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O X}$$

$$\mathbf{B}_{170} = -x_{43}\,\mathbf{a}_1 - y_{43}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{43}\right)\mathbf{a}_3 = -x_{43}a\,\hat{\mathbf{x}} - y_{43}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{43}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O X}$$

$$\mathbf{B}_{171} = \left(\tfrac{1}{2} + x_{43}\right)\mathbf{a}_1 - y_{43}\,\mathbf{a}_2 + z_{43}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{43}\right)a\,\hat{\mathbf{x}} - y_{43}b\,\hat{\mathbf{y}} + z_{43}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O X}$$

$$\mathbf{B}_{172} = \left(\tfrac{1}{2} - x_{43}\right)\mathbf{a}_1 + y_{43}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{43}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{43}\right)a\,\hat{\mathbf{x}} + y_{43}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{43}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O X}$$

$$\mathbf{B}_{173} = x_{44}\,\mathbf{a}_1 + y_{44}\,\mathbf{a}_2 + z_{44}\,\mathbf{a}_3 = x_{44}a\,\hat{\mathbf{x}} + y_{44}b\,\hat{\mathbf{y}} + z_{44}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XI}$$

$$\mathbf{B}_{174} = -x_{44}\,\mathbf{a}_1 - y_{44}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{44}\right)\mathbf{a}_3 = -x_{44}a\,\hat{\mathbf{x}} - y_{44}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{44}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XI}$$

$$\mathbf{B}_{175} = \left(\tfrac{1}{2} + x_{44}\right)\mathbf{a}_1 - y_{44}\,\mathbf{a}_2 + z_{44}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{44}\right)a\,\hat{\mathbf{x}} - y_{44}b\,\hat{\mathbf{y}} + z_{44}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XI}$$

$$\mathbf{B}_{176} = \left(\tfrac{1}{2} - x_{44}\right)\mathbf{a}_1 + y_{44}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{44}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{44}\right)a\,\hat{\mathbf{x}} + y_{44}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{44}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XI}$$

$$\mathbf{B}_{177} = x_{45}\,\mathbf{a}_1 + y_{45}\,\mathbf{a}_2 + z_{45}\,\mathbf{a}_3 = x_{45}a\,\hat{\mathbf{x}} + y_{45}b\,\hat{\mathbf{y}} + z_{45}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XII}$$

$$\mathbf{B}_{178} = -x_{45}\,\mathbf{a}_1 - y_{45}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{45}\right)\mathbf{a}_3 = -x_{45}a\,\hat{\mathbf{x}} - y_{45}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{45}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XII}$$

$$\mathbf{B}_{179} = \left(\tfrac{1}{2} + x_{45}\right)\mathbf{a}_1 - y_{45}\,\mathbf{a}_2 + z_{45}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{45}\right)a\,\hat{\mathbf{x}} - y_{45}b\,\hat{\mathbf{y}} + z_{45}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XII}$$

$$\mathbf{B}_{180} = \left(\tfrac{1}{2} - x_{45}\right)\mathbf{a}_1 + y_{45}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{45}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{45}\right)a\,\hat{\mathbf{x}} + y_{45}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{45}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XII}$$

$$\mathbf{B}_{181} = x_{46}\,\mathbf{a}_1 + y_{46}\,\mathbf{a}_2 + z_{46}\,\mathbf{a}_3 = x_{46}a\,\hat{\mathbf{x}} + y_{46}b\,\hat{\mathbf{y}} + z_{46}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XIII}$$

$$\mathbf{B}_{182} = -x_{46}\,\mathbf{a}_1 - y_{46}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{46}\right)\mathbf{a}_3 = -x_{46}a\,\hat{\mathbf{x}} - y_{46}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{46}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XIII}$$

$$\mathbf{B}_{183} = \left(\tfrac{1}{2} + x_{46}\right)\mathbf{a}_1 - y_{46}\,\mathbf{a}_2 + z_{46}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{46}\right)a\,\hat{\mathbf{x}} - y_{46}b\,\hat{\mathbf{y}} + z_{46}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XIII}$$

$$\mathbf{B}_{184} = \left(\tfrac{1}{2} - x_{46}\right)\mathbf{a}_1 + y_{46}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{46}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{46}\right)a\,\hat{\mathbf{x}} + y_{46}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{46}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XIII}$$

$$\mathbf{B}_{185} = x_{47}\,\mathbf{a}_1 + y_{47}\,\mathbf{a}_2 + z_{47}\,\mathbf{a}_3 = x_{47}a\,\hat{\mathbf{x}} + y_{47}b\,\hat{\mathbf{y}} + z_{47}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XIV}$$

$$\mathbf{B}_{186} = -x_{47}\,\mathbf{a}_1 - y_{47}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{47}\right)\mathbf{a}_3 = -x_{47}a\,\hat{\mathbf{x}} - y_{47}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{47}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XIV}$$

$$\mathbf{B}_{187} = \left(\tfrac{1}{2} + x_{47}\right)\mathbf{a}_1 - y_{47}\,\mathbf{a}_2 + z_{47}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{47}\right)a\,\hat{\mathbf{x}} - y_{47}b\,\hat{\mathbf{y}} + z_{47}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XIV}$$

$$\mathbf{B}_{188} = \left(\tfrac{1}{2} - x_{47}\right)\mathbf{a}_1 + y_{47}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{47}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{47}\right)a\,\hat{\mathbf{x}} + y_{47}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{47}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XIV}$$

$$\mathbf{B}_{189} = x_{48}\,\mathbf{a}_1 + y_{48}\,\mathbf{a}_2 + z_{48}\,\mathbf{a}_3 = x_{48}a\,\hat{\mathbf{x}} + y_{48}b\,\hat{\mathbf{y}} + z_{48}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XV}$$

$$\mathbf{B}_{190} = -x_{48}\,\mathbf{a}_1 - y_{48}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{48}\right)\mathbf{a}_3 = -x_{48}a\,\hat{\mathbf{x}} - y_{48}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{48}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XV}$$

$$\mathbf{B}_{191} = \left(\tfrac{1}{2} + x_{48}\right)\mathbf{a}_1 - y_{48}\,\mathbf{a}_2 + z_{48}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{48}\right)a\,\hat{\mathbf{x}} - y_{48}b\,\hat{\mathbf{y}} + z_{48}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XV}$$

$$\mathbf{B}_{192} = \left(\tfrac{1}{2} - x_{48}\right)\mathbf{a}_1 + y_{48}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{48}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{48}\right)a\,\hat{\mathbf{x}} + y_{48}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{48}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XV}$$

$$\mathbf{B}_{193} = x_{49}\,\mathbf{a}_1 + y_{49}\,\mathbf{a}_2 + z_{49}\,\mathbf{a}_3 = x_{49}a\,\hat{\mathbf{x}} + y_{49}b\,\hat{\mathbf{y}} + z_{49}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XVI}$$

$$\mathbf{B}_{194} = -x_{49}\,\mathbf{a}_1 - y_{49}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{49}\right)\mathbf{a}_3 = -x_{49}a\,\hat{\mathbf{x}} - y_{49}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{49}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XVI}$$

$$\mathbf{B}_{195} = \left(\tfrac{1}{2} + x_{49}\right)\mathbf{a}_1 - y_{49}\,\mathbf{a}_2 + z_{49}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{49}\right)a\,\hat{\mathbf{x}} - y_{49}b\,\hat{\mathbf{y}} + z_{49}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XVI}$$

$$\mathbf{B}_{196} = \left(\tfrac{1}{2} - x_{49}\right)\mathbf{a}_1 + y_{49}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{49}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{49}\right)a\,\hat{\mathbf{x}} + y_{49}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{49}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XVI}$$

$$\mathbf{B}_{197} = x_{50}\,\mathbf{a}_1 + y_{50}\,\mathbf{a}_2 + z_{50}\,\mathbf{a}_3 = x_{50}a\,\hat{\mathbf{x}} + y_{50}b\,\hat{\mathbf{y}} + z_{50}c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XVII}$$

$\mathbf{B}_{198}$	=	$-x_{50}\,\mathbf{a}_1 - y_{50}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{50}\right)\mathbf{a}_3$	=	$-x_{50}a\,\hat{\mathbf{x}} - y_{50}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{50}\right)c\,\hat{\mathbf{z}}$	(4*a*)	O XVII
$\mathbf{B}_{199}$	=	$\left(\frac{1}{2} + x_{50}\right)\mathbf{a}_1 - y_{50}\,\mathbf{a}_2 + z_{50}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{50}\right)a\,\hat{\mathbf{x}} - y_{50}b\,\hat{\mathbf{y}} + z_{50}c\,\hat{\mathbf{z}}$	(4*a*)	O XVII
$\mathbf{B}_{200}$	=	$\left(\frac{1}{2} - x_{50}\right)\mathbf{a}_1 + y_{50}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{50}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{50}\right)a\,\hat{\mathbf{x}} + y_{50}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{50}\right)c\,\hat{\mathbf{z}}$	(4*a*)	O XVII
$\mathbf{B}_{201}$	=	$x_{51}\,\mathbf{a}_1 + y_{51}\,\mathbf{a}_2 + z_{51}\,\mathbf{a}_3$	=	$x_{51}a\,\hat{\mathbf{x}} + y_{51}b\,\hat{\mathbf{y}} + z_{51}c\,\hat{\mathbf{z}}$	(4*a*)	O XVIII
$\mathbf{B}_{202}$	=	$-x_{51}\,\mathbf{a}_1 - y_{51}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{51}\right)\mathbf{a}_3$	=	$-x_{51}a\,\hat{\mathbf{x}} - y_{51}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{51}\right)c\,\hat{\mathbf{z}}$	(4*a*)	O XVIII
$\mathbf{B}_{203}$	=	$\left(\frac{1}{2} + x_{51}\right)\mathbf{a}_1 - y_{51}\,\mathbf{a}_2 + z_{51}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{51}\right)a\,\hat{\mathbf{x}} - y_{51}b\,\hat{\mathbf{y}} + z_{51}c\,\hat{\mathbf{z}}$	(4*a*)	O XVIII
$\mathbf{B}_{204}$	=	$\left(\frac{1}{2} - x_{51}\right)\mathbf{a}_1 + y_{51}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{51}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{51}\right)a\,\hat{\mathbf{x}} + y_{51}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{51}\right)c\,\hat{\mathbf{z}}$	(4*a*)	O XVIII
$\mathbf{B}_{205}$	=	$x_{52}\,\mathbf{a}_1 + y_{52}\,\mathbf{a}_2 + z_{52}\,\mathbf{a}_3$	=	$x_{52}a\,\hat{\mathbf{x}} + y_{52}b\,\hat{\mathbf{y}} + z_{52}c\,\hat{\mathbf{z}}$	(4*a*)	O XIX
$\mathbf{B}_{206}$	=	$-x_{52}\,\mathbf{a}_1 - y_{52}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{52}\right)\mathbf{a}_3$	=	$-x_{52}a\,\hat{\mathbf{x}} - y_{52}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{52}\right)c\,\hat{\mathbf{z}}$	(4*a*)	O XIX
$\mathbf{B}_{207}$	=	$\left(\frac{1}{2} + x_{52}\right)\mathbf{a}_1 - y_{52}\,\mathbf{a}_2 + z_{52}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{52}\right)a\,\hat{\mathbf{x}} - y_{52}b\,\hat{\mathbf{y}} + z_{52}c\,\hat{\mathbf{z}}$	(4*a*)	O XIX
$\mathbf{B}_{208}$	=	$\left(\frac{1}{2} - x_{52}\right)\mathbf{a}_1 + y_{52}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{52}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{52}\right)a\,\hat{\mathbf{x}} + y_{52}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{52}\right)c\,\hat{\mathbf{z}}$	(4*a*)	O XIX
$\mathbf{B}_{209}$	=	$x_{53}\,\mathbf{a}_1 + y_{53}\,\mathbf{a}_2 + z_{53}\,\mathbf{a}_3$	=	$x_{53}a\,\hat{\mathbf{x}} + y_{53}b\,\hat{\mathbf{y}} + z_{53}c\,\hat{\mathbf{z}}$	(4*a*)	O XX
$\mathbf{B}_{210}$	=	$-x_{53}\,\mathbf{a}_1 - y_{53}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{53}\right)\mathbf{a}_3$	=	$-x_{53}a\,\hat{\mathbf{x}} - y_{53}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{53}\right)c\,\hat{\mathbf{z}}$	(4*a*)	O XX
$\mathbf{B}_{211}$	=	$\left(\frac{1}{2} + x_{53}\right)\mathbf{a}_1 - y_{53}\,\mathbf{a}_2 + z_{53}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{53}\right)a\,\hat{\mathbf{x}} - y_{53}b\,\hat{\mathbf{y}} + z_{53}c\,\hat{\mathbf{z}}$	(4*a*)	O XX
$\mathbf{B}_{212}$	=	$\left(\frac{1}{2} - x_{53}\right)\mathbf{a}_1 + y_{53}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{53}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{53}\right)a\,\hat{\mathbf{x}} + y_{53}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{53}\right)c\,\hat{\mathbf{z}}$	(4*a*)	O XX
$\mathbf{B}_{213}$	=	$x_{54}\,\mathbf{a}_1 + y_{54}\,\mathbf{a}_2 + z_{54}\,\mathbf{a}_3$	=	$x_{54}a\,\hat{\mathbf{x}} + y_{54}b\,\hat{\mathbf{y}} + z_{54}c\,\hat{\mathbf{z}}$	(4*a*)	S I
$\mathbf{B}_{214}$	=	$-x_{54}\,\mathbf{a}_1 - y_{54}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{54}\right)\mathbf{a}_3$	=	$-x_{54}a\,\hat{\mathbf{x}} - y_{54}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{54}\right)c\,\hat{\mathbf{z}}$	(4*a*)	S I
$\mathbf{B}_{215}$	=	$\left(\frac{1}{2} + x_{54}\right)\mathbf{a}_1 - y_{54}\,\mathbf{a}_2 + z_{54}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{54}\right)a\,\hat{\mathbf{x}} - y_{54}b\,\hat{\mathbf{y}} + z_{54}c\,\hat{\mathbf{z}}$	(4*a*)	S I
$\mathbf{B}_{216}$	=	$\left(\frac{1}{2} - x_{54}\right)\mathbf{a}_1 + y_{54}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{54}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{54}\right)a\,\hat{\mathbf{x}} + y_{54}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{54}\right)c\,\hat{\mathbf{z}}$	(4*a*)	S I
$\mathbf{B}_{217}$	=	$x_{55}\,\mathbf{a}_1 + y_{55}\,\mathbf{a}_2 + z_{55}\,\mathbf{a}_3$	=	$x_{55}a\,\hat{\mathbf{x}} + y_{55}b\,\hat{\mathbf{y}} + z_{55}c\,\hat{\mathbf{z}}$	(4*a*)	S II
$\mathbf{B}_{218}$	=	$-x_{55}\,\mathbf{a}_1 - y_{55}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{55}\right)\mathbf{a}_3$	=	$-x_{55}a\,\hat{\mathbf{x}} - y_{55}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{55}\right)c\,\hat{\mathbf{z}}$	(4*a*)	S II
$\mathbf{B}_{219}$	=	$\left(\frac{1}{2} + x_{55}\right)\mathbf{a}_1 - y_{55}\,\mathbf{a}_2 + z_{55}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{55}\right)a\,\hat{\mathbf{x}} - y_{55}b\,\hat{\mathbf{y}} + z_{55}c\,\hat{\mathbf{z}}$	(4*a*)	S II
$\mathbf{B}_{220}$	=	$\left(\frac{1}{2} - x_{55}\right)\mathbf{a}_1 + y_{55}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{55}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{55}\right)a\,\hat{\mathbf{x}} + y_{55}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{55}\right)c\,\hat{\mathbf{z}}$	(4*a*)	S II

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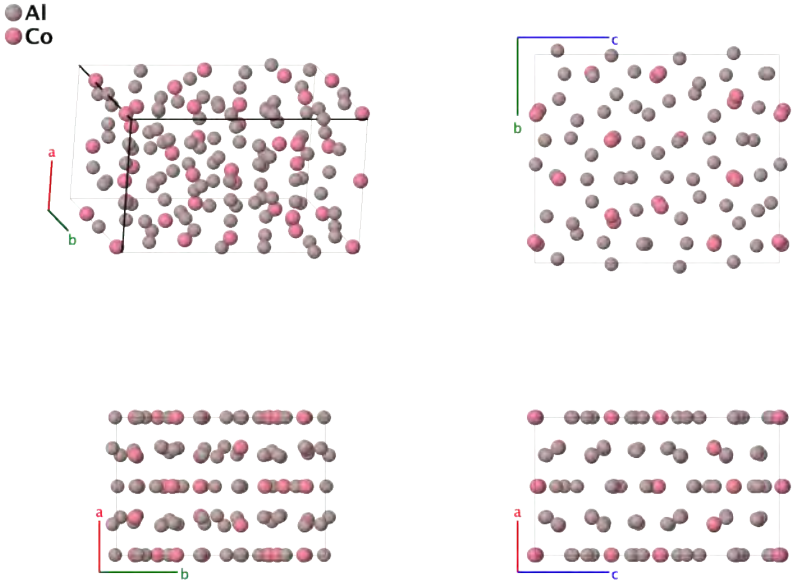
Geometry files:

- CIF: pp. 1582
- POSCAR: pp. 1583

Orthorhombic Co_4Al_{13} Structure: A13B4_oP102_31_17a11b_8a2b

http://aflow.org/prototype-encyclopedia/A13B4_oP102_31_17a11b_8a2b

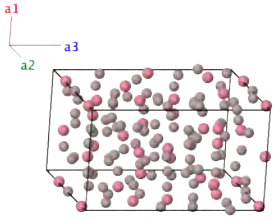


Prototype	:	$Al_{13}Co_4$
AFLOW prototype label	:	A13B4_oP102_31_17a11b_8a2b
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP102
Space group number	:	31
Space group symbol	:	$Pmn2_1$
AFLOW prototype command	:	`aflow --proto=A13B4_oP102_31_17a11b_8a2b --params=`$a, b/a, c/a, y_1, z_1, y_2, z_2, y_3, z_3, y_4, z_4, y_5, z_5, y_6, z_6, y_7, z_7, y_8, z_8, y_9, z_9, y_{10}, z_{10}, y_{11}, z_{11}, y_{12}, z_{12}, y_{13}, z_{13}, y_{14}, z_{14}, y_{15}, z_{15}, y_{16}, z_{16}, y_{17}, z_{17}, y_{18}, z_{18}, y_{19}, z_{19}, y_{20}, z_{20}, y_{21}, z_{21}, y_{22}, z_{22}, y_{23}, z_{23}, y_{24}, z_{24}, y_{25}, z_{25}, x_{26}, y_{26}, z_{26}, x_{27}, y_{27}, z_{27}, x_{28}, y_{28}, z_{28}, x_{29}, y_{29}, z_{29}, x_{30}, y_{30}, z_{30}, x_{31}, y_{31}, z_{31}, x_{32}, y_{32}, z_{32}, x_{33}, y_{33}, z_{33}, x_{34}, y_{34}, z_{34}, x_{35}, y_{35}, z_{35}, x_{36}, y_{36}, z_{36}, x_{37}, y_{37}, z_{37}, x_{38}, y_{38}, z_{38}$

- Space group $Pmn2_1$ #31 allows an arbitrary choice for the origin of the z-axis. We follow (Grin, 1994) and set the $z_{25} = 0$.
- If we allow a tolerance of 0.25 Å for AFLOW-SYM and 0.6 Å in the atomic positions for FINDSYM, the symmetry is classified as *Pnnm* #58.

Simple Orthorhombic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= a\,\hat{\mathbf{x}}\\ \mathbf{a}_2 &= b\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$y_1 b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	$(2a)$	Al I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\,\hat{\mathbf{z}}$	$(2a)$	Al I
$\mathbf{B}_3$	=	$y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(2a)$	Al II
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	$(2a)$	Al II
$\mathbf{B}_5$	=	$y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(2a)$	Al III
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	$(2a)$	Al III
$\mathbf{B}_7$	=	$y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(2a)$	Al IV
$\mathbf{B}_8$	=	$\frac{1}{2}\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	$(2a)$	Al IV
$\mathbf{B}_9$	=	$y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(2a)$	Al V
$\mathbf{B}_{10}$	=	$\frac{1}{2}\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	$(2a)$	Al V
$\mathbf{B}_{11}$	=	$y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(2a)$	Al VI
$\mathbf{B}_{12}$	=	$\frac{1}{2}\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\,\hat{\mathbf{z}}$	$(2a)$	Al VI
$\mathbf{B}_{13}$	=	$y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	$(2a)$	Al VII
$\mathbf{B}_{14}$	=	$\frac{1}{2}\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right)c\,\hat{\mathbf{z}}$	$(2a)$	Al VII
$\mathbf{B}_{15}$	=	$y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	$(2a)$	Al VIII
$\mathbf{B}_{16}$	=	$\frac{1}{2}\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right)c\,\hat{\mathbf{z}}$	$(2a)$	Al VIII
$\mathbf{B}_{17}$	=	$y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	$(2a)$	Al IX
$\mathbf{B}_{18}$	=	$\frac{1}{2}\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right)c\,\hat{\mathbf{z}}$	$(2a)$	Al IX
$\mathbf{B}_{19}$	=	$y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$y_{10} b\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}}$	$(2a)$	Al X
$\mathbf{B}_{20}$	=	$\frac{1}{2}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{10}\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{10}\right)c\,\hat{\mathbf{z}}$	$(2a)$	Al X
$\mathbf{B}_{21}$	=	$y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$y_{11} b\,\hat{\mathbf{y}} + z_{11} c\,\hat{\mathbf{z}}$	$(2a)$	Al XI
$\mathbf{B}_{22}$	=	$\frac{1}{2}\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{11}\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{11}\right)c\,\hat{\mathbf{z}}$	$(2a)$	Al XI
$\mathbf{B}_{23}$	=	$y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3$	=	$y_{12} b\,\hat{\mathbf{y}} + z_{12} c\,\hat{\mathbf{z}}$	$(2a)$	Al XII
$\mathbf{B}_{24}$	=	$\frac{1}{2}\,\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{12}\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - y_{12} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{12}\right)c\,\hat{\mathbf{z}}$	$(2a)$	Al XII
$\mathbf{B}_{25}$	=	$y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3$	=	$y_{13} b\,\hat{\mathbf{y}} + z_{13} c\,\hat{\mathbf{z}}$	$(2a)$	Al XIII
$\mathbf{B}_{26}$	=	$\frac{1}{2}\,\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{13}\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - y_{13} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{13}\right)c\,\hat{\mathbf{z}}$	$(2a)$	Al XIII
$\mathbf{B}_{27}$	=	$y_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3$	=	$y_{14} b\,\hat{\mathbf{y}} + z_{14} c\,\hat{\mathbf{z}}$	$(2a)$	Al XIV

$$
\begin{aligned}
\mathbf{B}_{28} &= \tfrac{1}{2}\,\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{14}\right)\mathbf{a}_3 &&= \tfrac{1}{2}a\,\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{14}\right)c\,\hat{\mathbf{z}} && (2a) && \text{Al XIV}\\
\mathbf{B}_{29} &= y_{15}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 &&= y_{15}b\,\hat{\mathbf{y}} + z_{15}c\,\hat{\mathbf{z}} && (2a) && \text{Al XV}\\
\mathbf{B}_{30} &= \tfrac{1}{2}\,\mathbf{a}_1 - y_{15}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{15}\right)\mathbf{a}_3 &&= \tfrac{1}{2}a\,\hat{\mathbf{x}} - y_{15}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{15}\right)c\,\hat{\mathbf{z}} && (2a) && \text{Al XV}\\
\mathbf{B}_{31} &= y_{16}\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 &&= y_{16}b\,\hat{\mathbf{y}} + z_{16}c\,\hat{\mathbf{z}} && (2a) && \text{Al XVI}\\
\mathbf{B}_{32} &= \tfrac{1}{2}\,\mathbf{a}_1 - y_{16}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{16}\right)\mathbf{a}_3 &&= \tfrac{1}{2}a\,\hat{\mathbf{x}} - y_{16}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{16}\right)c\,\hat{\mathbf{z}} && (2a) && \text{Al XVI}\\
\mathbf{B}_{33} &= y_{17}\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 &&= y_{17}b\,\hat{\mathbf{y}} + z_{17}c\,\hat{\mathbf{z}} && (2a) && \text{Al XVII}\\
\mathbf{B}_{34} &= \tfrac{1}{2}\,\mathbf{a}_1 - y_{17}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{17}\right)\mathbf{a}_3 &&= \tfrac{1}{2}a\,\hat{\mathbf{x}} - y_{17}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{17}\right)c\,\hat{\mathbf{z}} && (2a) && \text{Al XVII}\\
\mathbf{B}_{35} &= y_{18}\,\mathbf{a}_2 + z_{18}\,\mathbf{a}_3 &&= y_{18}b\,\hat{\mathbf{y}} + z_{18}c\,\hat{\mathbf{z}} && (2a) && \text{Co I}\\
\mathbf{B}_{36} &= \tfrac{1}{2}\,\mathbf{a}_1 - y_{18}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{18}\right)\mathbf{a}_3 &&= \tfrac{1}{2}a\,\hat{\mathbf{x}} - y_{18}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{18}\right)c\,\hat{\mathbf{z}} && (2a) && \text{Co I}\\
\mathbf{B}_{37} &= y_{19}\,\mathbf{a}_2 + z_{19}\,\mathbf{a}_3 &&= y_{19}b\,\hat{\mathbf{y}} + z_{19}c\,\hat{\mathbf{z}} && (2a) && \text{Co II}\\
\mathbf{B}_{38} &= \tfrac{1}{2}\,\mathbf{a}_1 - y_{19}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{19}\right)\mathbf{a}_3 &&= \tfrac{1}{2}a\,\hat{\mathbf{x}} - y_{19}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{19}\right)c\,\hat{\mathbf{z}} && (2a) && \text{Co II}\\
\mathbf{B}_{39} &= y_{20}\,\mathbf{a}_2 + z_{20}\,\mathbf{a}_3 &&= y_{20}b\,\hat{\mathbf{y}} + z_{20}c\,\hat{\mathbf{z}} && (2a) && \text{Co III}\\
\mathbf{B}_{40} &= \tfrac{1}{2}\,\mathbf{a}_1 - y_{20}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{20}\right)\mathbf{a}_3 &&= \tfrac{1}{2}a\,\hat{\mathbf{x}} - y_{20}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{20}\right)c\,\hat{\mathbf{z}} && (2a) && \text{Co III}\\
\mathbf{B}_{41} &= y_{21}\,\mathbf{a}_2 + z_{21}\,\mathbf{a}_3 &&= y_{21}b\,\hat{\mathbf{y}} + z_{21}c\,\hat{\mathbf{z}} && (2a) && \text{Co IV}\\
\mathbf{B}_{42} &= \tfrac{1}{2}\,\mathbf{a}_1 - y_{21}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{21}\right)\mathbf{a}_3 &&= \tfrac{1}{2}a\,\hat{\mathbf{x}} - y_{21}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{21}\right)c\,\hat{\mathbf{z}} && (2a) && \text{Co IV}\\
\mathbf{B}_{43} &= y_{22}\,\mathbf{a}_2 + z_{22}\,\mathbf{a}_3 &&= y_{22}b\,\hat{\mathbf{y}} + z_{22}c\,\hat{\mathbf{z}} && (2a) && \text{Co V}\\
\mathbf{B}_{44} &= \tfrac{1}{2}\,\mathbf{a}_1 - y_{22}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{22}\right)\mathbf{a}_3 &&= \tfrac{1}{2}a\,\hat{\mathbf{x}} - y_{22}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{22}\right)c\,\hat{\mathbf{z}} && (2a) && \text{Co V}\\
\mathbf{B}_{45} &= y_{23}\,\mathbf{a}_2 + z_{23}\,\mathbf{a}_3 &&= y_{23}b\,\hat{\mathbf{y}} + z_{23}c\,\hat{\mathbf{z}} && (2a) && \text{Co VI}\\
\mathbf{B}_{46} &= \tfrac{1}{2}\,\mathbf{a}_1 - y_{23}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{23}\right)\mathbf{a}_3 &&= \tfrac{1}{2}a\,\hat{\mathbf{x}} - y_{23}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{23}\right)c\,\hat{\mathbf{z}} && (2a) && \text{Co VI}\\
\mathbf{B}_{47} &= y_{24}\,\mathbf{a}_2 + z_{24}\,\mathbf{a}_3 &&= y_{24}b\,\hat{\mathbf{y}} + z_{24}c\,\hat{\mathbf{z}} && (2a) && \text{Co VII}\\
\mathbf{B}_{48} &= \tfrac{1}{2}\,\mathbf{a}_1 - y_{24}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{24}\right)\mathbf{a}_3 &&= \tfrac{1}{2}a\,\hat{\mathbf{x}} - y_{24}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{24}\right)c\,\hat{\mathbf{z}} && (2a) && \text{Co VII}\\
\mathbf{B}_{49} &= y_{25}\,\mathbf{a}_2 + z_{25}\,\mathbf{a}_3 &&= y_{25}b\,\hat{\mathbf{y}} + z_{25}c\,\hat{\mathbf{z}} && (2a) && \text{Co VIII}\\
\mathbf{B}_{50} &= \tfrac{1}{2}\,\mathbf{a}_1 - y_{25}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{25}\right)\mathbf{a}_3 &&= \tfrac{1}{2}a\,\hat{\mathbf{x}} - y_{25}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{25}\right)c\,\hat{\mathbf{z}} && (2a) && \text{Co VIII}\\
\mathbf{B}_{51} &= x_{26}\,\mathbf{a}_1 + y_{26}\,\mathbf{a}_2 + z_{26}\,\mathbf{a}_3 &&= x_{26}a\,\hat{\mathbf{x}} + y_{26}b\,\hat{\mathbf{y}} + z_{26}c\,\hat{\mathbf{z}} && (4b) && \text{Al XVIII}\\
\mathbf{B}_{52} &= \left(\tfrac{1}{2} - x_{26}\right)\mathbf{a}_1 - y_{26}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{26}\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_{26}\right)a\,\hat{\mathbf{x}} - y_{26}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{26}\right)c\,\hat{\mathbf{z}} && (4b) && \text{Al XVIII}\\
\mathbf{B}_{53} &= \left(\tfrac{1}{2} + x_{26}\right)\mathbf{a}_1 - y_{26}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{26}\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_{26}\right)a\,\hat{\mathbf{x}} - y_{26}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{26}\right)c\,\hat{\mathbf{z}} && (4b) && \text{Al XVIII}\\
\mathbf{B}_{54} &= -x_{26}\,\mathbf{a}_1 + y_{26}\,\mathbf{a}_2 + z_{26}\,\mathbf{a}_3 &&= -x_{26}a\,\hat{\mathbf{x}} + y_{26}b\,\hat{\mathbf{y}} + z_{26}c\,\hat{\mathbf{z}} && (4b) && \text{Al XVIII}\\
\mathbf{B}_{55} &= x_{27}\,\mathbf{a}_1 + y_{27}\,\mathbf{a}_2 + z_{27}\,\mathbf{a}_3 &&= x_{27}a\,\hat{\mathbf{x}} + y_{27}b\,\hat{\mathbf{y}} + z_{27}c\,\hat{\mathbf{z}} && (4b) && \text{Al XIX}\\
\mathbf{B}_{56} &= \left(\tfrac{1}{2} - x_{27}\right)\mathbf{a}_1 - y_{27}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{27}\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_{27}\right)a\,\hat{\mathbf{x}} - y_{27}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{27}\right)c\,\hat{\mathbf{z}} && (4b) && \text{Al XIX}\\
\mathbf{B}_{57} &= \left(\tfrac{1}{2} + x_{27}\right)\mathbf{a}_1 - y_{27}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{27}\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_{27}\right)a\,\hat{\mathbf{x}} - y_{27}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{27}\right)c\,\hat{\mathbf{z}} && (4b) && \text{Al XIX}\\
\mathbf{B}_{58} &= -x_{27}\,\mathbf{a}_1 + y_{27}\,\mathbf{a}_2 + z_{27}\,\mathbf{a}_3 &&= -x_{27}a\,\hat{\mathbf{x}} + y_{27}b\,\hat{\mathbf{y}} + z_{27}c\,\hat{\mathbf{z}} && (4b) && \text{Al XIX}\\
\mathbf{B}_{59} &= x_{28}\,\mathbf{a}_1 + y_{28}\,\mathbf{a}_2 + z_{28}\,\mathbf{a}_3 &&= x_{28}a\,\hat{\mathbf{x}} + y_{28}b\,\hat{\mathbf{y}} + z_{28}c\,\hat{\mathbf{z}} && (4b) && \text{Al XX}\\
\mathbf{B}_{60} &= \left(\tfrac{1}{2} - x_{28}\right)\mathbf{a}_1 - y_{28}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{28}\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_{28}\right)a\,\hat{\mathbf{x}} - y_{28}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{28}\right)c\,\hat{\mathbf{z}} && (4b) && \text{Al XX}\\
\mathbf{B}_{61} &= \left(\tfrac{1}{2} + x_{28}\right)\mathbf{a}_1 - y_{28}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{28}\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_{28}\right)a\,\hat{\mathbf{x}} - y_{28}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{28}\right)c\,\hat{\mathbf{z}} && (4b) && \text{Al XX}\\
\mathbf{B}_{62} &= -x_{28}\,\mathbf{a}_1 + y_{28}\,\mathbf{a}_2 + z_{28}\,\mathbf{a}_3 &&= -x_{28}a\,\hat{\mathbf{x}} + y_{28}b\,\hat{\mathbf{y}} + z_{28}c\,\hat{\mathbf{z}} && (4b) && \text{Al XX}\\
\mathbf{B}_{63} &= x_{29}\,\mathbf{a}_1 + y_{29}\,\mathbf{a}_2 + z_{29}\,\mathbf{a}_3 &&= x_{29}a\,\hat{\mathbf{x}} + y_{29}b\,\hat{\mathbf{y}} + z_{29}c\,\hat{\mathbf{z}} && (4b) && \text{Al XXI}
\end{aligned}
$$

$$\mathbf{B}_{64} = \left(\tfrac{1}{2} - x_{29}\right)\mathbf{a}_1 - y_{29}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{29}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{29}\right)a\,\hat{\mathbf{x}} - y_{29}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{29}\right)c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXI}$$

$$\mathbf{B}_{65} = \left(\tfrac{1}{2} + x_{29}\right)\mathbf{a}_1 - y_{29}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{29}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{29}\right)a\,\hat{\mathbf{x}} - y_{29}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{29}\right)c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXI}$$

$$\mathbf{B}_{66} = -x_{29}\,\mathbf{a}_1 + y_{29}\,\mathbf{a}_2 + z_{29}\,\mathbf{a}_3 = -x_{29}a\,\hat{\mathbf{x}} + y_{29}b\,\hat{\mathbf{y}} + z_{29}c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXI}$$

$$\mathbf{B}_{67} = x_{30}\,\mathbf{a}_1 + y_{30}\,\mathbf{a}_2 + z_{30}\,\mathbf{a}_3 = x_{30}a\,\hat{\mathbf{x}} + y_{30}b\,\hat{\mathbf{y}} + z_{30}c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXII}$$

$$\mathbf{B}_{68} = \left(\tfrac{1}{2} - x_{30}\right)\mathbf{a}_1 - y_{30}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{30}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{30}\right)a\,\hat{\mathbf{x}} - y_{30}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{30}\right)c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXII}$$

$$\mathbf{B}_{69} = \left(\tfrac{1}{2} + x_{30}\right)\mathbf{a}_1 - y_{30}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{30}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{30}\right)a\,\hat{\mathbf{x}} - y_{30}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{30}\right)c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXII}$$

$$\mathbf{B}_{70} = -x_{30}\,\mathbf{a}_1 + y_{30}\,\mathbf{a}_2 + z_{30}\,\mathbf{a}_3 = -x_{30}a\,\hat{\mathbf{x}} + y_{30}b\,\hat{\mathbf{y}} + z_{30}c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXII}$$

$$\mathbf{B}_{71} = x_{31}\,\mathbf{a}_1 + y_{31}\,\mathbf{a}_2 + z_{31}\,\mathbf{a}_3 = x_{31}a\,\hat{\mathbf{x}} + y_{31}b\,\hat{\mathbf{y}} + z_{31}c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXIII}$$

$$\mathbf{B}_{72} = \left(\tfrac{1}{2} - x_{31}\right)\mathbf{a}_1 - y_{31}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{31}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{31}\right)a\,\hat{\mathbf{x}} - y_{31}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{31}\right)c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXIII}$$

$$\mathbf{B}_{73} = \left(\tfrac{1}{2} + x_{31}\right)\mathbf{a}_1 - y_{31}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{31}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{31}\right)a\,\hat{\mathbf{x}} - y_{31}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{31}\right)c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXIII}$$

$$\mathbf{B}_{74} = -x_{31}\,\mathbf{a}_1 + y_{31}\,\mathbf{a}_2 + z_{31}\,\mathbf{a}_3 = -x_{31}a\,\hat{\mathbf{x}} + y_{31}b\,\hat{\mathbf{y}} + z_{31}c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXIII}$$

$$\mathbf{B}_{75} = x_{32}\,\mathbf{a}_1 + y_{32}\,\mathbf{a}_2 + z_{32}\,\mathbf{a}_3 = x_{32}a\,\hat{\mathbf{x}} + y_{32}b\,\hat{\mathbf{y}} + z_{32}c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXIV}$$

$$\mathbf{B}_{76} = \left(\tfrac{1}{2} - x_{32}\right)\mathbf{a}_1 - y_{32}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{32}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{32}\right)a\,\hat{\mathbf{x}} - y_{32}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{32}\right)c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXIV}$$

$$\mathbf{B}_{77} = \left(\tfrac{1}{2} + x_{32}\right)\mathbf{a}_1 - y_{32}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{32}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{32}\right)a\,\hat{\mathbf{x}} - y_{32}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{32}\right)c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXIV}$$

$$\mathbf{B}_{78} = -x_{32}\,\mathbf{a}_1 + y_{32}\,\mathbf{a}_2 + z_{32}\,\mathbf{a}_3 = -x_{32}a\,\hat{\mathbf{x}} + y_{32}b\,\hat{\mathbf{y}} + z_{32}c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXIV}$$

$$\mathbf{B}_{79} = x_{33}\,\mathbf{a}_1 + y_{33}\,\mathbf{a}_2 + z_{33}\,\mathbf{a}_3 = x_{33}a\,\hat{\mathbf{x}} + y_{33}b\,\hat{\mathbf{y}} + z_{33}c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXV}$$

$$\mathbf{B}_{80} = \left(\tfrac{1}{2} - x_{33}\right)\mathbf{a}_1 - y_{33}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{33}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{33}\right)a\,\hat{\mathbf{x}} - y_{33}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{33}\right)c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXV}$$

$$\mathbf{B}_{81} = \left(\tfrac{1}{2} + x_{33}\right)\mathbf{a}_1 - y_{33}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{33}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{33}\right)a\,\hat{\mathbf{x}} - y_{33}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{33}\right)c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXV}$$

$$\mathbf{B}_{82} = -x_{33}\,\mathbf{a}_1 + y_{33}\,\mathbf{a}_2 + z_{33}\,\mathbf{a}_3 = -x_{33}a\,\hat{\mathbf{x}} + y_{33}b\,\hat{\mathbf{y}} + z_{33}c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXV}$$

$$\mathbf{B}_{83} = x_{34}\,\mathbf{a}_1 + y_{34}\,\mathbf{a}_2 + z_{34}\,\mathbf{a}_3 = x_{34}a\,\hat{\mathbf{x}} + y_{34}b\,\hat{\mathbf{y}} + z_{34}c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXVI}$$

$$\mathbf{B}_{84} = \left(\tfrac{1}{2} - x_{34}\right)\mathbf{a}_1 - y_{34}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{34}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{34}\right)a\,\hat{\mathbf{x}} - y_{34}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{34}\right)c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXVI}$$

$$\mathbf{B}_{85} = \left(\tfrac{1}{2} + x_{34}\right)\mathbf{a}_1 - y_{34}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{34}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{34}\right)a\,\hat{\mathbf{x}} - y_{34}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{34}\right)c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXVI}$$

$$\mathbf{B}_{86} = -x_{34}\,\mathbf{a}_1 + y_{34}\,\mathbf{a}_2 + z_{34}\,\mathbf{a}_3 = -x_{34}a\,\hat{\mathbf{x}} + y_{34}b\,\hat{\mathbf{y}} + z_{34}c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXVI}$$

$$\mathbf{B}_{87} = x_{35}\,\mathbf{a}_1 + y_{35}\,\mathbf{a}_2 + z_{35}\,\mathbf{a}_3 = x_{35}a\,\hat{\mathbf{x}} + y_{35}b\,\hat{\mathbf{y}} + z_{35}c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXVII}$$

$$\mathbf{B}_{88} = \left(\tfrac{1}{2} - x_{35}\right)\mathbf{a}_1 - y_{35}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{35}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{35}\right)a\,\hat{\mathbf{x}} - y_{35}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{35}\right)c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXVII}$$

$$\mathbf{B}_{89} = \left(\tfrac{1}{2} + x_{35}\right)\mathbf{a}_1 - y_{35}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{35}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{35}\right)a\,\hat{\mathbf{x}} - y_{35}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{35}\right)c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXVII}$$

$$\mathbf{B}_{90} = -x_{35}\,\mathbf{a}_1 + y_{35}\,\mathbf{a}_2 + z_{35}\,\mathbf{a}_3 = -x_{35}a\,\hat{\mathbf{x}} + y_{35}b\,\hat{\mathbf{y}} + z_{35}c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXVII}$$

$$\mathbf{B}_{91} = x_{36}\,\mathbf{a}_1 + y_{36}\,\mathbf{a}_2 + z_{36}\,\mathbf{a}_3 = x_{36}a\,\hat{\mathbf{x}} + y_{36}b\,\hat{\mathbf{y}} + z_{36}c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXVIII}$$

$$\mathbf{B}_{92} = \left(\tfrac{1}{2} - x_{36}\right)\mathbf{a}_1 - y_{36}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{36}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{36}\right)a\,\hat{\mathbf{x}} - y_{36}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{36}\right)c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXVIII}$$

$$\mathbf{B}_{93} = \left(\tfrac{1}{2} + x_{36}\right)\mathbf{a}_1 - y_{36}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{36}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{36}\right)a\,\hat{\mathbf{x}} - y_{36}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{36}\right)c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXVIII}$$

$$\mathbf{B}_{94} = -x_{36}\,\mathbf{a}_1 + y_{36}\,\mathbf{a}_2 + z_{36}\,\mathbf{a}_3 = -x_{36}a\,\hat{\mathbf{x}} + y_{36}b\,\hat{\mathbf{y}} + z_{36}c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Al XXVIII}$$

$$\mathbf{B}_{95} = x_{37}\,\mathbf{a}_1 + y_{37}\,\mathbf{a}_2 + z_{37}\,\mathbf{a}_3 = x_{37}a\,\hat{\mathbf{x}} + y_{37}b\,\hat{\mathbf{y}} + z_{37}c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Co IX}$$

$$\mathbf{B}_{96} = \left(\tfrac{1}{2} - x_{37}\right)\mathbf{a}_1 - y_{37}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{37}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{37}\right)a\,\hat{\mathbf{x}} - y_{37}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{37}\right)c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Co IX}$$

$$\mathbf{B}_{97} = \left(\tfrac{1}{2} + x_{37}\right)\mathbf{a}_1 - y_{37}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{37}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{37}\right)a\,\hat{\mathbf{x}} - y_{37}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{37}\right)c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Co IX}$$

$$\mathbf{B}_{98} = -x_{37}\,\mathbf{a}_1 + y_{37}\,\mathbf{a}_2 + z_{37}\,\mathbf{a}_3 = -x_{37}a\,\hat{\mathbf{x}} + y_{37}b\,\hat{\mathbf{y}} + z_{37}c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Co IX}$$

$$\mathbf{B}_{99} = x_{38}\,\mathbf{a}_1 + y_{38}\,\mathbf{a}_2 + z_{38}\,\mathbf{a}_3 = x_{38}a\,\hat{\mathbf{x}} + y_{38}b\,\hat{\mathbf{y}} + z_{38}c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Co X}$$

$$\mathbf{B}_{100} = \left(\tfrac{1}{2} - x_{38}\right)\mathbf{a}_1 - y_{38}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{38}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{38}\right)a\,\hat{\mathbf{x}} - y_{38}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{38}\right)c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Co X}$$

$$\mathbf{B}_{101} = \left(\tfrac{1}{2} + x_{38}\right)\mathbf{a}_1 - y_{38}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{38}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{38}\right)a\,\hat{\mathbf{x}} - y_{38}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{38}\right)c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Co X}$$

$$\mathbf{B}_{102} = -x_{38}\,\mathbf{a}_1 + y_{38}\,\mathbf{a}_2 + z_{38}\,\mathbf{a}_3 = -x_{38}a\,\hat{\mathbf{x}} + y_{38}b\,\hat{\mathbf{y}} + z_{38}c\,\hat{\mathbf{z}} \quad (4b) \quad \text{Co X}$$

References:

- J. Grin, U. Burkhard, M. Ellner, and K. Peters, *Crystal structure of orthorhombic Co_4Al_{13}*, J. Alloys Compd. **206**, 243–247 (1994), doi:10.1016/0925-8388(94)90043-4.

Found in:

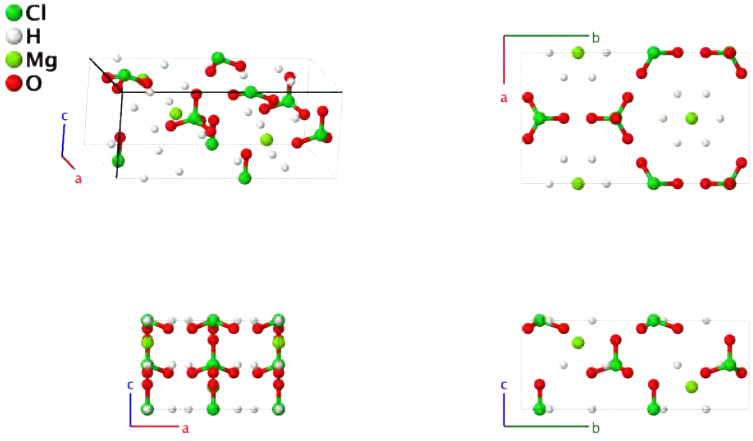
- R. Addou, E. Gaudry, T. Deniozou, M. Heggen, M. Feuerbacher, P. Gille, Y. Grin, R. Widmer, O. Gröning, V. Fournée, J.-M. Dubois, and J. Ledieu, *Structure investigation of the (100) surface of the orthorhombic $Al_{13}Co_4$ crystal*, Phys. Rev. B **80**, 014203 (2009), doi:10.1103/PhysRevB.80.014203.

Geometry files:

- CIF: pp. 1584
- POSCAR: pp. 1585

$Mg(ClO_4)_2{\cdot}6H_2O$ ($H4_{11}$) Structure: A2B6CD8_oP34_31_2a_2a2b_a_4a2b

http://aflow.org/prototype-encyclopedia/A2B6CD8_oP34_31_2a_2a2b_a_4a2b



Prototype	:	$Cl_2(H_2O)_6MgO_8$
AFLOW prototype label	:	A2B6CD8_oP34_31_2a_2a2b_a_4a2b
***Strukturbericht* designation**	:	$H4_{11}$
Pearson symbol	:	oP34
Space group number	:	31
Space group symbol	:	$Pmn2_1$
AFLOW prototype command	:	`aflow --proto=A2B6CD8_oP34_31_2a_2a2b_a_4a2b --params=`$a, b/a, c/a, y_1, z_1, y_2, z_2, y_3, z_3, y_4, z_4, y_5, z_5, y_6, z_6, y_7, z_7, y_8, z_8, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}$

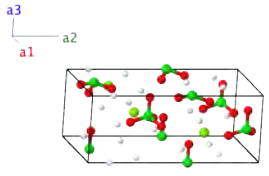
Other compounds with this structure

- $Co(ClO_4)_2{\cdot}6H_2O$, $Fe(ClO_4)_2{\cdot}6H_2O$, $Mn(ClO_4)_2{\cdot}6H_2O$, $Ni(ClO_4)_2{\cdot}6H_2O$, $Zn(ClO_4)_2{\cdot}6H_2O$, $Mg(BF_4)_2{\cdot}6H_2O$, $Co(BF_4)_2{\cdot}6H_2O$, $Fe(BF_4)_2{\cdot}6H_2O$, $Mn(BF_4)_2{\cdot}6H_2O$, $Ni(BF_4)_2{\cdot}6H_2O$, and $Zn(BF_4)_2{\cdot}6H_2O$

- (Gottfried, 1937) writes $z_{13} = 0.408$ for the coordinate of the (H_2O-II) molecule, but West uses $z_{13} = 0$, which gives a symmetric arrangement of water molecules around the chlorine atom. We use West's value here.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$\mathbf{B}_1$	=	$y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$y_1 b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(2*a*)	Cl I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	(2*a*)	Cl I
$\mathbf{B}_3$	=	$y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(2*a*)	Cl II
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	(2*a*)	Cl II
$\mathbf{B}_5$	=	$y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(2*a*)	H_2O I
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	(2*a*)	H_2O I
$\mathbf{B}_7$	=	$y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(2*a*)	H_2O II
$\mathbf{B}_8$	=	$\frac{1}{2}\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	(2*a*)	H_2O II
$\mathbf{B}_9$	=	$y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(2*a*)	Mg
$\mathbf{B}_{10}$	=	$\frac{1}{2}\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	(2*a*)	Mg
$\mathbf{B}_{11}$	=	$y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(2*a*)	O I
$\mathbf{B}_{12}$	=	$\frac{1}{2}\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	(2*a*)	O I
$\mathbf{B}_{13}$	=	$y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(2*a*)	O II
$\mathbf{B}_{14}$	=	$\frac{1}{2}\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right) c\,\hat{\mathbf{z}}$	(2*a*)	O II
$\mathbf{B}_{15}$	=	$y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(2*a*)	O III
$\mathbf{B}_{16}$	=	$\frac{1}{2}\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right) c\,\hat{\mathbf{z}}$	(2*a*)	O III
$\mathbf{B}_{17}$	=	$y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(2*a*)	O IV
$\mathbf{B}_{18}$	=	$\frac{1}{2}\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right) c\,\hat{\mathbf{z}}$	(2*a*)	O IV
$\mathbf{B}_{19}$	=	$x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$x_{10}a\,\hat{\mathbf{x}} + y_{10}b\,\hat{\mathbf{y}} + z_{10}c\,\hat{\mathbf{z}}$	(4*b*)	H_2O III
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} - x_{10}\right)\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{10}\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{10}\right)a\,\hat{\mathbf{x}} - y_{10}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{10}\right)c\,\hat{\mathbf{z}}$	(4*b*)	H_2O III
$\mathbf{B}_{21}$	=	$\left(\frac{1}{2} + x_{10}\right)\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{10}\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{10}\right)a\,\hat{\mathbf{x}} - y_{10}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{10}\right)c\,\hat{\mathbf{z}}$	(4*b*)	H_2O III
$\mathbf{B}_{22}$	=	$-x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$-x_{10}a\,\hat{\mathbf{x}} + y_{10}b\,\hat{\mathbf{y}} + z_{10}c\,\hat{\mathbf{z}}$	(4*b*)	H_2O III
$\mathbf{B}_{23}$	=	$x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$x_{11}a\,\hat{\mathbf{x}} + y_{11}b\,\hat{\mathbf{y}} + z_{11}c\,\hat{\mathbf{z}}$	(4*b*)	H_2O IV
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} - x_{11}\right)\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{11}\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{11}\right)a\,\hat{\mathbf{x}} - y_{11}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{11}\right)c\,\hat{\mathbf{z}}$	(4*b*)	H_2O IV
$\mathbf{B}_{25}$	=	$\left(\frac{1}{2} + x_{11}\right)\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{11}\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{11}\right)a\,\hat{\mathbf{x}} - y_{11}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{11}\right)c\,\hat{\mathbf{z}}$	(4*b*)	H_2O IV
$\mathbf{B}_{26}$	=	$-x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$-x_{11}a\,\hat{\mathbf{x}} + y_{11}b\,\hat{\mathbf{y}} + z_{11}c\,\hat{\mathbf{z}}$	(4*b*)	H_2O IV
$\mathbf{B}_{27}$	=	$x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3$	=	$x_{12}a\,\hat{\mathbf{x}} + y_{12}b\,\hat{\mathbf{y}} + z_{12}c\,\hat{\mathbf{z}}$	(4*b*)	O V
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} - x_{12}\right)\,\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{12}\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{12}\right)a\,\hat{\mathbf{x}} - y_{12}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{12}\right)c\,\hat{\mathbf{z}}$	(4*b*)	O V
$\mathbf{B}_{29}$	=	$\left(\frac{1}{2} + x_{12}\right)\,\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{12}\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{12}\right)a\,\hat{\mathbf{x}} - y_{12}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{12}\right)c\,\hat{\mathbf{z}}$	(4*b*)	O V
$\mathbf{B}_{30}$	=	$-x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3$	=	$-x_{12}a\,\hat{\mathbf{x}} + y_{12}b\,\hat{\mathbf{y}} + z_{12}c\,\hat{\mathbf{z}}$	(4*b*)	O V
$\mathbf{B}_{31}$	=	$x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3$	=	$x_{13}a\,\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} + z_{13}c\,\hat{\mathbf{z}}$	(4*b*)	O VI
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2} - x_{13}\right)\,\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{13}\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{13}\right)a\,\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{13}\right)c\,\hat{\mathbf{z}}$	(4*b*)	O VI
$\mathbf{B}_{33}$	=	$\left(\frac{1}{2} + x_{13}\right)\,\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{13}\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{13}\right)a\,\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{13}\right)c\,\hat{\mathbf{z}}$	(4*b*)	O VI
$\mathbf{B}_{34}$	=	$-x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3$	=	$-x_{13}a\,\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} + z_{13}c\,\hat{\mathbf{z}}$	(4*b*)	O VI

References:

- C. D. West, *Crystal Structures of Hydrated Compounds II. Structure Type $Mg(ClO_4)_2{\cdot}6H_2O$*, Zeitschrift für

Kristallographie - Crystalline Materials **91**, 480–493 (1935), doi:10.1524/zkri.1935.91.1.480.
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Found in:

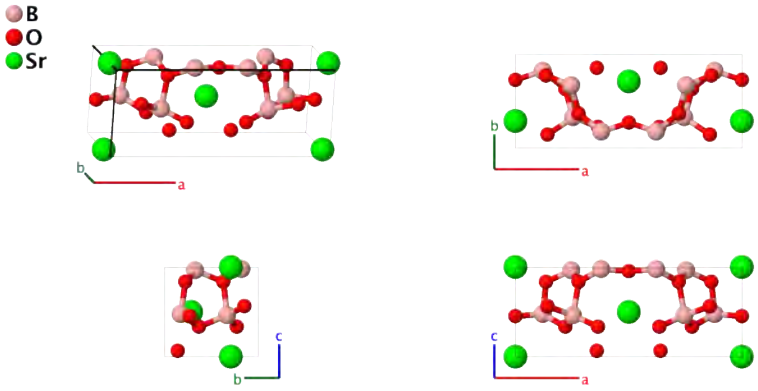
- K. Robertson and D. Bish, *Stability of phases in the $Mg(ClO_4)_2{\cdot}nH_2O$ system and implications for perchlorate occurrences on Mars*, J. Geophys. Res. **116**, E07006 (2011), doi:10.1029/2010JE003754.

Geometry files:

- CIF: pp. 1586
- POSCAR: pp. 1586

B_4SrO_7 Structure: A4B7C_oP24_31_2b_a3b_a

http://aflow.org/prototype-encyclopedia/A4B7C_oP24_31_2b_a3b_a

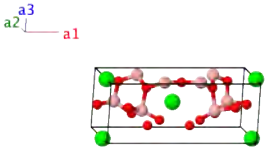


Prototype	:	B_4O_7Sr
AFLOW prototype label	:	A4B7C_oP24_31_2b_a3b_a
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP24
Space group number	:	31
Space group symbol	:	$Pmn2_1$
AFLOW prototype command	:	`aflow --proto=A4B7C_oP24_31_2b_a3b_a --params=`$a, b/a, c/a, y_1, z_1, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7$

- Space group $Pmn2_1$ #31 allows an arbitrary choice of the zero of the z-axis. Here it is chosen so that $z_1 = 0$ for the strontium atom.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$y_1 b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	$(2a)$	O I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\,\hat{\mathbf{z}}$	$(2a)$	O I
$\mathbf{B}_3$	=	$y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(2a)$	Sr
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	$(2a)$	Sr
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(4b)$	B I
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	$(4b)$	B I

$$
\begin{aligned}
\mathbf{B}_{7} &= \left(\tfrac{1}{2}+x_3\right)\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_3\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_3\right)a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_3\right)c\,\hat{\mathbf{z}} && (4b) && \text{B I}\\
\mathbf{B}_{8} &= -x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3 &&= -x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}} && (4b) && \text{B I}\\
\mathbf{B}_{9} &= x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3 &&= x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}} && (4b) && \text{B II}\\
\mathbf{B}_{10} &= \left(\tfrac{1}{2}-x_4\right)\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_4\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_4\right)a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_4\right)c\,\hat{\mathbf{z}} && (4b) && \text{B II}\\
\mathbf{B}_{11} &= \left(\tfrac{1}{2}+x_4\right)\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_4\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_4\right)a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_4\right)c\,\hat{\mathbf{z}} && (4b) && \text{B II}\\
\mathbf{B}_{12} &= -x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3 &&= -x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}} && (4b) && \text{B II}\\
\mathbf{B}_{13} &= x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3 &&= x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}} && (4b) && \text{O II}\\
\mathbf{B}_{14} &= \left(\tfrac{1}{2}-x_5\right)\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_5\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_5\right)a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_5\right)c\,\hat{\mathbf{z}} && (4b) && \text{O II}\\
\mathbf{B}_{15} &= \left(\tfrac{1}{2}+x_5\right)\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_5\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_5\right)a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_5\right)c\,\hat{\mathbf{z}} && (4b) && \text{O II}\\
\mathbf{B}_{16} &= -x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3 &&= -x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}} && (4b) && \text{O II}\\
\mathbf{B}_{17} &= x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 &&= x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}} && (4b) && \text{O III}\\
\mathbf{B}_{18} &= \left(\tfrac{1}{2}-x_6\right)\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_6\right)c\,\hat{\mathbf{z}} && (4b) && \text{O III}\\
\mathbf{B}_{19} &= \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_6\right)c\,\hat{\mathbf{z}} && (4b) && \text{O III}\\
\mathbf{B}_{20} &= -x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 &&= -x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}} && (4b) && \text{O III}\\
\mathbf{B}_{21} &= x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 &&= x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}} && (4b) && \text{O IV}\\
\mathbf{B}_{22} &= \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_7\right)c\,\hat{\mathbf{z}} && (4b) && \text{O IV}\\
\mathbf{B}_{23} &= \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_7\right)c\,\hat{\mathbf{z}} && (4b) && \text{O IV}\\
\mathbf{B}_{24} &= -x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 &&= -x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}} && (4b) && \text{O IV}
\end{aligned}
$$

References:

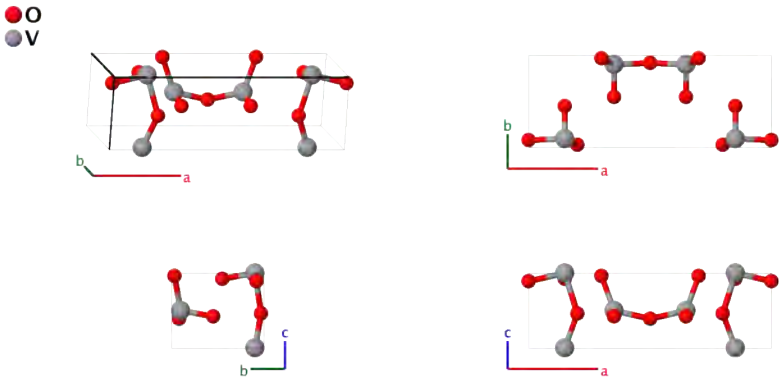
- J. Krogh-Moe, *The Crystal Structure of Strontium Diborate, $SrO{\cdot}2B_2O_3$*, Acta Chem. Scand. **18**, 2055–2060 (1964), doi:10.3891/acta.chem.scand.18-2055.

Geometry files:

- CIF: pp. 1586
- POSCAR: pp. 1587

$D8_7$ (Shcherbinaite, V_2O_5) (*obsolete*) Structure: A5B2_oP14_31_a2b_b

http://aflow.org/prototype-encyclopedia/A5B2_oP14_31_a2b_b

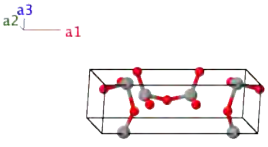


Prototype	:	O_5V_2
AFLOW prototype label	:	A5B2_oP14_31_a2b_b
***Strukturbericht* designation**	:	$D8_7$
Pearson symbol	:	oP14
Space group number	:	31
Space group symbol	:	$Pmn2_1$
AFLOW prototype command	:	`aflow --proto=A5B2_oP14_31_a2b_b --params=`$a, b/a, c/a, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4$

- This structure was reinvestigated by (Enjalbert, 1986) and others because of "[dissatisfaction with] some aspects of the crystal structure." They proposed a revised structure in space group *Pmmn*. We list this structure here for the historical record.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$y_1 b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	($2a$)	O I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\,\hat{\mathbf{z}}$	($2a$)	O I
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	($4b$)	O II
$\mathbf{B}_4$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	($4b$)	O II
$\mathbf{B}_5$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	($4b$)	O II

$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*b*)	O II
$\mathbf{B}_7$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*b*)	O III
$\mathbf{B}_8$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(4*b*)	O III
$\mathbf{B}_9$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(4*b*)	O III
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*b*)	O III
$\mathbf{B}_{11}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(4*b*)	V
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(4*b*)	V
$\mathbf{B}_{13}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(4*b*)	V
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(4*b*)	V

References:

- J. A. A. Ketelaar, *Die Kristallstruktur des Vanadinpentoxyds*, Zeitschrift für Kristallographie - Crystalline Materials **95**, 9–27 (1936), doi:10.1524/zkri.1936.95.1.9.
- J. A. A. Ketelaar, *Crystal Structure and Shape of Colloidal Particles of Vanadium Pentoxide*, Nature **137**, 316 (1936), doi:10.1038/137316a0.

Found in:

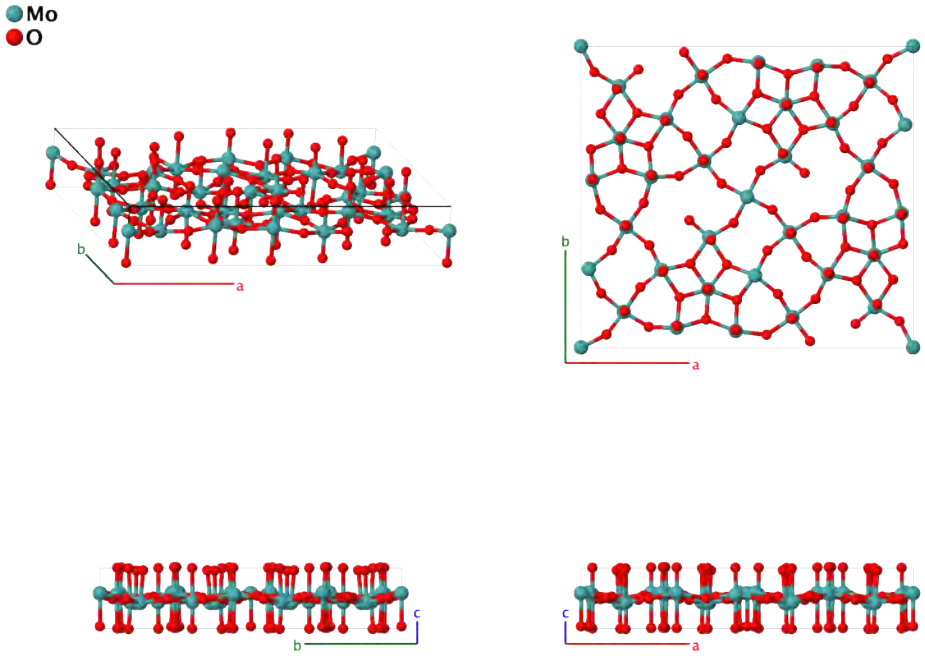
- R. Enjalbert and J. Galy, *A Refinement of the Structure of V_2O_5*, Acta Crystallogr. C **42**, 1467–1469 (1986), doi:10.1107/S0108270186091825.

Geometry files:

- CIF: pp. 1587
- POSCAR: pp. 1587

$Mo_{17}O_{47}$ Structure: A17B47_oP128_32_a8c_a23c

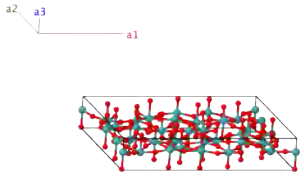
http://aflow.org/prototype-encyclopedia/A17B47_oP128_32_a8c_a23c



Prototype	:	$Mo_{17}O_{47}$
AFLOW prototype label	:	A17B47_oP128_32_a8c_a23c
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP128
Space group number	:	32
Space group symbol	:	*Pba*2
AFLOW prototype command	:	`aflow --proto=A17B47_oP128_32_a8c_a23c --params=`$a, b/a, c/a, z_1, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}, x_{17}, y_{17}, z_{17}, x_{18}, y_{18}, z_{18}, x_{19}, y_{19}, z_{19}, x_{20}, y_{20}, z_{20}, x_{21}, y_{21}, z_{21}, x_{22}, y_{22}, z_{22}, x_{23}, y_{23}, z_{23}, x_{24}, y_{24}, z_{24}, x_{25}, y_{25}, z_{25}, x_{26}, y_{26}, z_{26}, x_{27}, y_{27}, z_{27}, x_{28}, y_{28}, z_{28}, x_{29}, y_{29}, z_{29}, x_{30}, y_{30}, z_{30}, x_{31}, y_{31}, z_{31}, x_{32}, y_{32}, z_{32}, x_{33}, y_{33}, z_{33}$

Simple Orthorhombic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= a\,\hat{\mathbf{x}}\\ \mathbf{a}_2 &= b\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$z_1\,\mathbf{a}_3$	$z_1 c\,\hat{\mathbf{z}}$	$(2a)$	Mo I
$\mathbf{B}_2$	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	$(2a)$	Mo I
$\mathbf{B}_3$	$z_2\,\mathbf{a}_3$	$z_2 c\,\hat{\mathbf{z}}$	$(2a)$	O I
$\mathbf{B}_4$	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(2a)$	O I
$\mathbf{B}_5$	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(4c)$	Mo II
$\mathbf{B}_6$	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$-x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(4c)$	Mo II
$\mathbf{B}_7$	$\left(\frac{1}{2}+x_3\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_3\right)\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$\left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_3\right)b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(4c)$	Mo II
$\mathbf{B}_8$	$\left(\frac{1}{2}-x_3\right)\mathbf{a}_1 + \left(\frac{1}{2}+y_3\right)\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$\left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_3\right)b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(4c)$	Mo II
$\mathbf{B}_9$	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(4c)$	Mo III
$\mathbf{B}_{10}$	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(4c)$	Mo III
$\mathbf{B}_{11}$	$\left(\frac{1}{2}+x_4\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_4\right)b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(4c)$	Mo III
$\mathbf{B}_{12}$	$\left(\frac{1}{2}-x_4\right)\mathbf{a}_1 + \left(\frac{1}{2}+y_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$\left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_4\right)b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(4c)$	Mo III
$\mathbf{B}_{13}$	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(4c)$	Mo IV
$\mathbf{B}_{14}$	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(4c)$	Mo IV
$\mathbf{B}_{15}$	$\left(\frac{1}{2}+x_5\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_5\right)\mathbf{a}_2 + z_5\,\mathbf{a}_3$	$\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_5\right)b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(4c)$	Mo IV
$\mathbf{B}_{16}$	$\left(\frac{1}{2}-x_5\right)\mathbf{a}_1 + \left(\frac{1}{2}+y_5\right)\mathbf{a}_2 + z_5\,\mathbf{a}_3$	$\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_5\right)b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(4c)$	Mo IV
$\mathbf{B}_{17}$	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(4c)$	Mo V
$\mathbf{B}_{18}$	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	$-x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(4c)$	Mo V
$\mathbf{B}_{19}$	$\left(\frac{1}{2}+x_6\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_6\right)\mathbf{a}_2 + z_6\,\mathbf{a}_3$	$\left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_6\right)b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(4c)$	Mo V
$\mathbf{B}_{20}$	$\left(\frac{1}{2}-x_6\right)\mathbf{a}_1 + \left(\frac{1}{2}+y_6\right)\mathbf{a}_2 + z_6\,\mathbf{a}_3$	$\left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_6\right)b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(4c)$	Mo V
$\mathbf{B}_{21}$	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	$x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	$(4c)$	Mo VI
$\mathbf{B}_{22}$	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	$-x_7 a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	$(4c)$	Mo VI
$\mathbf{B}_{23}$	$\left(\frac{1}{2}+x_7\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_7\right)\mathbf{a}_2 + z_7\,\mathbf{a}_3$	$\left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_7\right)b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	$(4c)$	Mo VI
$\mathbf{B}_{24}$	$\left(\frac{1}{2}-x_7\right)\mathbf{a}_1 + \left(\frac{1}{2}+y_7\right)\mathbf{a}_2 + z_7\,\mathbf{a}_3$	$\left(\frac{1}{2}-x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_7\right)b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	$(4c)$	Mo VI
$\mathbf{B}_{25}$	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	$x_8 a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	$(4c)$	Mo VII
$\mathbf{B}_{26}$	$-x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	$-x_8 a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	$(4c)$	Mo VII
$\mathbf{B}_{27}$	$\left(\frac{1}{2}+x_8\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_8\right)\mathbf{a}_2 + z_8\,\mathbf{a}_3$	$\left(\frac{1}{2}+x_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_8\right)b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	$(4c)$	Mo VII

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\begin{aligned}
\mathbf{B}_{28} &= \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_8\right)\mathbf{a}_2+z_8\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_8\right)b\,\hat{\mathbf{y}}+z_8c\,\hat{\mathbf{z}} && (4c) && \text{Mo VII}\\
\mathbf{B}_{29} &= x_9\,\mathbf{a}_1+y_9\,\mathbf{a}_2+z_9\,\mathbf{a}_3 &&= x_9a\,\hat{\mathbf{x}}+y_9b\,\hat{\mathbf{y}}+z_9c\,\hat{\mathbf{z}} && (4c) && \text{Mo VIII}\\
\mathbf{B}_{30} &= -x_9\,\mathbf{a}_1-y_9\,\mathbf{a}_2+z_9\,\mathbf{a}_3 &&= -x_9a\,\hat{\mathbf{x}}-y_9b\,\hat{\mathbf{y}}+z_9c\,\hat{\mathbf{z}} && (4c) && \text{Mo VIII}\\
\mathbf{B}_{31} &= \left(\tfrac{1}{2}+x_9\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_9\right)\mathbf{a}_2+z_9\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_9\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_9\right)b\,\hat{\mathbf{y}}+z_9c\,\hat{\mathbf{z}} && (4c) && \text{Mo VIII}\\
\mathbf{B}_{32} &= \left(\tfrac{1}{2}-x_9\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_9\right)\mathbf{a}_2+z_9\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_9\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_9\right)b\,\hat{\mathbf{y}}+z_9c\,\hat{\mathbf{z}} && (4c) && \text{Mo VIII}\\
\mathbf{B}_{33} &= x_{10}\,\mathbf{a}_1+y_{10}\,\mathbf{a}_2+z_{10}\,\mathbf{a}_3 &&= x_{10}a\,\hat{\mathbf{x}}+y_{10}b\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}} && (4c) && \text{Mo IX}\\
\mathbf{B}_{34} &= -x_{10}\,\mathbf{a}_1-y_{10}\,\mathbf{a}_2+z_{10}\,\mathbf{a}_3 &&= -x_{10}a\,\hat{\mathbf{x}}-y_{10}b\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}} && (4c) && \text{Mo IX}\\
\mathbf{B}_{35} &= \left(\tfrac{1}{2}+x_{10}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{10}\right)\mathbf{a}_2+z_{10}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_{10}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{10}\right)b\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}} && (4c) && \text{Mo IX}\\
\mathbf{B}_{36} &= \left(\tfrac{1}{2}-x_{10}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{10}\right)\mathbf{a}_2+z_{10}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_{10}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{10}\right)b\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}} && (4c) && \text{Mo IX}\\
\mathbf{B}_{37} &= x_{11}\,\mathbf{a}_1+y_{11}\,\mathbf{a}_2+z_{11}\,\mathbf{a}_3 &&= x_{11}a\,\hat{\mathbf{x}}+y_{11}b\,\hat{\mathbf{y}}+z_{11}c\,\hat{\mathbf{z}} && (4c) && \text{O II}\\
\mathbf{B}_{38} &= -x_{11}\,\mathbf{a}_1-y_{11}\,\mathbf{a}_2+z_{11}\,\mathbf{a}_3 &&= -x_{11}a\,\hat{\mathbf{x}}-y_{11}b\,\hat{\mathbf{y}}+z_{11}c\,\hat{\mathbf{z}} && (4c) && \text{O II}\\
\mathbf{B}_{39} &= \left(\tfrac{1}{2}+x_{11}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{11}\right)\mathbf{a}_2+z_{11}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_{11}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{11}\right)b\,\hat{\mathbf{y}}+z_{11}c\,\hat{\mathbf{z}} && (4c) && \text{O II}\\
\mathbf{B}_{40} &= \left(\tfrac{1}{2}-x_{11}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{11}\right)\mathbf{a}_2+z_{11}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_{11}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{11}\right)b\,\hat{\mathbf{y}}+z_{11}c\,\hat{\mathbf{z}} && (4c) && \text{O II}\\
\mathbf{B}_{41} &= x_{12}\,\mathbf{a}_1+y_{12}\,\mathbf{a}_2+z_{12}\,\mathbf{a}_3 &&= x_{12}a\,\hat{\mathbf{x}}+y_{12}b\,\hat{\mathbf{y}}+z_{12}c\,\hat{\mathbf{z}} && (4c) && \text{O III}\\
\mathbf{B}_{42} &= -x_{12}\,\mathbf{a}_1-y_{12}\,\mathbf{a}_2+z_{12}\,\mathbf{a}_3 &&= -x_{12}a\,\hat{\mathbf{x}}-y_{12}b\,\hat{\mathbf{y}}+z_{12}c\,\hat{\mathbf{z}} && (4c) && \text{O III}\\
\mathbf{B}_{43} &= \left(\tfrac{1}{2}+x_{12}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{12}\right)\mathbf{a}_2+z_{12}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_{12}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{12}\right)b\,\hat{\mathbf{y}}+z_{12}c\,\hat{\mathbf{z}} && (4c) && \text{O III}\\
\mathbf{B}_{44} &= \left(\tfrac{1}{2}-x_{12}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{12}\right)\mathbf{a}_2+z_{12}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_{12}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{12}\right)b\,\hat{\mathbf{y}}+z_{12}c\,\hat{\mathbf{z}} && (4c) && \text{O III}\\
\mathbf{B}_{45} &= x_{13}\,\mathbf{a}_1+y_{13}\,\mathbf{a}_2+z_{13}\,\mathbf{a}_3 &&= x_{13}a\,\hat{\mathbf{x}}+y_{13}b\,\hat{\mathbf{y}}+z_{13}c\,\hat{\mathbf{z}} && (4c) && \text{O IV}\\
\mathbf{B}_{46} &= -x_{13}\,\mathbf{a}_1-y_{13}\,\mathbf{a}_2+z_{13}\,\mathbf{a}_3 &&= -x_{13}a\,\hat{\mathbf{x}}-y_{13}b\,\hat{\mathbf{y}}+z_{13}c\,\hat{\mathbf{z}} && (4c) && \text{O IV}\\
\mathbf{B}_{47} &= \left(\tfrac{1}{2}+x_{13}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{13}\right)\mathbf{a}_2+z_{13}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_{13}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{13}\right)b\,\hat{\mathbf{y}}+z_{13}c\,\hat{\mathbf{z}} && (4c) && \text{O IV}\\
\mathbf{B}_{48} &= \left(\tfrac{1}{2}-x_{13}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{13}\right)\mathbf{a}_2+z_{13}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_{13}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{13}\right)b\,\hat{\mathbf{y}}+z_{13}c\,\hat{\mathbf{z}} && (4c) && \text{O IV}\\
\mathbf{B}_{49} &= x_{14}\,\mathbf{a}_1+y_{14}\,\mathbf{a}_2+z_{14}\,\mathbf{a}_3 &&= x_{14}a\,\hat{\mathbf{x}}+y_{14}b\,\hat{\mathbf{y}}+z_{14}c\,\hat{\mathbf{z}} && (4c) && \text{O V}\\
\mathbf{B}_{50} &= -x_{14}\,\mathbf{a}_1-y_{14}\,\mathbf{a}_2+z_{14}\,\mathbf{a}_3 &&= -x_{14}a\,\hat{\mathbf{x}}-y_{14}b\,\hat{\mathbf{y}}+z_{14}c\,\hat{\mathbf{z}} && (4c) && \text{O V}\\
\mathbf{B}_{51} &= \left(\tfrac{1}{2}+x_{14}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{14}\right)\mathbf{a}_2+z_{14}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_{14}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{14}\right)b\,\hat{\mathbf{y}}+z_{14}c\,\hat{\mathbf{z}} && (4c) && \text{O V}\\
\mathbf{B}_{52} &= \left(\tfrac{1}{2}-x_{14}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{14}\right)\mathbf{a}_2+z_{14}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_{14}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{14}\right)b\,\hat{\mathbf{y}}+z_{14}c\,\hat{\mathbf{z}} && (4c) && \text{O V}\\
\mathbf{B}_{53} &= x_{15}\,\mathbf{a}_1+y_{15}\,\mathbf{a}_2+z_{15}\,\mathbf{a}_3 &&= x_{15}a\,\hat{\mathbf{x}}+y_{15}b\,\hat{\mathbf{y}}+z_{15}c\,\hat{\mathbf{z}} && (4c) && \text{O VI}\\
\mathbf{B}_{54} &= -x_{15}\,\mathbf{a}_1-y_{15}\,\mathbf{a}_2+z_{15}\,\mathbf{a}_3 &&= -x_{15}a\,\hat{\mathbf{x}}-y_{15}b\,\hat{\mathbf{y}}+z_{15}c\,\hat{\mathbf{z}} && (4c) && \text{O VI}\\
\mathbf{B}_{55} &= \left(\tfrac{1}{2}+x_{15}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{15}\right)\mathbf{a}_2+z_{15}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_{15}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{15}\right)b\,\hat{\mathbf{y}}+z_{15}c\,\hat{\mathbf{z}} && (4c) && \text{O VI}\\
\mathbf{B}_{56} &= \left(\tfrac{1}{2}-x_{15}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{15}\right)\mathbf{a}_2+z_{15}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_{15}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{15}\right)b\,\hat{\mathbf{y}}+z_{15}c\,\hat{\mathbf{z}} && (4c) && \text{O VI}\\
\mathbf{B}_{57} &= x_{16}\,\mathbf{a}_1+y_{16}\,\mathbf{a}_2+z_{16}\,\mathbf{a}_3 &&= x_{16}a\,\hat{\mathbf{x}}+y_{16}b\,\hat{\mathbf{y}}+z_{16}c\,\hat{\mathbf{z}} && (4c) && \text{O VII}\\
\mathbf{B}_{58} &= -x_{16}\,\mathbf{a}_1-y_{16}\,\mathbf{a}_2+z_{16}\,\mathbf{a}_3 &&= -x_{16}a\,\hat{\mathbf{x}}-y_{16}b\,\hat{\mathbf{y}}+z_{16}c\,\hat{\mathbf{z}} && (4c) && \text{O VII}\\
\mathbf{B}_{59} &= \left(\tfrac{1}{2}+x_{16}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{16}\right)\mathbf{a}_2+z_{16}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_{16}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{16}\right)b\,\hat{\mathbf{y}}+z_{16}c\,\hat{\mathbf{z}} && (4c) && \text{O VII}\\
\mathbf{B}_{60} &= \left(\tfrac{1}{2}-x_{16}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{16}\right)\mathbf{a}_2+z_{16}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_{16}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{16}\right)b\,\hat{\mathbf{y}}+z_{16}c\,\hat{\mathbf{z}} && (4c) && \text{O VII}\\
\mathbf{B}_{61} &= x_{17}\,\mathbf{a}_1+y_{17}\,\mathbf{a}_2+z_{17}\,\mathbf{a}_3 &&= x_{17}a\,\hat{\mathbf{x}}+y_{17}b\,\hat{\mathbf{y}}+z_{17}c\,\hat{\mathbf{z}} && (4c) && \text{O VIII}\\
\mathbf{B}_{62} &= -x_{17}\,\mathbf{a}_1-y_{17}\,\mathbf{a}_2+z_{17}\,\mathbf{a}_3 &&= -x_{17}a\,\hat{\mathbf{x}}-y_{17}b\,\hat{\mathbf{y}}+z_{17}c\,\hat{\mathbf{z}} && (4c) && \text{O VIII}\\
\mathbf{B}_{63} &= \left(\tfrac{1}{2}+x_{17}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{17}\right)\mathbf{a}_2+z_{17}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_{17}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{17}\right)b\,\hat{\mathbf{y}}+z_{17}c\,\hat{\mathbf{z}} && (4c) && \text{O VIII}
\end{aligned}
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\begin{aligned}
\mathbf{B}_{64} &= \left(\tfrac{1}{2}-x_{17}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{17}\right)\mathbf{a}_2+z_{17}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_{17}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{17}\right)b\,\hat{\mathbf{y}}+z_{17}c\,\hat{\mathbf{z}} && (4c) && \text{O VIII}\\
\mathbf{B}_{65} &= x_{18}\,\mathbf{a}_1+y_{18}\,\mathbf{a}_2+z_{18}\,\mathbf{a}_3 &&= x_{18}a\,\hat{\mathbf{x}}+y_{18}b\,\hat{\mathbf{y}}+z_{18}c\,\hat{\mathbf{z}} && (4c) && \text{O IX}\\
\mathbf{B}_{66} &= -x_{18}\,\mathbf{a}_1-y_{18}\,\mathbf{a}_2+z_{18}\,\mathbf{a}_3 &&= -x_{18}a\,\hat{\mathbf{x}}-y_{18}b\,\hat{\mathbf{y}}+z_{18}c\,\hat{\mathbf{z}} && (4c) && \text{O IX}\\
\mathbf{B}_{67} &= \left(\tfrac{1}{2}+x_{18}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{18}\right)\mathbf{a}_2+z_{18}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_{18}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{18}\right)b\,\hat{\mathbf{y}}+z_{18}c\,\hat{\mathbf{z}} && (4c) && \text{O IX}\\
\mathbf{B}_{68} &= \left(\tfrac{1}{2}-x_{18}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{18}\right)\mathbf{a}_2+z_{18}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_{18}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{18}\right)b\,\hat{\mathbf{y}}+z_{18}c\,\hat{\mathbf{z}} && (4c) && \text{O IX}\\
\mathbf{B}_{69} &= x_{19}\,\mathbf{a}_1+y_{19}\,\mathbf{a}_2+z_{19}\,\mathbf{a}_3 &&= x_{19}a\,\hat{\mathbf{x}}+y_{19}b\,\hat{\mathbf{y}}+z_{19}c\,\hat{\mathbf{z}} && (4c) && \text{O X}\\
\mathbf{B}_{70} &= -x_{19}\,\mathbf{a}_1-y_{19}\,\mathbf{a}_2+z_{19}\,\mathbf{a}_3 &&= -x_{19}a\,\hat{\mathbf{x}}-y_{19}b\,\hat{\mathbf{y}}+z_{19}c\,\hat{\mathbf{z}} && (4c) && \text{O X}\\
\mathbf{B}_{71} &= \left(\tfrac{1}{2}+x_{19}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{19}\right)\mathbf{a}_2+z_{19}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_{19}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{19}\right)b\,\hat{\mathbf{y}}+z_{19}c\,\hat{\mathbf{z}} && (4c) && \text{O X}\\
\mathbf{B}_{72} &= \left(\tfrac{1}{2}-x_{19}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{19}\right)\mathbf{a}_2+z_{19}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_{19}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{19}\right)b\,\hat{\mathbf{y}}+z_{19}c\,\hat{\mathbf{z}} && (4c) && \text{O X}\\
\mathbf{B}_{73} &= x_{20}\,\mathbf{a}_1+y_{20}\,\mathbf{a}_2+z_{20}\,\mathbf{a}_3 &&= x_{20}a\,\hat{\mathbf{x}}+y_{20}b\,\hat{\mathbf{y}}+z_{20}c\,\hat{\mathbf{z}} && (4c) && \text{O XI}\\
\mathbf{B}_{74} &= -x_{20}\,\mathbf{a}_1-y_{20}\,\mathbf{a}_2+z_{20}\,\mathbf{a}_3 &&= -x_{20}a\,\hat{\mathbf{x}}-y_{20}b\,\hat{\mathbf{y}}+z_{20}c\,\hat{\mathbf{z}} && (4c) && \text{O XI}\\
\mathbf{B}_{75} &= \left(\tfrac{1}{2}+x_{20}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{20}\right)\mathbf{a}_2+z_{20}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_{20}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{20}\right)b\,\hat{\mathbf{y}}+z_{20}c\,\hat{\mathbf{z}} && (4c) && \text{O XI}\\
\mathbf{B}_{76} &= \left(\tfrac{1}{2}-x_{20}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{20}\right)\mathbf{a}_2+z_{20}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_{20}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{20}\right)b\,\hat{\mathbf{y}}+z_{20}c\,\hat{\mathbf{z}} && (4c) && \text{O XI}\\
\mathbf{B}_{77} &= x_{21}\,\mathbf{a}_1+y_{21}\,\mathbf{a}_2+z_{21}\,\mathbf{a}_3 &&= x_{21}a\,\hat{\mathbf{x}}+y_{21}b\,\hat{\mathbf{y}}+z_{21}c\,\hat{\mathbf{z}} && (4c) && \text{O XII}\\
\mathbf{B}_{78} &= -x_{21}\,\mathbf{a}_1-y_{21}\,\mathbf{a}_2+z_{21}\,\mathbf{a}_3 &&= -x_{21}a\,\hat{\mathbf{x}}-y_{21}b\,\hat{\mathbf{y}}+z_{21}c\,\hat{\mathbf{z}} && (4c) && \text{O XII}\\
\mathbf{B}_{79} &= \left(\tfrac{1}{2}+x_{21}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{21}\right)\mathbf{a}_2+z_{21}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_{21}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{21}\right)b\,\hat{\mathbf{y}}+z_{21}c\,\hat{\mathbf{z}} && (4c) && \text{O XII}\\
\mathbf{B}_{80} &= \left(\tfrac{1}{2}-x_{21}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{21}\right)\mathbf{a}_2+z_{21}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_{21}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{21}\right)b\,\hat{\mathbf{y}}+z_{21}c\,\hat{\mathbf{z}} && (4c) && \text{O XII}\\
\mathbf{B}_{81} &= x_{22}\,\mathbf{a}_1+y_{22}\,\mathbf{a}_2+z_{22}\,\mathbf{a}_3 &&= x_{22}a\,\hat{\mathbf{x}}+y_{22}b\,\hat{\mathbf{y}}+z_{22}c\,\hat{\mathbf{z}} && (4c) && \text{O XIII}\\
\mathbf{B}_{82} &= -x_{22}\,\mathbf{a}_1-y_{22}\,\mathbf{a}_2+z_{22}\,\mathbf{a}_3 &&= -x_{22}a\,\hat{\mathbf{x}}-y_{22}b\,\hat{\mathbf{y}}+z_{22}c\,\hat{\mathbf{z}} && (4c) && \text{O XIII}\\
\mathbf{B}_{83} &= \left(\tfrac{1}{2}+x_{22}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{22}\right)\mathbf{a}_2+z_{22}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_{22}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{22}\right)b\,\hat{\mathbf{y}}+z_{22}c\,\hat{\mathbf{z}} && (4c) && \text{O XIII}\\
\mathbf{B}_{84} &= \left(\tfrac{1}{2}-x_{22}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{22}\right)\mathbf{a}_2+z_{22}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_{22}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{22}\right)b\,\hat{\mathbf{y}}+z_{22}c\,\hat{\mathbf{z}} && (4c) && \text{O XIII}\\
\mathbf{B}_{85} &= x_{23}\,\mathbf{a}_1+y_{23}\,\mathbf{a}_2+z_{23}\,\mathbf{a}_3 &&= x_{23}a\,\hat{\mathbf{x}}+y_{23}b\,\hat{\mathbf{y}}+z_{23}c\,\hat{\mathbf{z}} && (4c) && \text{O XIV}\\
\mathbf{B}_{86} &= -x_{23}\,\mathbf{a}_1-y_{23}\,\mathbf{a}_2+z_{23}\,\mathbf{a}_3 &&= -x_{23}a\,\hat{\mathbf{x}}-y_{23}b\,\hat{\mathbf{y}}+z_{23}c\,\hat{\mathbf{z}} && (4c) && \text{O XIV}\\
\mathbf{B}_{87} &= \left(\tfrac{1}{2}+x_{23}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{23}\right)\mathbf{a}_2+z_{23}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_{23}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{23}\right)b\,\hat{\mathbf{y}}+z_{23}c\,\hat{\mathbf{z}} && (4c) && \text{O XIV}\\
\mathbf{B}_{88} &= \left(\tfrac{1}{2}-x_{23}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{23}\right)\mathbf{a}_2+z_{23}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_{23}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{23}\right)b\,\hat{\mathbf{y}}+z_{23}c\,\hat{\mathbf{z}} && (4c) && \text{O XIV}\\
\mathbf{B}_{89} &= x_{24}\,\mathbf{a}_1+y_{24}\,\mathbf{a}_2+z_{24}\,\mathbf{a}_3 &&= x_{24}a\,\hat{\mathbf{x}}+y_{24}b\,\hat{\mathbf{y}}+z_{24}c\,\hat{\mathbf{z}} && (4c) && \text{O XV}\\
\mathbf{B}_{90} &= -x_{24}\,\mathbf{a}_1-y_{24}\,\mathbf{a}_2+z_{24}\,\mathbf{a}_3 &&= -x_{24}a\,\hat{\mathbf{x}}-y_{24}b\,\hat{\mathbf{y}}+z_{24}c\,\hat{\mathbf{z}} && (4c) && \text{O XV}\\
\mathbf{B}_{91} &= \left(\tfrac{1}{2}+x_{24}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{24}\right)\mathbf{a}_2+z_{24}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_{24}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{24}\right)b\,\hat{\mathbf{y}}+z_{24}c\,\hat{\mathbf{z}} && (4c) && \text{O XV}\\
\mathbf{B}_{92} &= \left(\tfrac{1}{2}-x_{24}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{24}\right)\mathbf{a}_2+z_{24}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_{24}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{24}\right)b\,\hat{\mathbf{y}}+z_{24}c\,\hat{\mathbf{z}} && (4c) && \text{O XV}\\
\mathbf{B}_{93} &= x_{25}\,\mathbf{a}_1+y_{25}\,\mathbf{a}_2+z_{25}\,\mathbf{a}_3 &&= x_{25}a\,\hat{\mathbf{x}}+y_{25}b\,\hat{\mathbf{y}}+z_{25}c\,\hat{\mathbf{z}} && (4c) && \text{O XVI}\\
\mathbf{B}_{94} &= -x_{25}\,\mathbf{a}_1-y_{25}\,\mathbf{a}_2+z_{25}\,\mathbf{a}_3 &&= -x_{25}a\,\hat{\mathbf{x}}-y_{25}b\,\hat{\mathbf{y}}+z_{25}c\,\hat{\mathbf{z}} && (4c) && \text{O XVI}\\
\mathbf{B}_{95} &= \left(\tfrac{1}{2}+x_{25}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{25}\right)\mathbf{a}_2+z_{25}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_{25}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{25}\right)b\,\hat{\mathbf{y}}+z_{25}c\,\hat{\mathbf{z}} && (4c) && \text{O XVI}\\
\mathbf{B}_{96} &= \left(\tfrac{1}{2}-x_{25}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{25}\right)\mathbf{a}_2+z_{25}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_{25}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{25}\right)b\,\hat{\mathbf{y}}+z_{25}c\,\hat{\mathbf{z}} && (4c) && \text{O XVI}\\
\mathbf{B}_{97} &= x_{26}\,\mathbf{a}_1+y_{26}\,\mathbf{a}_2+z_{26}\,\mathbf{a}_3 &&= x_{26}a\,\hat{\mathbf{x}}+y_{26}b\,\hat{\mathbf{y}}+z_{26}c\,\hat{\mathbf{z}} && (4c) && \text{O XVII}\\
\mathbf{B}_{98} &= -x_{26}\,\mathbf{a}_1-y_{26}\,\mathbf{a}_2+z_{26}\,\mathbf{a}_3 &&= -x_{26}a\,\hat{\mathbf{x}}-y_{26}b\,\hat{\mathbf{y}}+z_{26}c\,\hat{\mathbf{z}} && (4c) && \text{O XVII}\\
\mathbf{B}_{99} &= \left(\tfrac{1}{2}+x_{26}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{26}\right)\mathbf{a}_2+z_{26}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_{26}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{26}\right)b\,\hat{\mathbf{y}}+z_{26}c\,\hat{\mathbf{z}} && (4c) && \text{O XVII}
\end{aligned}
$$

$\mathbf{B}_{100}$	=	$\left(\frac{1}{2}-x_{26}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{26}\right)\mathbf{a}_2+z_{26}\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{26}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{26}\right)b\,\hat{\mathbf{y}}+z_{26}c\,\hat{\mathbf{z}}$	(4*c*)	O XVII
$\mathbf{B}_{101}$	=	$x_{27}\mathbf{a}_1+y_{27}\mathbf{a}_2+z_{27}\mathbf{a}_3$	=	$x_{27}a\,\hat{\mathbf{x}}+y_{27}b\,\hat{\mathbf{y}}+z_{27}c\,\hat{\mathbf{z}}$	(4*c*)	O XVIII
$\mathbf{B}_{102}$	=	$-x_{27}\mathbf{a}_1-y_{27}\mathbf{a}_2+z_{27}\mathbf{a}_3$	=	$-x_{27}a\,\hat{\mathbf{x}}-y_{27}b\,\hat{\mathbf{y}}+z_{27}c\,\hat{\mathbf{z}}$	(4*c*)	O XVIII
$\mathbf{B}_{103}$	=	$\left(\frac{1}{2}+x_{27}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{27}\right)\mathbf{a}_2+z_{27}\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{27}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{27}\right)b\,\hat{\mathbf{y}}+z_{27}c\,\hat{\mathbf{z}}$	(4*c*)	O XVIII
$\mathbf{B}_{104}$	=	$\left(\frac{1}{2}-x_{27}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{27}\right)\mathbf{a}_2+z_{27}\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{27}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{27}\right)b\,\hat{\mathbf{y}}+z_{27}c\,\hat{\mathbf{z}}$	(4*c*)	O XVIII
$\mathbf{B}_{105}$	=	$x_{28}\mathbf{a}_1+y_{28}\mathbf{a}_2+z_{28}\mathbf{a}_3$	=	$x_{28}a\,\hat{\mathbf{x}}+y_{28}b\,\hat{\mathbf{y}}+z_{28}c\,\hat{\mathbf{z}}$	(4*c*)	O XIX
$\mathbf{B}_{106}$	=	$-x_{28}\mathbf{a}_1-y_{28}\mathbf{a}_2+z_{28}\mathbf{a}_3$	=	$-x_{28}a\,\hat{\mathbf{x}}-y_{28}b\,\hat{\mathbf{y}}+z_{28}c\,\hat{\mathbf{z}}$	(4*c*)	O XIX
$\mathbf{B}_{107}$	=	$\left(\frac{1}{2}+x_{28}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{28}\right)\mathbf{a}_2+z_{28}\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{28}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{28}\right)b\,\hat{\mathbf{y}}+z_{28}c\,\hat{\mathbf{z}}$	(4*c*)	O XIX
$\mathbf{B}_{108}$	=	$\left(\frac{1}{2}-x_{28}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{28}\right)\mathbf{a}_2+z_{28}\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{28}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{28}\right)b\,\hat{\mathbf{y}}+z_{28}c\,\hat{\mathbf{z}}$	(4*c*)	O XIX
$\mathbf{B}_{109}$	=	$x_{29}\mathbf{a}_1+y_{29}\mathbf{a}_2+z_{29}\mathbf{a}_3$	=	$x_{29}a\,\hat{\mathbf{x}}+y_{29}b\,\hat{\mathbf{y}}+z_{29}c\,\hat{\mathbf{z}}$	(4*c*)	O XX
$\mathbf{B}_{110}$	=	$-x_{29}\mathbf{a}_1-y_{29}\mathbf{a}_2+z_{29}\mathbf{a}_3$	=	$-x_{29}a\,\hat{\mathbf{x}}-y_{29}b\,\hat{\mathbf{y}}+z_{29}c\,\hat{\mathbf{z}}$	(4*c*)	O XX
$\mathbf{B}_{111}$	=	$\left(\frac{1}{2}+x_{29}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{29}\right)\mathbf{a}_2+z_{29}\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{29}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{29}\right)b\,\hat{\mathbf{y}}+z_{29}c\,\hat{\mathbf{z}}$	(4*c*)	O XX
$\mathbf{B}_{112}$	=	$\left(\frac{1}{2}-x_{29}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{29}\right)\mathbf{a}_2+z_{29}\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{29}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{29}\right)b\,\hat{\mathbf{y}}+z_{29}c\,\hat{\mathbf{z}}$	(4*c*)	O XX
$\mathbf{B}_{113}$	=	$x_{30}\mathbf{a}_1+y_{30}\mathbf{a}_2+z_{30}\mathbf{a}_3$	=	$x_{30}a\,\hat{\mathbf{x}}+y_{30}b\,\hat{\mathbf{y}}+z_{30}c\,\hat{\mathbf{z}}$	(4*c*)	O XXI
$\mathbf{B}_{114}$	=	$-x_{30}\mathbf{a}_1-y_{30}\mathbf{a}_2+z_{30}\mathbf{a}_3$	=	$-x_{30}a\,\hat{\mathbf{x}}-y_{30}b\,\hat{\mathbf{y}}+z_{30}c\,\hat{\mathbf{z}}$	(4*c*)	O XXI
$\mathbf{B}_{115}$	=	$\left(\frac{1}{2}+x_{30}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{30}\right)\mathbf{a}_2+z_{30}\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{30}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{30}\right)b\,\hat{\mathbf{y}}+z_{30}c\,\hat{\mathbf{z}}$	(4*c*)	O XXI
$\mathbf{B}_{116}$	=	$\left(\frac{1}{2}-x_{30}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{30}\right)\mathbf{a}_2+z_{30}\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{30}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{30}\right)b\,\hat{\mathbf{y}}+z_{30}c\,\hat{\mathbf{z}}$	(4*c*)	O XXI
$\mathbf{B}_{117}$	=	$x_{31}\mathbf{a}_1+y_{31}\mathbf{a}_2+z_{31}\mathbf{a}_3$	=	$x_{31}a\,\hat{\mathbf{x}}+y_{31}b\,\hat{\mathbf{y}}+z_{31}c\,\hat{\mathbf{z}}$	(4*c*)	O XXII
$\mathbf{B}_{118}$	=	$-x_{31}\mathbf{a}_1-y_{31}\mathbf{a}_2+z_{31}\mathbf{a}_3$	=	$-x_{31}a\,\hat{\mathbf{x}}-y_{31}b\,\hat{\mathbf{y}}+z_{31}c\,\hat{\mathbf{z}}$	(4*c*)	O XXII
$\mathbf{B}_{119}$	=	$\left(\frac{1}{2}+x_{31}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{31}\right)\mathbf{a}_2+z_{31}\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{31}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{31}\right)b\,\hat{\mathbf{y}}+z_{31}c\,\hat{\mathbf{z}}$	(4*c*)	O XXII
$\mathbf{B}_{120}$	=	$\left(\frac{1}{2}-x_{31}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{31}\right)\mathbf{a}_2+z_{31}\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{31}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{31}\right)b\,\hat{\mathbf{y}}+z_{31}c\,\hat{\mathbf{z}}$	(4*c*)	O XXII
$\mathbf{B}_{121}$	=	$x_{32}\mathbf{a}_1+y_{32}\mathbf{a}_2+z_{32}\mathbf{a}_3$	=	$x_{32}a\,\hat{\mathbf{x}}+y_{32}b\,\hat{\mathbf{y}}+z_{32}c\,\hat{\mathbf{z}}$	(4*c*)	O XXIII
$\mathbf{B}_{122}$	=	$-x_{32}\mathbf{a}_1-y_{32}\mathbf{a}_2+z_{32}\mathbf{a}_3$	=	$-x_{32}a\,\hat{\mathbf{x}}-y_{32}b\,\hat{\mathbf{y}}+z_{32}c\,\hat{\mathbf{z}}$	(4*c*)	O XXIII
$\mathbf{B}_{123}$	=	$\left(\frac{1}{2}+x_{32}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{32}\right)\mathbf{a}_2+z_{32}\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{32}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{32}\right)b\,\hat{\mathbf{y}}+z_{32}c\,\hat{\mathbf{z}}$	(4*c*)	O XXIII
$\mathbf{B}_{124}$	=	$\left(\frac{1}{2}-x_{32}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{32}\right)\mathbf{a}_2+z_{32}\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{32}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{32}\right)b\,\hat{\mathbf{y}}+z_{32}c\,\hat{\mathbf{z}}$	(4*c*)	O XXIII
$\mathbf{B}_{125}$	=	$x_{33}\mathbf{a}_1+y_{33}\mathbf{a}_2+z_{33}\mathbf{a}_3$	=	$x_{33}a\,\hat{\mathbf{x}}+y_{33}b\,\hat{\mathbf{y}}+z_{33}c\,\hat{\mathbf{z}}$	(4*c*)	O XXIV
$\mathbf{B}_{126}$	=	$-x_{33}\mathbf{a}_1-y_{33}\mathbf{a}_2+z_{33}\mathbf{a}_3$	=	$-x_{33}a\,\hat{\mathbf{x}}-y_{33}b\,\hat{\mathbf{y}}+z_{33}c\,\hat{\mathbf{z}}$	(4*c*)	O XXIV
$\mathbf{B}_{127}$	=	$\left(\frac{1}{2}+x_{33}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{33}\right)\mathbf{a}_2+z_{33}\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{33}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{33}\right)b\,\hat{\mathbf{y}}+z_{33}c\,\hat{\mathbf{z}}$	(4*c*)	O XXIV
$\mathbf{B}_{128}$	=	$\left(\frac{1}{2}-x_{33}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{33}\right)\mathbf{a}_2+z_{33}\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{33}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{33}\right)b\,\hat{\mathbf{y}}+z_{33}c\,\hat{\mathbf{z}}$	(4*c*)	O XXIV

References:

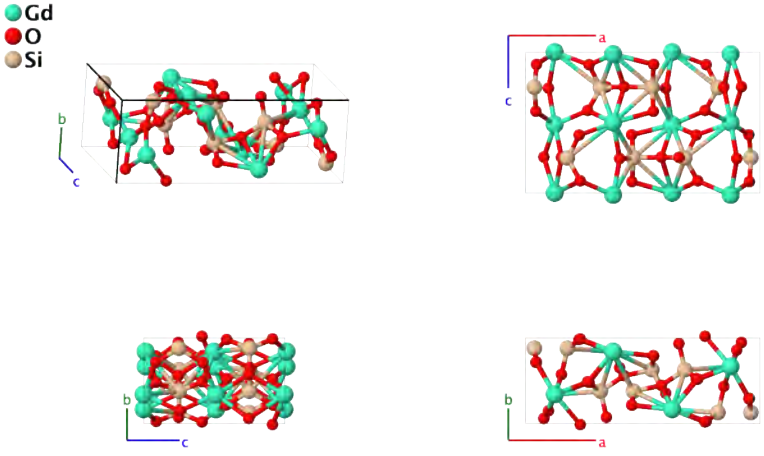
- L. Kihlborg, *Least Squares Refinement of the Structure of $Mo_{17}O_{47}$*, Acta Chem. Scand. **17**, 1485–1487 (1963), doi:10.3891/acta.chem.scand.17-1485.

Geometry files:

- CIF: pp. 1587
- POSCAR: pp. 1588

Possible δ-$Gd_2Si_2O_7$ Structure: A2B7C2_oP44_33_2a_7a_2a

http://aflow.org/prototype-encyclopedia/A2B7C2_oP44_33_2a_7a_2a



Prototype	:	$Gd_2O_7Si_2$
AFLOW prototype label	:	A2B7C2_oP44_33_2a_7a_2a
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP44
Space group number	:	33
Space group symbol	:	$Pna2_1$
AFLOW prototype command	:	`aflow --proto=A2B7C2_oP44_33_2a_7a_2a --params=`$a, b/a, c/a, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}$

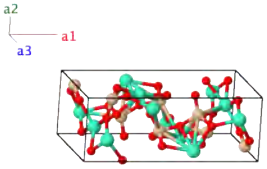
Other compounds with this structure

- δ-$Ho_2O_7Si_2$, δ-$Dy_2O_7Si_2$, and δ-$Y_2O_7Si_2$

- (Smolin, 1970) found that some structures of $RE_2O_7Si_2$ (RE = Ho, Dy, Gd, Y) were in the orthorhombic $Pna2_1$ #33 space group, in which case this would be the prototype of δ-$RE_2O_7Si_2$ (Becerro, 2004). However, (Dias, 1990) found δ-$Y_2O_7Si_2$ to be in the centro-symmetric *Pnma* #62 space group. This was supported by (Becerro, 2004), who found only one yttrium site in the δ-structure. In addition, if we allow a small amount of uncertainty (0.2 Å) in positions, AFLOW-SYM and FINDSYM place this structure in the *Pnma* group. Nevertheless we have found no work explicitly stating that the structure of (Smolin, 1970) is in error, and indeed (Christensen, 1994) found δ-$Y_2O_7Si_2$ in space group $Pna2_1$. Given this ambiguity, we list $Gd_2O_7Si_2$ as a possible prototype for the δ-phase pyrosilicates.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	($4a$)	Gd I
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	($4a$)	Gd I
$\mathbf{B}_3$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_1\right)\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_1\right) b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	($4a$)	Gd I
$\mathbf{B}_4$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_1\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_1\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	($4a$)	Gd I
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	($4a$)	Gd II
$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	($4a$)	Gd II
$\mathbf{B}_7$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right) b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	($4a$)	Gd II
$\mathbf{B}_8$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	($4a$)	Gd II
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	($4a$)	O I
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	($4a$)	O I
$\mathbf{B}_{11}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right) b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	($4a$)	O I
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	($4a$)	O I
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	($4a$)	O II
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	($4a$)	O II
$\mathbf{B}_{15}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right) b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	($4a$)	O II
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	($4a$)	O II
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	($4a$)	O III
$\mathbf{B}_{18}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	($4a$)	O III
$\mathbf{B}_{19}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right) b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	($4a$)	O III
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	($4a$)	O III
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	($4a$)	O IV
$\mathbf{B}_{22}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	($4a$)	O IV
$\mathbf{B}_{23}$	=	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right) b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	($4a$)	O IV
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	($4a$)	O IV
$\mathbf{B}_{25}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	($4a$)	O V
$\mathbf{B}_{26}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right) c\,\hat{\mathbf{z}}$	($4a$)	O V
$\mathbf{B}_{27}$	=	$\left(\frac{1}{2} + x_7\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_7\right)\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_7\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_7\right) b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	($4a$)	O V
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} - x_7\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_7\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right) c\,\hat{\mathbf{z}}$	($4a$)	O V

$\mathbf{B}_{29}$	=	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(4*a*)	O VI
$\mathbf{B}_{30}$	=	$-x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right) c\,\hat{\mathbf{z}}$	(4*a*)	O VI
$\mathbf{B}_{31}$	=	$\left(\frac{1}{2} + x_8\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_8\right)\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_8\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_8\right) b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(4*a*)	O VI
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2} - x_8\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_8\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_8\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_8\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right) c\,\hat{\mathbf{z}}$	(4*a*)	O VI
$\mathbf{B}_{33}$	=	$x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(4*a*)	O VII
$\mathbf{B}_{34}$	=	$-x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right) c\,\hat{\mathbf{z}}$	(4*a*)	O VII
$\mathbf{B}_{35}$	=	$\left(\frac{1}{2} + x_9\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_9\right)\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_9\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_9\right) b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(4*a*)	O VII
$\mathbf{B}_{36}$	=	$\left(\frac{1}{2} - x_9\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_9\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_9\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_9\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right) c\,\hat{\mathbf{z}}$	(4*a*)	O VII
$\mathbf{B}_{37}$	=	$x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$x_{10} a\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}}$	(4*a*)	Si I
$\mathbf{B}_{38}$	=	$-x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{10}\right)\mathbf{a}_3$	=	$-x_{10} a\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{10}\right) c\,\hat{\mathbf{z}}$	(4*a*)	Si I
$\mathbf{B}_{39}$	=	$\left(\frac{1}{2} + x_{10}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{10}\right)\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{10}\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{10}\right) b\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}}$	(4*a*)	Si I
$\mathbf{B}_{40}$	=	$\left(\frac{1}{2} - x_{10}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{10}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{10}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{10}\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{10}\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{10}\right) c\,\hat{\mathbf{z}}$	(4*a*)	Si I
$\mathbf{B}_{41}$	=	$x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$x_{11} a\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + z_{11} c\,\hat{\mathbf{z}}$	(4*a*)	Si II
$\mathbf{B}_{42}$	=	$-x_{11}\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{11}\right)\mathbf{a}_3$	=	$-x_{11} a\,\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{11}\right) c\,\hat{\mathbf{z}}$	(4*a*)	Si II
$\mathbf{B}_{43}$	=	$\left(\frac{1}{2} + x_{11}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{11}\right)\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{11}\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{11}\right) b\,\hat{\mathbf{y}} + z_{11} c\,\hat{\mathbf{z}}$	(4*a*)	Si II
$\mathbf{B}_{44}$	=	$\left(\frac{1}{2} - x_{11}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{11}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{11}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{11}\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{11}\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{11}\right) c\,\hat{\mathbf{z}}$	(4*a*)	Si II

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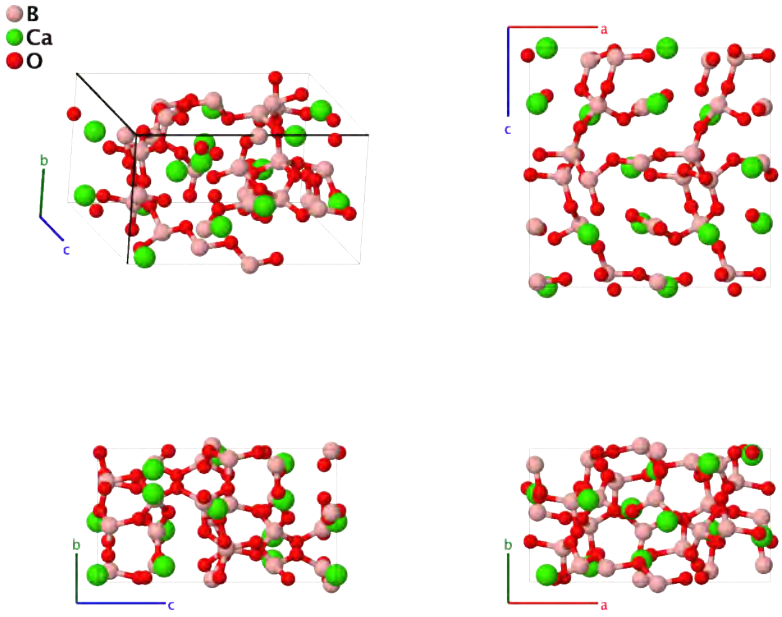
Geometry files:

- CIF: pp. 1589
- POSCAR: pp. 1589

CaB_2O_4 (III) Structure: A2BC4_oP84_33_6a_3a_12a

http://aflow.org/prototype-encyclopedia/A2BC4_oP84_33_6a_3a_12a

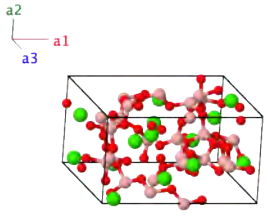


Prototype	:	B_2CaO_4
AFLOW prototype label	:	A2BC4_oP84_33_6a_3a_12a
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP84
Space group number	:	33
Space group symbol	:	$Pna2_1$
AFLOW prototype command	:	`aflow --proto=A2BC4_oP84_33_6a_3a_12a --params=`$a, b/a, c/a, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}, x_{17}, y_{17}, z_{17}, x_{18}, y_{18}, z_{18}, x_{19}, y_{19}, z_{19}, x_{20}, y_{20}, z_{20}, x_{21}, y_{21}, z_{21}$

- CaB_2O_4 exists in at least four phases (Marezio, 1969):
- I - The ground state, stable below 1.2 GPa, *Strukturbericht* $E3_2$.
- II – Orthorhombic high pressure structure, stable between 1.2 and 1.5 GPa, presumably calciborite.
- III – Orthorhombic high pressure structure, stable between 1.5 and 2.5 GPa (this structure).
- IV – Cubic high pressure structure, stable between 2.5 and 4.0 GPa.

Simple Orthorhombic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= a\,\hat{\mathbf{x}}\\ \mathbf{a}_2 &= b\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	$x_1 a\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	$(4a)$	B I
$\mathbf{B}_2$	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	$-x_1 a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	$(4a)$	B I
$\mathbf{B}_3$	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_1\right)\mathbf{a}_2 + z_1\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_1\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_1\right) b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	$(4a)$	B I
$\mathbf{B}_4$	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_1\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_1\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	$(4a)$	B I
$\mathbf{B}_5$	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	$x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(4a)$	B II
$\mathbf{B}_6$	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	$-x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	$(4a)$	B II
$\mathbf{B}_7$	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 + z_2\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right) b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(4a)$	B II
$\mathbf{B}_8$	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	$(4a)$	B II
$\mathbf{B}_9$	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(4a)$	B III
$\mathbf{B}_{10}$	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	$-x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	$(4a)$	B III
$\mathbf{B}_{11}$	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_3\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right) b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(4a)$	B III
$\mathbf{B}_{12}$	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_3\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	$(4a)$	B III
$\mathbf{B}_{13}$	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(4a)$	B IV
$\mathbf{B}_{14}$	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	$(4a)$	B IV
$\mathbf{B}_{15}$	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right) b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(4a)$	B IV
$\mathbf{B}_{16}$	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	$(4a)$	B IV
$\mathbf{B}_{17}$	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(4a)$	B V
$\mathbf{B}_{18}$	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	$(4a)$	B V
$\mathbf{B}_{19}$	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + z_5\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right) b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(4a)$	B V
$\mathbf{B}_{20}$	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_5\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	$(4a)$	B V
$\mathbf{B}_{21}$	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(4a)$	B VI
$\mathbf{B}_{22}$	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	$-x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	$(4a)$	B VI
$\mathbf{B}_{23}$	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + z_6\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right) b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(4a)$	B VI

$$
\begin{array}{lcccccc}
\mathbf{B}_{24} & = & \left(\frac{1}{2}-x_6\right)\mathbf{a}_1+\left(\frac{1}{2}+y_6\right)\mathbf{a}_2+\left(\frac{1}{2}+z_6\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_6\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}} & (4a) & \text{B VI}\\
\mathbf{B}_{25} & = & x_7\,\mathbf{a}_1+y_7\,\mathbf{a}_2+z_7\,\mathbf{a}_3 & = & x_7a\,\hat{\mathbf{x}}+y_7b\,\hat{\mathbf{y}}+z_7c\,\hat{\mathbf{z}} & (4a) & \text{Ca I}\\
\mathbf{B}_{26} & = & -x_7\,\mathbf{a}_1-y_7\,\mathbf{a}_2+\left(\frac{1}{2}+z_7\right)\mathbf{a}_3 & = & -x_7a\,\hat{\mathbf{x}}-y_7b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}} & (4a) & \text{Ca I}\\
\mathbf{B}_{27} & = & \left(\frac{1}{2}+x_7\right)\mathbf{a}_1+\left(\frac{1}{2}-y_7\right)\mathbf{a}_2+z_7\,\mathbf{a}_3 & = & \left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_7\right)b\,\hat{\mathbf{y}}+z_7c\,\hat{\mathbf{z}} & (4a) & \text{Ca I}\\
\mathbf{B}_{28} & = & \left(\frac{1}{2}-x_7\right)\mathbf{a}_1+\left(\frac{1}{2}+y_7\right)\mathbf{a}_2+\left(\frac{1}{2}+z_7\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_7\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_7\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}} & (4a) & \text{Ca I}\\
\mathbf{B}_{29} & = & x_8\,\mathbf{a}_1+y_8\,\mathbf{a}_2+z_8\,\mathbf{a}_3 & = & x_8a\,\hat{\mathbf{x}}+y_8b\,\hat{\mathbf{y}}+z_8c\,\hat{\mathbf{z}} & (4a) & \text{Ca II}\\
\mathbf{B}_{30} & = & -x_8\,\mathbf{a}_1-y_8\,\mathbf{a}_2+\left(\frac{1}{2}+z_8\right)\mathbf{a}_3 & = & -x_8a\,\hat{\mathbf{x}}-y_8b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_8\right)c\,\hat{\mathbf{z}} & (4a) & \text{Ca II}\\
\mathbf{B}_{31} & = & \left(\frac{1}{2}+x_8\right)\mathbf{a}_1+\left(\frac{1}{2}-y_8\right)\mathbf{a}_2+z_8\,\mathbf{a}_3 & = & \left(\frac{1}{2}+x_8\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_8\right)b\,\hat{\mathbf{y}}+z_8c\,\hat{\mathbf{z}} & (4a) & \text{Ca II}\\
\mathbf{B}_{32} & = & \left(\frac{1}{2}-x_8\right)\mathbf{a}_1+\left(\frac{1}{2}+y_8\right)\mathbf{a}_2+\left(\frac{1}{2}+z_8\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_8\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_8\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_8\right)c\,\hat{\mathbf{z}} & (4a) & \text{Ca II}\\
\mathbf{B}_{33} & = & x_9\,\mathbf{a}_1+y_9\,\mathbf{a}_2+z_9\,\mathbf{a}_3 & = & x_9a\,\hat{\mathbf{x}}+y_9b\,\hat{\mathbf{y}}+z_9c\,\hat{\mathbf{z}} & (4a) & \text{Ca III}\\
\mathbf{B}_{34} & = & -x_9\,\mathbf{a}_1-y_9\,\mathbf{a}_2+\left(\frac{1}{2}+z_9\right)\mathbf{a}_3 & = & -x_9a\,\hat{\mathbf{x}}-y_9b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_9\right)c\,\hat{\mathbf{z}} & (4a) & \text{Ca III}\\
\mathbf{B}_{35} & = & \left(\frac{1}{2}+x_9\right)\mathbf{a}_1+\left(\frac{1}{2}-y_9\right)\mathbf{a}_2+z_9\,\mathbf{a}_3 & = & \left(\frac{1}{2}+x_9\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_9\right)b\,\hat{\mathbf{y}}+z_9c\,\hat{\mathbf{z}} & (4a) & \text{Ca III}\\
\mathbf{B}_{36} & = & \left(\frac{1}{2}-x_9\right)\mathbf{a}_1+\left(\frac{1}{2}+y_9\right)\mathbf{a}_2+\left(\frac{1}{2}+z_9\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_9\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_9\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_9\right)c\,\hat{\mathbf{z}} & (4a) & \text{Ca III}\\
\mathbf{B}_{37} & = & x_{10}\,\mathbf{a}_1+y_{10}\,\mathbf{a}_2+z_{10}\,\mathbf{a}_3 & = & x_{10}a\,\hat{\mathbf{x}}+y_{10}b\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}} & (4a) & \text{O I}\\
\mathbf{B}_{38} & = & -x_{10}\,\mathbf{a}_1-y_{10}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{10}\right)\mathbf{a}_3 & = & -x_{10}a\,\hat{\mathbf{x}}-y_{10}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{10}\right)c\,\hat{\mathbf{z}} & (4a) & \text{O I}\\
\mathbf{B}_{39} & = & \left(\frac{1}{2}+x_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{10}\right)\mathbf{a}_2+z_{10}\,\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{10}\right)b\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}} & (4a) & \text{O I}\\
\mathbf{B}_{40} & = & \left(\frac{1}{2}-x_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{10}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{10}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{10}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{10}\right)c\,\hat{\mathbf{z}} & (4a) & \text{O I}\\
\mathbf{B}_{41} & = & x_{11}\,\mathbf{a}_1+y_{11}\,\mathbf{a}_2+z_{11}\,\mathbf{a}_3 & = & x_{11}a\,\hat{\mathbf{x}}+y_{11}b\,\hat{\mathbf{y}}+z_{11}c\,\hat{\mathbf{z}} & (4a) & \text{O II}\\
\mathbf{B}_{42} & = & -x_{11}\,\mathbf{a}_1-y_{11}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{11}\right)\mathbf{a}_3 & = & -x_{11}a\,\hat{\mathbf{x}}-y_{11}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{11}\right)c\,\hat{\mathbf{z}} & (4a) & \text{O II}\\
\mathbf{B}_{43} & = & \left(\frac{1}{2}+x_{11}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{11}\right)\mathbf{a}_2+z_{11}\,\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{11}\right)b\,\hat{\mathbf{y}}+z_{11}c\,\hat{\mathbf{z}} & (4a) & \text{O II}\\
\mathbf{B}_{44} & = & \left(\frac{1}{2}-x_{11}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{11}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{11}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{11}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{11}\right)c\,\hat{\mathbf{z}} & (4a) & \text{O II}\\
\mathbf{B}_{45} & = & x_{12}\,\mathbf{a}_1+y_{12}\,\mathbf{a}_2+z_{12}\,\mathbf{a}_3 & = & x_{12}a\,\hat{\mathbf{x}}+y_{12}b\,\hat{\mathbf{y}}+z_{12}c\,\hat{\mathbf{z}} & (4a) & \text{O III}\\
\mathbf{B}_{46} & = & -x_{12}\,\mathbf{a}_1-y_{12}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{12}\right)\mathbf{a}_3 & = & -x_{12}a\,\hat{\mathbf{x}}-y_{12}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{12}\right)c\,\hat{\mathbf{z}} & (4a) & \text{O III}\\
\mathbf{B}_{47} & = & \left(\frac{1}{2}+x_{12}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{12}\right)\mathbf{a}_2+z_{12}\,\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{12}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{12}\right)b\,\hat{\mathbf{y}}+z_{12}c\,\hat{\mathbf{z}} & (4a) & \text{O III}\\
\mathbf{B}_{48} & = & \left(\frac{1}{2}-x_{12}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{12}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{12}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{12}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{12}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{12}\right)c\,\hat{\mathbf{z}} & (4a) & \text{O III}\\
\mathbf{B}_{49} & = & x_{13}\,\mathbf{a}_1+y_{13}\,\mathbf{a}_2+z_{13}\,\mathbf{a}_3 & = & x_{13}a\,\hat{\mathbf{x}}+y_{13}b\,\hat{\mathbf{y}}+z_{13}c\,\hat{\mathbf{z}} & (4a) & \text{O IV}\\
\mathbf{B}_{50} & = & -x_{13}\,\mathbf{a}_1-y_{13}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{13}\right)\mathbf{a}_3 & = & -x_{13}a\,\hat{\mathbf{x}}-y_{13}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{13}\right)c\,\hat{\mathbf{z}} & (4a) & \text{O IV}\\
\mathbf{B}_{51} & = & \left(\frac{1}{2}+x_{13}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{13}\right)\mathbf{a}_2+z_{13}\,\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{13}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{13}\right)b\,\hat{\mathbf{y}}+z_{13}c\,\hat{\mathbf{z}} & (4a) & \text{O IV}\\
\mathbf{B}_{52} & = & \left(\frac{1}{2}-x_{13}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{13}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{13}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{13}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{13}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{13}\right)c\,\hat{\mathbf{z}} & (4a) & \text{O IV}\\
\mathbf{B}_{53} & = & x_{14}\,\mathbf{a}_1+y_{14}\,\mathbf{a}_2+z_{14}\,\mathbf{a}_3 & = & x_{14}a\,\hat{\mathbf{x}}+y_{14}b\,\hat{\mathbf{y}}+z_{14}c\,\hat{\mathbf{z}} & (4a) & \text{O V}
\end{array}
$$

$$
\begin{array}{lcccccc}
\mathbf{B}_{54} & = & -x_{14}\,\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{14}\right)\mathbf{a}_3 & = & -x_{14}a\,\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{14}\right)c\,\hat{\mathbf{z}} & (4a) & \text{O V} \\
\mathbf{B}_{55} & = & \left(\frac{1}{2} + x_{14}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{14}\right)\mathbf{a}_2 + z_{14}\,\mathbf{a}_3 & = & \left(\frac{1}{2} + x_{14}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{14}\right)b\,\hat{\mathbf{y}} + z_{14}c\,\hat{\mathbf{z}} & (4a) & \text{O V} \\
\mathbf{B}_{56} & = & \left(\frac{1}{2} - x_{14}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{14}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{14}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_{14}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{14}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{14}\right)c\,\hat{\mathbf{z}} & (4a) & \text{O V} \\
\mathbf{B}_{57} & = & x_{15}\,\mathbf{a}_1 + y_{15}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 & = & x_{15}a\,\hat{\mathbf{x}} + y_{15}b\,\hat{\mathbf{y}} + z_{15}c\,\hat{\mathbf{z}} & (4a) & \text{O VI} \\
\mathbf{B}_{58} & = & -x_{15}\,\mathbf{a}_1 - y_{15}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{15}\right)\mathbf{a}_3 & = & -x_{15}a\,\hat{\mathbf{x}} - y_{15}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{15}\right)c\,\hat{\mathbf{z}} & (4a) & \text{O VI} \\
\mathbf{B}_{59} & = & \left(\frac{1}{2} + x_{15}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{15}\right)\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 & = & \left(\frac{1}{2} + x_{15}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{15}\right)b\,\hat{\mathbf{y}} + z_{15}c\,\hat{\mathbf{z}} & (4a) & \text{O VI} \\
\mathbf{B}_{60} & = & \left(\frac{1}{2} - x_{15}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{15}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{15}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_{15}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{15}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{15}\right)c\,\hat{\mathbf{z}} & (4a) & \text{O VI} \\
\mathbf{B}_{61} & = & x_{16}\,\mathbf{a}_1 + y_{16}\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 & = & x_{16}a\,\hat{\mathbf{x}} + y_{16}b\,\hat{\mathbf{y}} + z_{16}c\,\hat{\mathbf{z}} & (4a) & \text{O VII} \\
\mathbf{B}_{62} & = & -x_{16}\,\mathbf{a}_1 - y_{16}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{16}\right)\mathbf{a}_3 & = & -x_{16}a\,\hat{\mathbf{x}} - y_{16}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{16}\right)c\,\hat{\mathbf{z}} & (4a) & \text{O VII} \\
\mathbf{B}_{63} & = & \left(\frac{1}{2} + x_{16}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{16}\right)\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 & = & \left(\frac{1}{2} + x_{16}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{16}\right)b\,\hat{\mathbf{y}} + z_{16}c\,\hat{\mathbf{z}} & (4a) & \text{O VII} \\
\mathbf{B}_{64} & = & \left(\frac{1}{2} - x_{16}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{16}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{16}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_{16}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{16}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{16}\right)c\,\hat{\mathbf{z}} & (4a) & \text{O VII} \\
\mathbf{B}_{65} & = & x_{17}\,\mathbf{a}_1 + y_{17}\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 & = & x_{17}a\,\hat{\mathbf{x}} + y_{17}b\,\hat{\mathbf{y}} + z_{17}c\,\hat{\mathbf{z}} & (4a) & \text{O VIII} \\
\mathbf{B}_{66} & = & -x_{17}\,\mathbf{a}_1 - y_{17}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{17}\right)\mathbf{a}_3 & = & -x_{17}a\,\hat{\mathbf{x}} - y_{17}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{17}\right)c\,\hat{\mathbf{z}} & (4a) & \text{O VIII} \\
\mathbf{B}_{67} & = & \left(\frac{1}{2} + x_{17}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{17}\right)\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 & = & \left(\frac{1}{2} + x_{17}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{17}\right)b\,\hat{\mathbf{y}} + z_{17}c\,\hat{\mathbf{z}} & (4a) & \text{O VIII} \\
\mathbf{B}_{68} & = & \left(\frac{1}{2} - x_{17}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{17}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{17}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_{17}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{17}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{17}\right)c\,\hat{\mathbf{z}} & (4a) & \text{O VIII} \\
\mathbf{B}_{69} & = & x_{18}\,\mathbf{a}_1 + y_{18}\,\mathbf{a}_2 + z_{18}\,\mathbf{a}_3 & = & x_{18}a\,\hat{\mathbf{x}} + y_{18}b\,\hat{\mathbf{y}} + z_{18}c\,\hat{\mathbf{z}} & (4a) & \text{O IX} \\
\mathbf{B}_{70} & = & -x_{18}\,\mathbf{a}_1 - y_{18}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{18}\right)\mathbf{a}_3 & = & -x_{18}a\,\hat{\mathbf{x}} - y_{18}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{18}\right)c\,\hat{\mathbf{z}} & (4a) & \text{O IX} \\
\mathbf{B}_{71} & = & \left(\frac{1}{2} + x_{18}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{18}\right)\mathbf{a}_2 + z_{18}\,\mathbf{a}_3 & = & \left(\frac{1}{2} + x_{18}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{18}\right)b\,\hat{\mathbf{y}} + z_{18}c\,\hat{\mathbf{z}} & (4a) & \text{O IX} \\
\mathbf{B}_{72} & = & \left(\frac{1}{2} - x_{18}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{18}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{18}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_{18}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{18}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{18}\right)c\,\hat{\mathbf{z}} & (4a) & \text{O IX} \\
\mathbf{B}_{73} & = & x_{19}\,\mathbf{a}_1 + y_{19}\,\mathbf{a}_2 + z_{19}\,\mathbf{a}_3 & = & x_{19}a\,\hat{\mathbf{x}} + y_{19}b\,\hat{\mathbf{y}} + z_{19}c\,\hat{\mathbf{z}} & (4a) & \text{O X} \\
\mathbf{B}_{74} & = & -x_{19}\,\mathbf{a}_1 - y_{19}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{19}\right)\mathbf{a}_3 & = & -x_{19}a\,\hat{\mathbf{x}} - y_{19}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{19}\right)c\,\hat{\mathbf{z}} & (4a) & \text{O X} \\
\mathbf{B}_{75} & = & \left(\frac{1}{2} + x_{19}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{19}\right)\mathbf{a}_2 + z_{19}\,\mathbf{a}_3 & = & \left(\frac{1}{2} + x_{19}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{19}\right)b\,\hat{\mathbf{y}} + z_{19}c\,\hat{\mathbf{z}} & (4a) & \text{O X} \\
\mathbf{B}_{76} & = & \left(\frac{1}{2} - x_{19}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{19}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{19}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_{19}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{19}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{19}\right)c\,\hat{\mathbf{z}} & (4a) & \text{O X} \\
\mathbf{B}_{77} & = & x_{20}\,\mathbf{a}_1 + y_{20}\,\mathbf{a}_2 + z_{20}\,\mathbf{a}_3 & = & x_{20}a\,\hat{\mathbf{x}} + y_{20}b\,\hat{\mathbf{y}} + z_{20}c\,\hat{\mathbf{z}} & (4a) & \text{O XI} \\
\mathbf{B}_{78} & = & -x_{20}\,\mathbf{a}_1 - y_{20}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{20}\right)\mathbf{a}_3 & = & -x_{20}a\,\hat{\mathbf{x}} - y_{20}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{20}\right)c\,\hat{\mathbf{z}} & (4a) & \text{O XI} \\
\mathbf{B}_{79} & = & \left(\frac{1}{2} + x_{20}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{20}\right)\mathbf{a}_2 + z_{20}\,\mathbf{a}_3 & = & \left(\frac{1}{2} + x_{20}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{20}\right)b\,\hat{\mathbf{y}} + z_{20}c\,\hat{\mathbf{z}} & (4a) & \text{O XI} \\
\mathbf{B}_{80} & = & \left(\frac{1}{2} - x_{20}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{20}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{20}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_{20}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{20}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{20}\right)c\,\hat{\mathbf{z}} & (4a) & \text{O XI} \\
\mathbf{B}_{81} & = & x_{21}\,\mathbf{a}_1 + y_{21}\,\mathbf{a}_2 + z_{21}\,\mathbf{a}_3 & = & x_{21}a\,\hat{\mathbf{x}} + y_{21}b\,\hat{\mathbf{y}} + z_{21}c\,\hat{\mathbf{z}} & (4a) & \text{O XII} \\
\mathbf{B}_{82} & = & -x_{21}\,\mathbf{a}_1 - y_{21}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{21}\right)\mathbf{a}_3 & = & -x_{21}a\,\hat{\mathbf{x}} - y_{21}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{21}\right)c\,\hat{\mathbf{z}} & (4a) & \text{O XII} \\
\mathbf{B}_{83} & = & \left(\frac{1}{2} + x_{21}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{21}\right)\mathbf{a}_2 + z_{21}\,\mathbf{a}_3 & = & \left(\frac{1}{2} + x_{21}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{21}\right)b\,\hat{\mathbf{y}} + z_{21}c\,\hat{\mathbf{z}} & (4a) & \text{O XII}
\end{array}
$$

$$\mathbf{B}_{84} = \left(\tfrac{1}{2} - x_{21}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{21}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{21}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{21}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{21}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{21}\right)c\,\hat{\mathbf{z}} \quad (4a) \quad \text{O XII}$$

References:

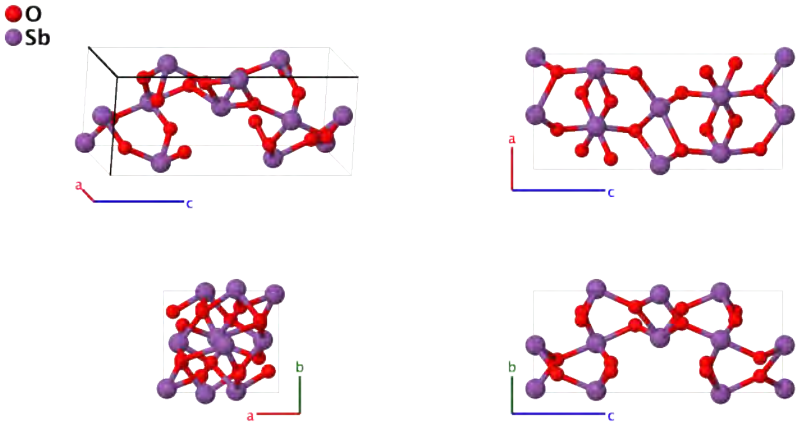
- M. Marezio, J. P. Remeika, and P. D. Dernier, *The crystal structure of the high-pressure phase CaB_2O_4(III)*, Acta Crystallogr. Sect. B Struct. Sci. **25**, 955–964 (1969), doi:10.1107/S0567740869003244.

Geometry files:

- CIF: pp. 1589
- POSCAR: pp. 1590

Cervantite (α-Sb_2O_4) Structure: A2B_oP24_33_4a_2a

http://aflow.org/prototype-encyclopedia/A2B_oP24_33_4a_2a



Prototype	:	O_2Sb
AFLOW prototype label	:	A2B_oP24_33_4a_2a
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP24
Space group number	:	33
Space group symbol	:	$Pna2_1$
AFLOW prototype command	:	`aflow --proto=A2B_oP24_33_4a_2a --params=`$a, b/a, c/a, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6$

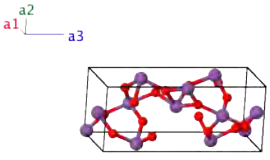
Other compounds with this structure

- $SbNbO_4$

- This is *not* the $D6_2$ structure of SbO_2, which was found to be incorrect (Thornton, 1977). There is also a naturally occurring monoclinic modification of cervantite, clinocervantite (β-Sb_2O_4).

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	($4a$)	O I
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	($4a$)	O I
$\mathbf{B}_3$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_1\right)\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_1\right) b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	($4a$)	O I

$\mathbf{B}_4$	=	$\left(\frac{1}{2}-x_1\right)\mathbf{a}_1+\left(\frac{1}{2}+y_1\right)\mathbf{a}_2+\left(\frac{1}{2}+z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_1\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_1\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_1\right)c\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1+y_2\,\mathbf{a}_2+z_2\,\mathbf{a}_3$	=	$x_2a\,\hat{\mathbf{x}}+y_2b\,\hat{\mathbf{y}}+z_2c\,\hat{\mathbf{z}}$	(4*a*)	O II
$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1-y_2\,\mathbf{a}_2+\left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}}-y_2b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_2\right)c\,\hat{\mathbf{z}}$	(4*a*)	O II
$\mathbf{B}_7$	=	$\left(\frac{1}{2}+x_2\right)\mathbf{a}_1+\left(\frac{1}{2}-y_2\right)\mathbf{a}_2+z_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_2\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_2\right)b\,\hat{\mathbf{y}}+z_2c\,\hat{\mathbf{z}}$	(4*a*)	O II
$\mathbf{B}_8$	=	$\left(\frac{1}{2}-x_2\right)\mathbf{a}_1+\left(\frac{1}{2}+y_2\right)\mathbf{a}_2+\left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_2\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_2\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_2\right)c\,\hat{\mathbf{z}}$	(4*a*)	O II
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1+y_3\,\mathbf{a}_2+z_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}}+y_3b\,\hat{\mathbf{y}}+z_3c\,\hat{\mathbf{z}}$	(4*a*)	O III
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1-y_3\,\mathbf{a}_2+\left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}}-y_3b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_3\right)c\,\hat{\mathbf{z}}$	(4*a*)	O III
$\mathbf{B}_{11}$	=	$\left(\frac{1}{2}+x_3\right)\mathbf{a}_1+\left(\frac{1}{2}-y_3\right)\mathbf{a}_2+z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_3\right)b\,\hat{\mathbf{y}}+z_3c\,\hat{\mathbf{z}}$	(4*a*)	O III
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2}-x_3\right)\mathbf{a}_1+\left(\frac{1}{2}+y_3\right)\mathbf{a}_2+\left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_3\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_3\right)c\,\hat{\mathbf{z}}$	(4*a*)	O III
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1+y_4\,\mathbf{a}_2+z_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}}+y_4b\,\hat{\mathbf{y}}+z_4c\,\hat{\mathbf{z}}$	(4*a*)	O IV
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1-y_4\,\mathbf{a}_2+\left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}}-y_4b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_4\right)c\,\hat{\mathbf{z}}$	(4*a*)	O IV
$\mathbf{B}_{15}$	=	$\left(\frac{1}{2}+x_4\right)\mathbf{a}_1+\left(\frac{1}{2}-y_4\right)\mathbf{a}_2+z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_4\right)b\,\hat{\mathbf{y}}+z_4c\,\hat{\mathbf{z}}$	(4*a*)	O IV
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2}-x_4\right)\mathbf{a}_1+\left(\frac{1}{2}+y_4\right)\mathbf{a}_2+\left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_4\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_4\right)c\,\hat{\mathbf{z}}$	(4*a*)	O IV
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1+y_5\,\mathbf{a}_2+z_5\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}}+y_5b\,\hat{\mathbf{y}}+z_5c\,\hat{\mathbf{z}}$	(4*a*)	Sb I
$\mathbf{B}_{18}$	=	$-x_5\,\mathbf{a}_1-y_5\,\mathbf{a}_2+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}}-y_5b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	(4*a*)	Sb I
$\mathbf{B}_{19}$	=	$\left(\frac{1}{2}+x_5\right)\mathbf{a}_1+\left(\frac{1}{2}-y_5\right)\mathbf{a}_2+z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_5\right)b\,\hat{\mathbf{y}}+z_5c\,\hat{\mathbf{z}}$	(4*a*)	Sb I
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2}-x_5\right)\mathbf{a}_1+\left(\frac{1}{2}+y_5\right)\mathbf{a}_2+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_5\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	(4*a*)	Sb I
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1+y_6\,\mathbf{a}_2+z_6\,\mathbf{a}_3$	=	$x_6a\,\hat{\mathbf{x}}+y_6b\,\hat{\mathbf{y}}+z_6c\,\hat{\mathbf{z}}$	(4*a*)	Sb II
$\mathbf{B}_{22}$	=	$-x_6\,\mathbf{a}_1-y_6\,\mathbf{a}_2+\left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	=	$-x_6a\,\hat{\mathbf{x}}-y_6b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	(4*a*)	Sb II
$\mathbf{B}_{23}$	=	$\left(\frac{1}{2}+x_6\right)\mathbf{a}_1+\left(\frac{1}{2}-y_6\right)\mathbf{a}_2+z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_6\right)b\,\hat{\mathbf{y}}+z_6c\,\hat{\mathbf{z}}$	(4*a*)	Sb II
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2}-x_6\right)\mathbf{a}_1+\left(\frac{1}{2}+y_6\right)\mathbf{a}_2+\left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_6\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	(4*a*)	Sb II

References:

- G. Thornton, *A Neutron Diffraction Study of α-Sb_2O_3*, Acta Crystallogr. Sect. B Struct. Sci. **33**, 1271–1273 (1977), doi:10.1107/S0567740877005822.

Found in:

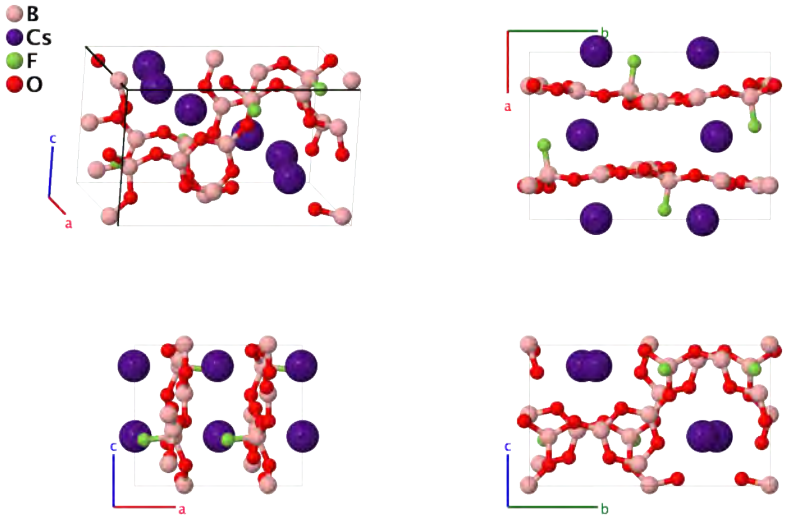
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1590
- POSCAR: pp. 1591

CsB_4O_6F Structure: A4BCD6_oP48_33_4a_a_a_6a

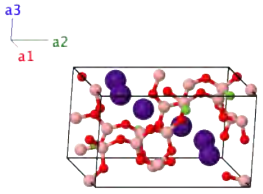
http://aflow.org/prototype-encyclopedia/A4BCD6_oP48_33_4a_a_a_6a



Prototype	:	B_4CsFO_6
AFLOW prototype label	:	A4BCD6_oP48_33_4a_a_a_6a
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP48
Space group number	:	33
Space group symbol	:	$Pna2_1$
AFLOW prototype command	:	`aflow --proto=A4BCD6_oP48_33_4a_a_a_6a --params=`$a, b/a, c/a, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}$

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	($4a$)	B I
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	($4a$)	B I
$\mathbf{B}_3$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_1\right)\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_1\right) b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	($4a$)	B I
$\mathbf{B}_4$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_1\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_1\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	($4a$)	B I

$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*a*)	B II
$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	(4*a*)	B II
$\mathbf{B}_7$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right) b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*a*)	B II
$\mathbf{B}_8$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	(4*a*)	B II
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*a*)	B III
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	(4*a*)	B III
$\mathbf{B}_{11}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right) b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*a*)	B III
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	(4*a*)	B III
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(4*a*)	B IV
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	(4*a*)	B IV
$\mathbf{B}_{15}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right) b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(4*a*)	B IV
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	(4*a*)	B IV
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(4*a*)	Cs
$\mathbf{B}_{18}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	(4*a*)	Cs
$\mathbf{B}_{19}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right) b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(4*a*)	Cs
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	(4*a*)	Cs
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(4*a*)	F
$\mathbf{B}_{22}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	(4*a*)	F
$\mathbf{B}_{23}$	=	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right) b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(4*a*)	F
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	(4*a*)	F
$\mathbf{B}_{25}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_{26}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right) c\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_{27}$	=	$\left(\frac{1}{2} + x_7\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_7\right)\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_7\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_7\right) b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} - x_7\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_7\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right) c\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_{29}$	=	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(4*a*)	O II
$\mathbf{B}_{30}$	=	$-x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right) c\,\hat{\mathbf{z}}$	(4*a*)	O II
$\mathbf{B}_{31}$	=	$\left(\frac{1}{2} + x_8\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_8\right)\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_8\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_8\right) b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(4*a*)	O II
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2} - x_8\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_8\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_8\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_8\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right) c\,\hat{\mathbf{z}}$	(4*a*)	O II
$\mathbf{B}_{33}$	=	$x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(4*a*)	O III
$\mathbf{B}_{34}$	=	$-x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right) c\,\hat{\mathbf{z}}$	(4*a*)	O III
$\mathbf{B}_{35}$	=	$\left(\frac{1}{2} + x_9\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_9\right)\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_9\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_9\right) b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(4*a*)	O III

$\mathbf{B}_{36}$	=	$\left(\frac{1}{2}-x_9\right)\mathbf{a}_1+\left(\frac{1}{2}+y_9\right)\mathbf{a}_2+\left(\frac{1}{2}+z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_9\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_9\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_9\right)c\,\hat{\mathbf{z}}$	(4*a*)	O III
$\mathbf{B}_{37}$	=	$x_{10}\,\mathbf{a}_1+y_{10}\,\mathbf{a}_2+z_{10}\,\mathbf{a}_3$	=	$x_{10}a\,\hat{\mathbf{x}}+y_{10}b\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}}$	(4*a*)	O IV
$\mathbf{B}_{38}$	=	$-x_{10}\,\mathbf{a}_1-y_{10}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{10}\right)\mathbf{a}_3$	=	$-x_{10}a\,\hat{\mathbf{x}}-y_{10}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{10}\right)c\,\hat{\mathbf{z}}$	(4*a*)	O IV
$\mathbf{B}_{39}$	=	$\left(\frac{1}{2}+x_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{10}\right)\mathbf{a}_2+z_{10}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{10}\right)b\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}}$	(4*a*)	O IV
$\mathbf{B}_{40}$	=	$\left(\frac{1}{2}-x_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{10}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{10}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{10}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{10}\right)c\,\hat{\mathbf{z}}$	(4*a*)	O IV
$\mathbf{B}_{41}$	=	$x_{11}\,\mathbf{a}_1+y_{11}\,\mathbf{a}_2+z_{11}\,\mathbf{a}_3$	=	$x_{11}a\,\hat{\mathbf{x}}+y_{11}b\,\hat{\mathbf{y}}+z_{11}c\,\hat{\mathbf{z}}$	(4*a*)	O V
$\mathbf{B}_{42}$	=	$-x_{11}\,\mathbf{a}_1-y_{11}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{11}\right)\mathbf{a}_3$	=	$-x_{11}a\,\hat{\mathbf{x}}-y_{11}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{11}\right)c\,\hat{\mathbf{z}}$	(4*a*)	O V
$\mathbf{B}_{43}$	=	$\left(\frac{1}{2}+x_{11}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{11}\right)\mathbf{a}_2+z_{11}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{11}\right)b\,\hat{\mathbf{y}}+z_{11}c\,\hat{\mathbf{z}}$	(4*a*)	O V
$\mathbf{B}_{44}$	=	$\left(\frac{1}{2}-x_{11}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{11}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{11}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{11}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{11}\right)c\,\hat{\mathbf{z}}$	(4*a*)	O V
$\mathbf{B}_{45}$	=	$x_{12}\,\mathbf{a}_1+y_{12}\,\mathbf{a}_2+z_{12}\,\mathbf{a}_3$	=	$x_{12}a\,\hat{\mathbf{x}}+y_{12}b\,\hat{\mathbf{y}}+z_{12}c\,\hat{\mathbf{z}}$	(4*a*)	O VI
$\mathbf{B}_{46}$	=	$-x_{12}\,\mathbf{a}_1-y_{12}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{12}\right)\mathbf{a}_3$	=	$-x_{12}a\,\hat{\mathbf{x}}-y_{12}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{12}\right)c\,\hat{\mathbf{z}}$	(4*a*)	O VI
$\mathbf{B}_{47}$	=	$\left(\frac{1}{2}+x_{12}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{12}\right)\mathbf{a}_2+z_{12}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{12}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{12}\right)b\,\hat{\mathbf{y}}+z_{12}c\,\hat{\mathbf{z}}$	(4*a*)	O VI
$\mathbf{B}_{48}$	=	$\left(\frac{1}{2}-x_{12}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{12}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{12}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{12}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{12}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{12}\right)c\,\hat{\mathbf{z}}$	(4*a*)	O VI

References:

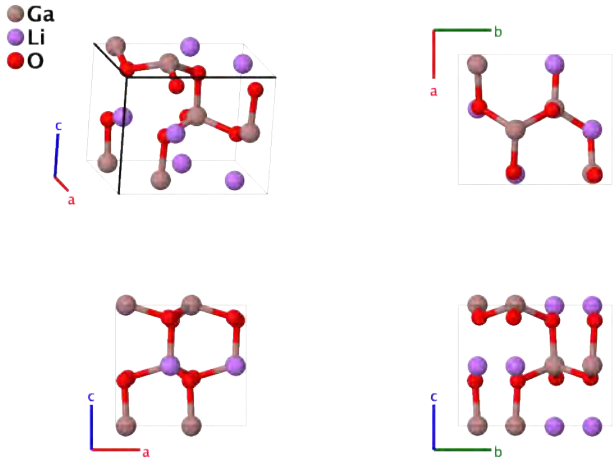
- X. Wang, Y. Wang, B. Zhang, F. Zhang, Z. Yang, and S. Pan, *CsB_4O_6F: A Congruent-Melting Deep-Ultraviolet Nonlinear Optical Material by Combining Superior Functional Units*, Angew. Chem. **129**, 14307–14311 (2017), doi:10.1002/ange.201708231.

Geometry files:

- CIF: pp. 1591
- POSCAR: pp. 1591

$LiGaO_2$ Structure: ABC2_oP16_33_a_a_2a

http://aflow.org/prototype-encyclopedia/ABC2_oP16_33_a_a_2a

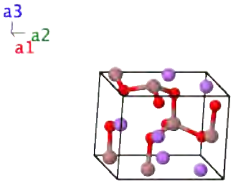


Prototype	:	$GaLiO_2$
AFLOW prototype label	:	ABC2_oP16_33_a_a_2a
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP16
Space group number	:	33
Space group symbol	:	$Pna2_1$
AFLOW prototype command	:	`aflow --proto=ABC2_oP16_33_a_a_2a --params=`$a, b/a, c/a, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4$

- Space group $Pna2_1$ #33 allows an arbitary choice of the zero of the z-axis. Here we set $z_1 = 0$ for gallium.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	$(4a)$	Ga
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	$(4a)$	Ga
$\mathbf{B}_3$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_1\right)\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_1\right) b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	$(4a)$	Ga
$\mathbf{B}_4$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_1\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_1\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	$(4a)$	Ga

$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*a*)	Li
$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	(4*a*)	Li
$\mathbf{B}_7$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right) b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*a*)	Li
$\mathbf{B}_8$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	(4*a*)	Li
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_{11}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right) b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(4*a*)	O II
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	(4*a*)	O II
$\mathbf{B}_{15}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right) b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(4*a*)	O II
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	(4*a*)	O II

References:

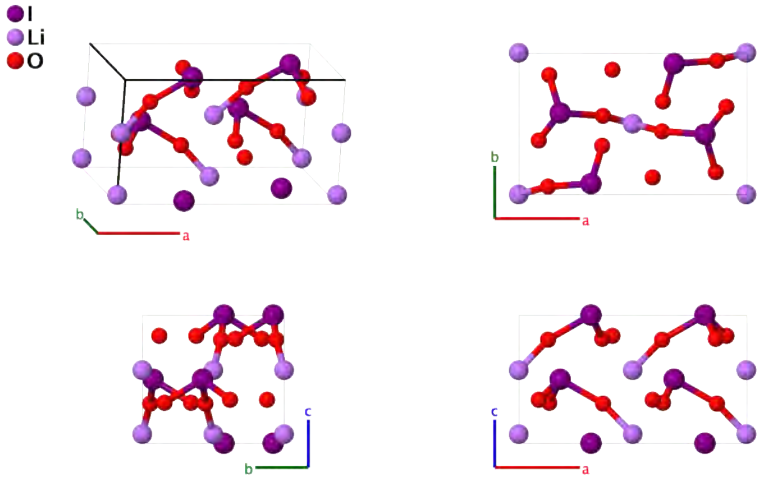
- M. Marezio, *The Crystal Structure of $LiGaO_2$*, Acta Cryst. **18**, 481–484 (1965), doi:10.1107/S0365110X65001068.

Geometry files:

- CIF: pp. 1591
- POSCAR: pp. 1592

γ-$LiIO_3$ Structure: ABC3_oP20_33_a_a_3a

http://aflow.org/prototype-encyclopedia/ABC3_oP20_33_a_a_3a

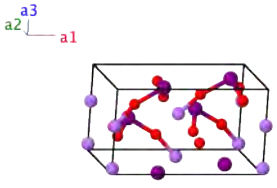


Prototype	:	$ILiO_3$
AFLOW prototype label	:	ABC3_oP20_33_a_a_3a
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP20
Space group number	:	33
Space group symbol	:	$Pna2_1$
AFLOW prototype command	:	`aflow --proto=ABC3_oP20_33_a_a_3a --params=`$a, b/a, c/a, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5$

- $LiIO_3$ is known to exist in three forms:
  - α-$LiIO_3$, stable below 473 K: (Zachariasen, 1931) originally determined that the structure of α-$LiIO_3$ was in space group $P6_322$ #182, which (Hermann, 1937) designated *Strukturbericht* $E2_3$. (Rosenzweig, 1966) subsequently determined that the true structure was in space group $P6_3$ #173.
  - β-$LiIO_3$, stable from 573 K up to the melting point at 708 K.
  - γ-$LiIO_3$, stable between the α- and β-phases. Here we show the structure at 515 K.
- Space group $Pna2_1$ #33 allows an arbitary choice of the zero of the z-axis. Here we set $z_1 = 0$ for iodine.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(4*a*)	I
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	(4*a*)	I
$\mathbf{B}_3$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_1\right)\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_1\right) b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(4*a*)	I
$\mathbf{B}_4$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_1\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_1\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	(4*a*)	I
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*a*)	Li
$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	(4*a*)	Li
$\mathbf{B}_7$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right) b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*a*)	Li
$\mathbf{B}_8$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	(4*a*)	Li
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_{11}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right) b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(4*a*)	O II
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	(4*a*)	O II
$\mathbf{B}_{15}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right) b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(4*a*)	O II
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	(4*a*)	O II
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(4*a*)	O III
$\mathbf{B}_{18}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	(4*a*)	O III
$\mathbf{B}_{19}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right) b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(4*a*)	O III
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	(4*a*)	O III

References:

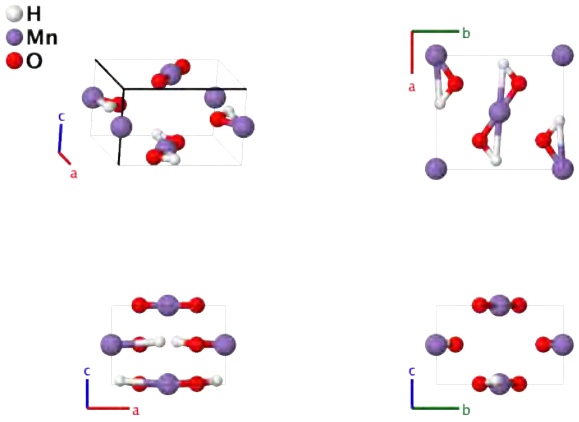
- R. Liminga, C. Svensson, J. Albertsson, and S. C. Abrahams, *Gamma-lithium iodate structure at 515 K and the α-$LiIO_3$ to γ-$LiIO_3$, γ-$LiIO_3$ to β-$LiIO_3$ phase transitions*, J. Chem. Phys. **77**, 4222–4226 (1982), doi:10.1063/1.444332.
- W. H. Zachariasen and F. A. Barta, *Crystal Structure of Lithium Iodate*, Phys. Rev. **37**, 1626–1630 (1931), doi:10.1103/PhysRev.37.1626.
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).
- A. Rosenzweig and B. Morosin, *A reinvestigation of the crystal structure of $LiIO_3$*, Acta Cryst. **20**, 758–761 (1966), doi:10.1107/S0365110X66001804.

Geometry files:

- CIF: pp. 1592
- POSCAR: pp. 1592

$MnF_{2-x}(OH)_x$ Structure: A2BC2_oP10_34_c_a_c

http://aflow.org/prototype-encyclopedia/A2BC2_oP10_34_c_a_c

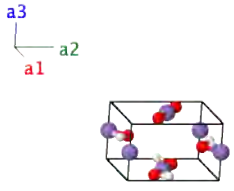


Prototype	:	$F_{2-x}H_xMnO_x$
AFLOW prototype label	:	A2BC2_oP10_34_c_a_c
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP10
Space group number	:	34
Space group symbol	:	*Pnn*2
AFLOW prototype command	:	`aflow --proto=A2BC2_oP10_34_c_a_c --params=`$a, b/a, c/a, z_1, x_2, y_2, z_2, x_3, y_3, z_3$

- We use the data (Yahia, 2013) give for the site occupancy of oxygen and hydrogen (0.399). The oxygen ($4c$) atom can be replaced by fluorine, in which case the hydrogen ($4c$) atom is not present. Approximately 80% of the first ($4c$) site is oxygen, with the remainder being hydrogen.
- Space group *Pnn*2 #34 allows an arbitrary choice of the origin of the z-coordinate. Here we use this to put the manganese atom at the origin, setting $z_1 = 0$.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$z_1\,\mathbf{a}_3$	=	$z_1 c\,\hat{\mathbf{z}}$	$(2a)$	Mn
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\,\hat{\mathbf{z}}$	$(2a)$	Mn

$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*c*)	H
$\mathbf{B}_4$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*c*)	H
$\mathbf{B}_5$	=	$\left(\frac{1}{2}+x_2\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_2\right)c\,\hat{\mathbf{z}}$	(4*c*)	H
$\mathbf{B}_6$	=	$\left(\frac{1}{2}-x_2\right)\mathbf{a}_1 + \left(\frac{1}{2}+y_2\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_2\right)c\,\hat{\mathbf{z}}$	(4*c*)	H
$\mathbf{B}_7$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*c*)	O
$\mathbf{B}_8$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*c*)	O
$\mathbf{B}_9$	=	$\left(\frac{1}{2}+x_3\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_3\right)c\,\hat{\mathbf{z}}$	(4*c*)	O
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2}-x_3\right)\mathbf{a}_1 + \left(\frac{1}{2}+y_3\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_3\right)c\,\hat{\mathbf{z}}$	(4*c*)	O

References:

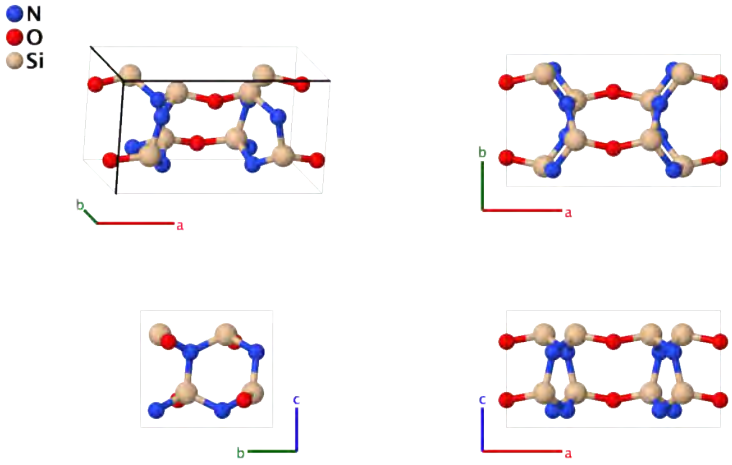
- H. B. Yahia, M. Shikano, H. Kobayashi, M. Avdeev, S. Liu, and C. D. Ling, *Synthesis and characterization of the crystal structure and magnetic properties of the hydroxyfluoride $MnF_{2-x}(OH)_x$ ($x \approx 0.8$)*, Phys. Chem. Chem. Phys. **15**, 13061–13069 (2013), doi:10.1039/C3CP50740H.

Geometry files:

- CIF: pp. 1592
- POSCAR: pp. 1593

Si_2N_2O Structure: A2BC2_oC20_36_b_a_b

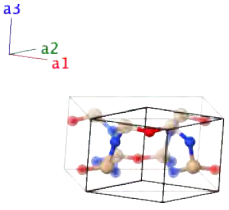
http://aflow.org/prototype-encyclopedia/A2BC2_oC20_36_b_a_b



Prototype	:	N_2OSi_2
AFLOW prototype label	:	A2BC2_oC20_36_b_a_b
***Strukturbericht* designation**	:	None
Pearson symbol	:	oC20
Space group number	:	36
Space group symbol	:	$Cmc2_1$
AFLOW prototype command	:	`aflow --proto=A2BC2_oC20_36_b_a_b --params=`$a, b/a, c/a, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3$

Base-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$y_1 b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	$(4a)$	O
$\mathbf{B}_2$	=	$y_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$-y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\,\hat{\mathbf{z}}$	$(4a)$	O
$\mathbf{B}_3$	=	$(x_2 - y_2)\,\mathbf{a}_1 + (x_2 + y_2)\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(8b)$	N
$\mathbf{B}_4$	=	$(-x_2 + y_2)\,\mathbf{a}_1 + (-x_2 - y_2)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	$(8b)$	N
$\mathbf{B}_5$	=	$(x_2 + y_2)\,\mathbf{a}_1 + (x_2 - y_2)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	$(8b)$	N
$\mathbf{B}_6$	=	$(-x_2 - y_2)\,\mathbf{a}_1 + (-x_2 + y_2)\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(8b)$	N

$$\mathbf{B}_7 = (x_3 - y_3)\,\mathbf{a}_1 + (x_3 + y_3)\,\mathbf{a}_2 + z_3\,\mathbf{a}_3 = x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}} \quad (8b) \quad \text{Si}$$

$$\mathbf{B}_8 = (-x_3 + y_3)\,\mathbf{a}_1 + (-x_3 - y_3)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_3\right)\mathbf{a}_3 = -x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_3\right) c\,\hat{\mathbf{z}} \quad (8b) \quad \text{Si}$$

$$\mathbf{B}_9 = (x_3 + y_3)\,\mathbf{a}_1 + (x_3 - y_3)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_3\right)\mathbf{a}_3 = x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_3\right) c\,\hat{\mathbf{z}} \quad (8b) \quad \text{Si}$$

$$\mathbf{B}_{10} = (-x_3 - y_3)\,\mathbf{a}_1 + (-x_3 + y_3)\,\mathbf{a}_2 + z_3\,\mathbf{a}_3 = -x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}} \quad (8b) \quad \text{Si}$$

References:

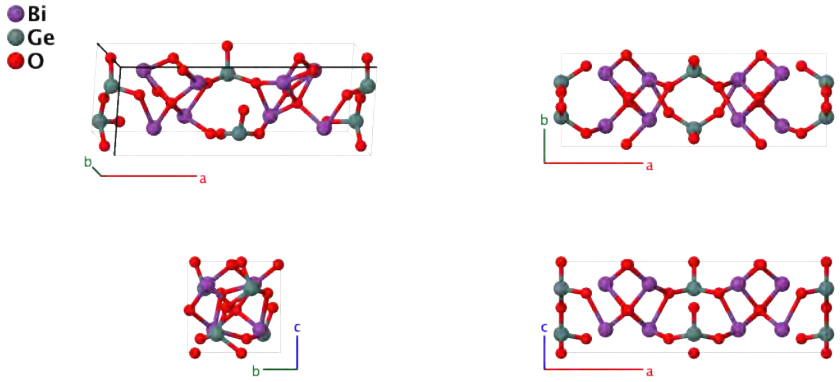
- I. Idrestedt and C. Brosset, *Structure of Si_2N_2O*, Acta Chem. Scand. **18**, 1879–1886 (1964), doi:10.3891/acta.chem.scand.18-1879.

Geometry files:

- CIF: pp. 1593
- POSCAR: pp. 1593

Bi_2GeO_5 Structure: A2BC5_oC32_36_b_a_a2b

http://aflow.org/prototype-encyclopedia/A2BC5_oC32_36_b_a_a2b



Prototype	:	Bi_2GeO_5
AFLOW prototype label	:	A2BC5_oC32_36_b_a_a2b
***Strukturbericht* designation**	:	None
Pearson symbol	:	oC32
Space group number	:	36
Space group symbol	:	$Cmc2_1$
AFLOW prototype command	:	`aflow --proto=A2BC5_oC32_36_b_a_a2b --params=`$a, b/a, c/a, y_1, z_1, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5$

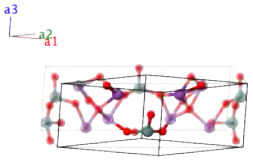
Other compounds with this structure

- Bi_2SiO_5

- Space group $Cmc2_1$ #36 allows an arbitrary choice of the zero of the z-axis. Here it is chosen so that $z_3 = 0$ for the bismuth atom.

Base-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$y_1 b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	($4a$)	Ge
$\mathbf{B}_2$	=	$y_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$-y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	($4a$)	Ge
$\mathbf{B}_3$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	($4a$)	O I
$\mathbf{B}_4$	=	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$-y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	($4a$)	O I
$\mathbf{B}_5$	=	$(x_3 - y_3)\,\mathbf{a}_1 + (x_3 + y_3)\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	($8b$)	Bi

$\mathbf{B}_6$	=	$(-x_3+y_3)\,\mathbf{a}_1+(-x_3-y_3)\,\mathbf{a}_2+\left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}}-y_3b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_3\right)c\,\hat{\mathbf{z}}$	(8*b*)	Bi
$\mathbf{B}_7$	=	$(x_3+y_3)\,\mathbf{a}_1+(x_3-y_3)\,\mathbf{a}_2+\left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}}-y_3b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_3\right)c\,\hat{\mathbf{z}}$	(8*b*)	Bi
$\mathbf{B}_8$	=	$(-x_3-y_3)\,\mathbf{a}_1+(-x_3+y_3)\,\mathbf{a}_2+z_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}}+y_3b\,\hat{\mathbf{y}}+z_3c\,\hat{\mathbf{z}}$	(8*b*)	Bi
$\mathbf{B}_9$	=	$(x_4-y_4)\,\mathbf{a}_1+(x_4+y_4)\,\mathbf{a}_2+z_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}}+y_4b\,\hat{\mathbf{y}}+z_4c\,\hat{\mathbf{z}}$	(8*b*)	O II
$\mathbf{B}_{10}$	=	$(-x_4+y_4)\,\mathbf{a}_1+(-x_4-y_4)\,\mathbf{a}_2+\left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}}-y_4b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_4\right)c\,\hat{\mathbf{z}}$	(8*b*)	O II
$\mathbf{B}_{11}$	=	$(x_4+y_4)\,\mathbf{a}_1+(x_4-y_4)\,\mathbf{a}_2+\left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}}-y_4b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_4\right)c\,\hat{\mathbf{z}}$	(8*b*)	O II
$\mathbf{B}_{12}$	=	$(-x_4-y_4)\,\mathbf{a}_1+(-x_4+y_4)\,\mathbf{a}_2+z_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}}+y_4b\,\hat{\mathbf{y}}+z_4c\,\hat{\mathbf{z}}$	(8*b*)	O II
$\mathbf{B}_{13}$	=	$(x_5-y_5)\,\mathbf{a}_1+(x_5+y_5)\,\mathbf{a}_2+z_5\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}}+y_5b\,\hat{\mathbf{y}}+z_5c\,\hat{\mathbf{z}}$	(8*b*)	O III
$\mathbf{B}_{14}$	=	$(-x_5+y_5)\,\mathbf{a}_1+(-x_5-y_5)\,\mathbf{a}_2+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}}-y_5b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	(8*b*)	O III
$\mathbf{B}_{15}$	=	$(x_5+y_5)\,\mathbf{a}_1+(x_5-y_5)\,\mathbf{a}_2+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}}-y_5b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	(8*b*)	O III
$\mathbf{B}_{16}$	=	$(-x_5-y_5)\,\mathbf{a}_1+(-x_5+y_5)\,\mathbf{a}_2+z_5\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}}+y_5b\,\hat{\mathbf{y}}+z_5c\,\hat{\mathbf{z}}$	(8*b*)	O III

References:

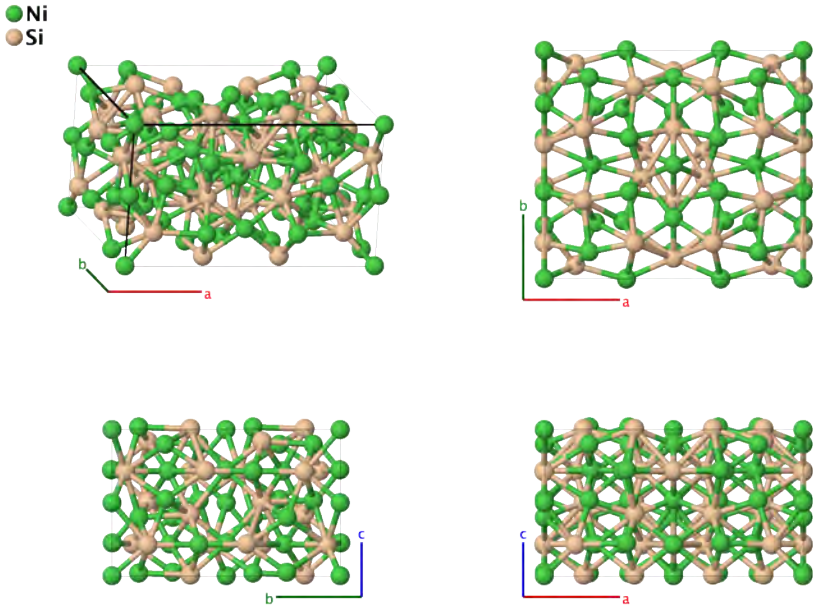
- B. Aurivillius, C.-I. Lindblom, and P. Sténson, *The Crystal Structure of Bi_2GeO_5*, Acta Chem. Scand. **18**, 1555–1557 (1964), doi:10.3891/acta.chem.scand.18-1555.

Geometry files:

- CIF: pp. 1593
- POSCAR: pp. 1594

Ni_3Si_2 Structure: A3B2_oC80_36_4a4b_2a3b

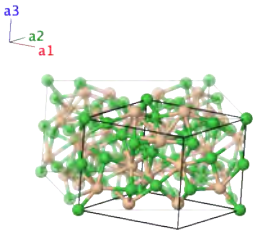
http://aflow.org/prototype-encyclopedia/A3B2_oC80_36_4a4b_2a3b



Prototype	:	Ni_3Si_2
AFLOW prototype label	:	A3B2_oC80_36_4a4b_2a3b
***Strukturbericht* designation**	:	None
Pearson symbol	:	oC80
Space group number	:	36
Space group symbol	:	$Cmc2_1$
AFLOW prototype command	:	`aflow --proto=A3B2_oC80_36_4a4b_2a3b --params=`$a, b/a, c/a, y_1, z_1, y_2, z_2, y_3, z_3, y_4, z_4, y_5, z_5, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}$

Base-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$y_1 b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	$(4a)$	Ni I
$\mathbf{B}_2$	=	$y_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$-y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\,\hat{\mathbf{z}}$	$(4a)$	Ni I

$$
\begin{array}{lcccccc}
\mathbf{B}_3 & = & -y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3 & = & y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}} & (4a) & \text{Ni II} \\
\mathbf{B}_4 & = & y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_2\right)\mathbf{a}_3 & = & -y_2 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_2\right) c\,\hat{\mathbf{z}} & (4a) & \text{Ni II} \\
\mathbf{B}_5 & = & -y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3 & = & y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}} & (4a) & \text{Ni III} \\
\mathbf{B}_6 & = & y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_3\right)\mathbf{a}_3 & = & -y_3 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_3\right) c\,\hat{\mathbf{z}} & (4a) & \text{Ni III} \\
\mathbf{B}_7 & = & -y_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3 & = & y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}} & (4a) & \text{Ni IV} \\
\mathbf{B}_8 & = & y_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_4\right)\mathbf{a}_3 & = & -y_4 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_4\right) c\,\hat{\mathbf{z}} & (4a) & \text{Ni IV} \\
\mathbf{B}_9 & = & -y_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3 & = & y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}} & (4a) & \text{Si I} \\
\mathbf{B}_{10} & = & y_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_3 & = & -y_5 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_5\right) c\,\hat{\mathbf{z}} & (4a) & \text{Si I} \\
\mathbf{B}_{11} & = & -y_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 & = & y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}} & (4a) & \text{Si II} \\
\mathbf{B}_{12} & = & y_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 & = & -y_6 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right) c\,\hat{\mathbf{z}} & (4a) & \text{Si II} \\
\mathbf{B}_{13} & = & (x_7 - y_7)\,\mathbf{a}_1 + (x_7 + y_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 & = & x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}} & (8b) & \text{Ni V} \\
\mathbf{B}_{14} & = & (-x_7 + y_7)\,\mathbf{a}_1 + (-x_7 - y_7)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_7\right)\mathbf{a}_3 & = & -x_7 a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_7\right) c\,\hat{\mathbf{z}} & (8b) & \text{Ni V} \\
\mathbf{B}_{15} & = & (x_7 + y_7)\,\mathbf{a}_1 + (x_7 - y_7)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_7\right)\mathbf{a}_3 & = & x_7 a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_7\right) c\,\hat{\mathbf{z}} & (8b) & \text{Ni V} \\
\mathbf{B}_{16} & = & (-x_7 - y_7)\,\mathbf{a}_1 + (-x_7 + y_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 & = & -x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}} & (8b) & \text{Ni V} \\
\mathbf{B}_{17} & = & (x_8 - y_8)\,\mathbf{a}_1 + (x_8 + y_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 & = & x_8 a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}} & (8b) & \text{Ni VI} \\
\mathbf{B}_{18} & = & (-x_8 + y_8)\,\mathbf{a}_1 + (-x_8 - y_8)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_8\right)\mathbf{a}_3 & = & -x_8 a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_8\right) c\,\hat{\mathbf{z}} & (8b) & \text{Ni VI} \\
\mathbf{B}_{19} & = & (x_8 + y_8)\,\mathbf{a}_1 + (x_8 - y_8)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_8\right)\mathbf{a}_3 & = & x_8 a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_8\right) c\,\hat{\mathbf{z}} & (8b) & \text{Ni VI} \\
\mathbf{B}_{20} & = & (-x_8 - y_8)\,\mathbf{a}_1 + (-x_8 + y_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 & = & -x_8 a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}} & (8b) & \text{Ni VI} \\
\mathbf{B}_{21} & = & (x_9 - y_9)\,\mathbf{a}_1 + (x_9 + y_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 & = & x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}} & (8b) & \text{Ni VII} \\
\mathbf{B}_{22} & = & (-x_9 + y_9)\,\mathbf{a}_1 + (-x_9 - y_9)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_3 & = & -x_9 a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_9\right) c\,\hat{\mathbf{z}} & (8b) & \text{Ni VII} \\
\mathbf{B}_{23} & = & (x_9 + y_9)\,\mathbf{a}_1 + (x_9 - y_9)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_3 & = & x_9 a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_9\right) c\,\hat{\mathbf{z}} & (8b) & \text{Ni VII} \\
\mathbf{B}_{24} & = & (-x_9 - y_9)\,\mathbf{a}_1 + (-x_9 + y_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 & = & -x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}} & (8b) & \text{Ni VII} \\
\mathbf{B}_{25} & = & (x_{10} - y_{10})\,\mathbf{a}_1 + (x_{10} + y_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 & = & x_{10} a\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}} & (8b) & \text{Ni VIII} \\
\mathbf{B}_{26} & = & (-x_{10} + y_{10})\,\mathbf{a}_1 + (-x_{10} - y_{10})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{10}\right)\mathbf{a}_3 & = & -x_{10} a\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{10}\right) c\,\hat{\mathbf{z}} & (8b) & \text{Ni VIII} \\
\mathbf{B}_{27} & = & (x_{10} + y_{10})\,\mathbf{a}_1 + (x_{10} - y_{10})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{10}\right)\mathbf{a}_3 & = & x_{10} a\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{10}\right) c\,\hat{\mathbf{z}} & (8b) & \text{Ni VIII} \\
\mathbf{B}_{28} & = & (-x_{10} - y_{10})\,\mathbf{a}_1 + (-x_{10} + y_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 & = & -x_{10} a\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}} & (8b) & \text{Ni VIII} \\
\mathbf{B}_{29} & = & (x_{11} - y_{11})\,\mathbf{a}_1 + (x_{11} + y_{11})\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 & = & x_{11} a\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + z_{11} c\,\hat{\mathbf{z}} & (8b) & \text{Si III} \\
\mathbf{B}_{30} & = & (-x_{11} + y_{11})\,\mathbf{a}_1 + (-x_{11} - y_{11})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{11}\right)\mathbf{a}_3 & = & -x_{11} a\,\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{11}\right) c\,\hat{\mathbf{z}} & (8b) & \text{Si III} \\
\mathbf{B}_{31} & = & (x_{11} + y_{11})\,\mathbf{a}_1 + (x_{11} - y_{11})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{11}\right)\mathbf{a}_3 & = & x_{11} a\,\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{11}\right) c\,\hat{\mathbf{z}} & (8b) & \text{Si III}
\end{array}
$$

$\mathbf{B}_{32}$	=	$(-x_{11}-y_{11})\,\mathbf{a}_1+(-x_{11}+y_{11})\,\mathbf{a}_2+z_{11}\,\mathbf{a}_3$	=	$-x_{11}a\,\hat{\mathbf{x}}+y_{11}b\,\hat{\mathbf{y}}+z_{11}c\,\hat{\mathbf{z}}$	(8*b*)	Si III
$\mathbf{B}_{33}$	=	$(x_{12}-y_{12})\,\mathbf{a}_1+(x_{12}+y_{12})\,\mathbf{a}_2+z_{12}\,\mathbf{a}_3$	=	$x_{12}a\,\hat{\mathbf{x}}+y_{12}b\,\hat{\mathbf{y}}+z_{12}c\,\hat{\mathbf{z}}$	(8*b*)	Si IV
$\mathbf{B}_{34}$	=	$(-x_{12}+y_{12})\,\mathbf{a}_1+(-x_{12}-y_{12})\,\mathbf{a}_2+\left(\frac{1}{2}+z_{12}\right)\mathbf{a}_3$	=	$-x_{12}a\,\hat{\mathbf{x}}-y_{12}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{12}\right)c\,\hat{\mathbf{z}}$	(8*b*)	Si IV
$\mathbf{B}_{35}$	=	$(x_{12}+y_{12})\,\mathbf{a}_1+(x_{12}-y_{12})\,\mathbf{a}_2+\left(\frac{1}{2}+z_{12}\right)\mathbf{a}_3$	=	$x_{12}a\,\hat{\mathbf{x}}-y_{12}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{12}\right)c\,\hat{\mathbf{z}}$	(8*b*)	Si IV
$\mathbf{B}_{36}$	=	$(-x_{12}-y_{12})\,\mathbf{a}_1+(-x_{12}+y_{12})\,\mathbf{a}_2+z_{12}\,\mathbf{a}_3$	=	$-x_{12}a\,\hat{\mathbf{x}}+y_{12}b\,\hat{\mathbf{y}}+z_{12}c\,\hat{\mathbf{z}}$	(8*b*)	Si IV
$\mathbf{B}_{37}$	=	$(x_{13}-y_{13})\,\mathbf{a}_1+(x_{13}+y_{13})\,\mathbf{a}_2+z_{13}\,\mathbf{a}_3$	=	$x_{13}a\,\hat{\mathbf{x}}+y_{13}b\,\hat{\mathbf{y}}+z_{13}c\,\hat{\mathbf{z}}$	(8*b*)	Si V
$\mathbf{B}_{38}$	=	$(-x_{13}+y_{13})\,\mathbf{a}_1+(-x_{13}-y_{13})\,\mathbf{a}_2+\left(\frac{1}{2}+z_{13}\right)\mathbf{a}_3$	=	$-x_{13}a\,\hat{\mathbf{x}}-y_{13}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{13}\right)c\,\hat{\mathbf{z}}$	(8*b*)	Si V
$\mathbf{B}_{39}$	=	$(x_{13}+y_{13})\,\mathbf{a}_1+(x_{13}-y_{13})\,\mathbf{a}_2+\left(\frac{1}{2}+z_{13}\right)\mathbf{a}_3$	=	$x_{13}a\,\hat{\mathbf{x}}-y_{13}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{13}\right)c\,\hat{\mathbf{z}}$	(8*b*)	Si V
$\mathbf{B}_{40}$	=	$(-x_{13}-y_{13})\,\mathbf{a}_1+(-x_{13}+y_{13})\,\mathbf{a}_2+z_{13}\,\mathbf{a}_3$	=	$-x_{13}a\,\hat{\mathbf{x}}+y_{13}b\,\hat{\mathbf{y}}+z_{13}c\,\hat{\mathbf{z}}$	(8*b*)	Si V

References:

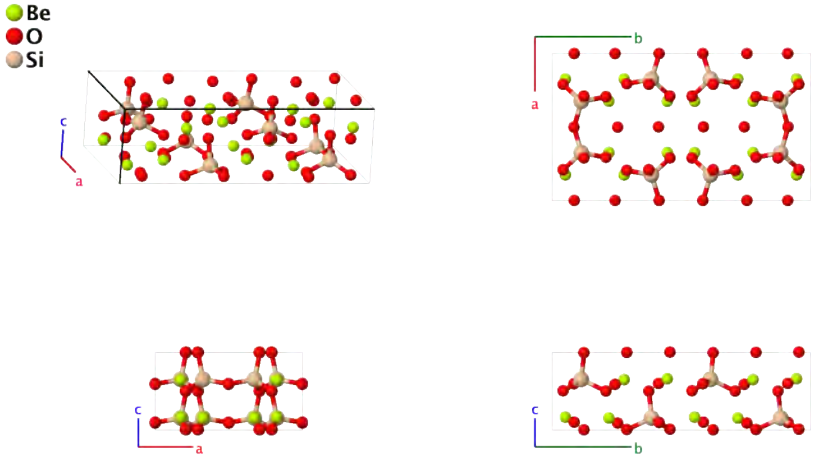
- G. Pilström, *The Crystal Structure of Ni_3Si_2 with some Notes on Ni_5Si_2*, Acta Chem. Scand. **15**, 893–902 (1961), doi:10.3891/acta.chem.scand.15-0893.

Geometry files:

- CIF: pp. 1594
- POSCAR: pp. 1594

Bertrandite ($Be_4Si_2O_7(OH)_2$, $S4_6$) Structure: A4B7C2D2_oC60_36_2b_a3b_2a_b

http://aflow.org/prototype-encyclopedia/A4B7C2D2_oC60_36_2b_a3b_2a_b

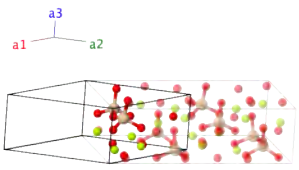


Prototype	:	$Be_4O_7(OH)_2Si_2$
AFLOW prototype label	:	A4B7C2D2_oC60_36_2b_a3b_2a_b
***Strukturbericht* designation**	:	$S4_6$
Pearson symbol	:	oC60
Space group number	:	36
Space group symbol	:	$Cmc2_1$
AFLOW prototype command	:	`aflow --proto=A4B7C2D2_oC60_36_2b_a3b_2a_b --params=`$a, b/a, c/a, y_1, z_1, y_2, z_2, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9$

- All sources agree that the space group for Bertrandite is $Cmc2_1$ #36 but differ slightly in details. We have chosen to use (Hazen, 1986) as the prototype over (Ito, 1932), which was originally designated $S4_6$, and (Solov'eva, 1965), which has unrealistically large Be-O distances.
- No paper locates the hydrogen atoms in the hydroxide ions, which may indicate that these ions can freely rotate.
- We use the data taken at 1 atmosphere.

Base-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$y_1 b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_2$	=	$y_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$-y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_3$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*a*)	OH I
$\mathbf{B}_4$	=	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$-y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	(4*a*)	OH I
$\mathbf{B}_5$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*a*)	OH II
$\mathbf{B}_6$	=	$y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$-y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	(4*a*)	OH II
$\mathbf{B}_7$	=	$(x_4 - y_4)\,\mathbf{a}_1 + (x_4 + y_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*b*)	Be I
$\mathbf{B}_8$	=	$(-x_4 + y_4)\,\mathbf{a}_1 + (-x_4 - y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	(8*b*)	Be I
$\mathbf{B}_9$	=	$(x_4 + y_4)\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	(8*b*)	Be I
$\mathbf{B}_{10}$	=	$(-x_4 - y_4)\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*b*)	Be I
$\mathbf{B}_{11}$	=	$(x_5 - y_5)\,\mathbf{a}_1 + (x_5 + y_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(8*b*)	Be II
$\mathbf{B}_{12}$	=	$(-x_5 + y_5)\,\mathbf{a}_1 + (-x_5 - y_5)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	(8*b*)	Be II
$\mathbf{B}_{13}$	=	$(x_5 + y_5)\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	(8*b*)	Be II
$\mathbf{B}_{14}$	=	$(-x_5 - y_5)\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(8*b*)	Be II
$\mathbf{B}_{15}$	=	$(x_6 - y_6)\,\mathbf{a}_1 + (x_6 + y_6)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(8*b*)	O II
$\mathbf{B}_{16}$	=	$(-x_6 + y_6)\,\mathbf{a}_1 + (-x_6 - y_6)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	(8*b*)	O II
$\mathbf{B}_{17}$	=	$(x_6 + y_6)\,\mathbf{a}_1 + (x_6 - y_6)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	(8*b*)	O II
$\mathbf{B}_{18}$	=	$(-x_6 - y_6)\,\mathbf{a}_1 + (-x_6 + y_6)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(8*b*)	O II
$\mathbf{B}_{19}$	=	$(x_7 - y_7)\,\mathbf{a}_1 + (x_7 + y_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(8*b*)	O III
$\mathbf{B}_{20}$	=	$(-x_7 + y_7)\,\mathbf{a}_1 + (-x_7 - y_7)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right) c\,\hat{\mathbf{z}}$	(8*b*)	O III
$\mathbf{B}_{21}$	=	$(x_7 + y_7)\,\mathbf{a}_1 + (x_7 - y_7)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right) c\,\hat{\mathbf{z}}$	(8*b*)	O III
$\mathbf{B}_{22}$	=	$(-x_7 - y_7)\,\mathbf{a}_1 + (-x_7 + y_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(8*b*)	O III
$\mathbf{B}_{23}$	=	$(x_8 - y_8)\,\mathbf{a}_1 + (x_8 + y_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(8*b*)	O IV
$\mathbf{B}_{24}$	=	$(-x_8 + y_8)\,\mathbf{a}_1 + (-x_8 - y_8)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right) c\,\hat{\mathbf{z}}$	(8*b*)	O IV
$\mathbf{B}_{25}$	=	$(x_8 + y_8)\,\mathbf{a}_1 + (x_8 - y_8)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right) c\,\hat{\mathbf{z}}$	(8*b*)	O IV
$\mathbf{B}_{26}$	=	$(-x_8 - y_8)\,\mathbf{a}_1 + (-x_8 + y_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(8*b*)	O IV
$\mathbf{B}_{27}$	=	$(x_9 - y_9)\,\mathbf{a}_1 + (x_9 + y_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(8*b*)	Si
$\mathbf{B}_{28}$	=	$(-x_9 + y_9)\,\mathbf{a}_1 + (-x_9 - y_9)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right) c\,\hat{\mathbf{z}}$	(8*b*)	Si
$\mathbf{B}_{29}$	=	$(x_9 + y_9)\,\mathbf{a}_1 + (x_9 - y_9)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right) c\,\hat{\mathbf{z}}$	(8*b*)	Si
$\mathbf{B}_{30}$	=	$(-x_9 - y_9)\,\mathbf{a}_1 + (-x_9 + y_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(8*b*)	Si

References:

- R. M. Hazen and A. Y. Au, *High-pressure crystal chemistry of phenakite (Be_2SiO_4) and bertrandite ($Be_4Si_2O_7(OH)_2$)*, Phys. Chem. Miner. **13**, 69–78 (1986), doi:10.1007/BF00311896.
- T. Ito and J. West, *The Structure of Bertrandite ($H_2Be_4Si_2O_9$)*, Zeitschrift für Kristallographie - Crystalline Materials **83**, 384–393 (1932), doi:10.1524/zkri.1932.83.1.384.
- L. P. Solov'eva and N. V. Belov, *Precise Determination of the Crystal Structure of Bertrandite $Be_4[Si_20_7](OH)_2$*, Sov. Phys. Crystallogr. **9**, 458–460 (1965).

Found in:

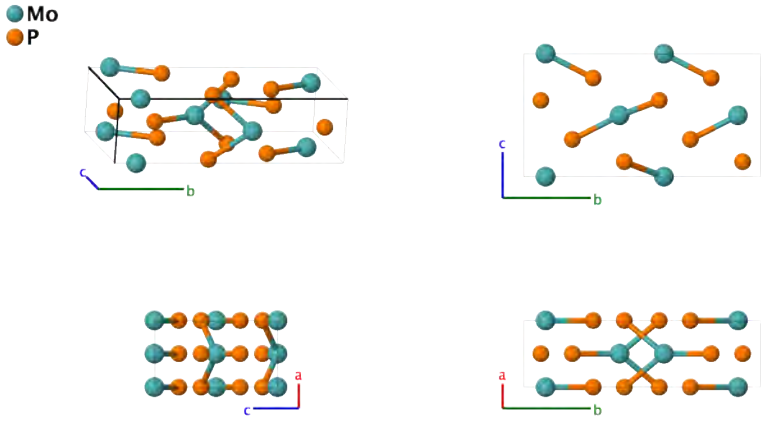
- B. Lafuente, R. T. Downs, H. Yang, and N. Stone, *The power of databases: the RRUFF project*, in *Highlights in Mineralogical Crystallography*, edited by T. Armbruster and R. M. Danisi (De Gruyter, Berlin, 2015), chap. 1, pp. 1–30.

Geometry files:

- CIF: pp. 1595
- POSCAR: pp. 1595

MoP_2 Structure: AB2_oC12_36_a_2a

http://aflow.org/prototype-encyclopedia/AB2_oC12_36_a_2a



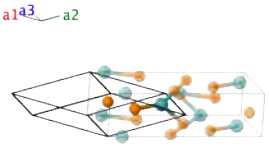
Prototype	:	MoP_2
AFLOW prototype label	:	AB2_oC12_36_a_2a
***Strukturbericht* designation**	:	None
Pearson symbol	:	oC12
Space group number	:	36
Space group symbol	:	$Cmc2_1$
AFLOW prototype command	:	`aflow --proto=AB2_oC12_36_a_2a --params=`$a, b/a, c/a, y_1, z_1, y_2, z_2, y_3, z_3$

Other compounds with this structure

- WP_2

Base-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$y_1 b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	$(4a)$	Mo
$\mathbf{B}_2$	=	$y_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$-y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	$(4a)$	Mo
$\mathbf{B}_3$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(4a)$	P I
$\mathbf{B}_4$	=	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$-y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	$(4a)$	P I
$\mathbf{B}_5$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(4a)$	P II
$\mathbf{B}_6$	=	$y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$-y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	$(4a)$	P II

References:

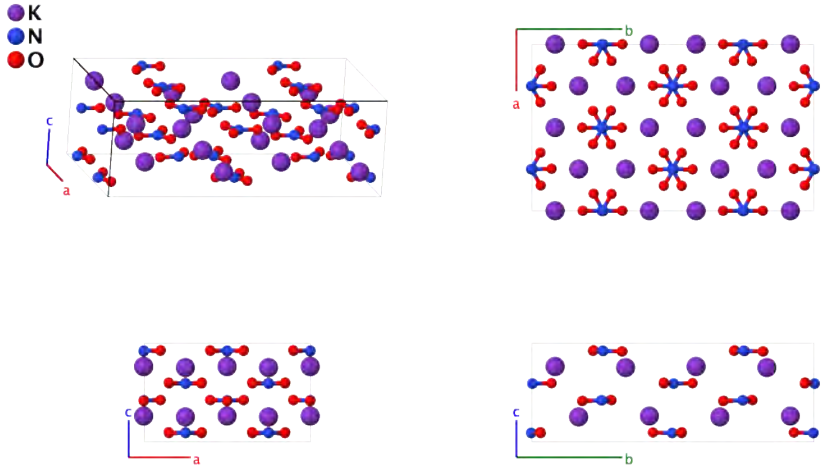
- S. Rundqvist and T. Lundström, *X-Ray Studies of Molybdenum and Tungsten Phosphides*, Acta Chem. Scand. **17**, 37–46 (1963), doi:10.3891/acta.chem.scand.17-0037.

Geometry files:

- CIF: pp. 1595
- POSCAR: pp. 1596

α-Potassium Nitrate (KNO_3) II Structure: ABC3_oC80_36_2ab_2ab_2a5b

http://aflow.org/prototype-encyclopedia/ABC3_oC80_36_2ab_2ab_2a5b

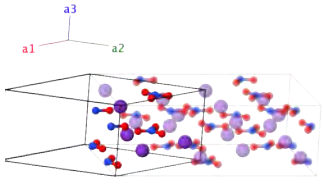


Prototype	:	KNO_3
AFLOW prototype label	:	ABC3_oC80_36_2ab_2ab_2a5b
***Strukturbericht* designation**	:	None
Pearson symbol	:	oC80
Space group number	:	36
Space group symbol	:	$Cmc2_1$
AFLOW prototype command	:	`aflow --proto=ABC3_oC80_36_2ab_2ab_2a5b --params=`$a, b/a, c/a, y_1, z_1, y_2, z_2, y_3, z_3, y_4, z_4, y_5, z_5, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}$

- Two possible structures have been identified for α-KNO_3: (Nimmo, 1973) proposed an orthorhombic structure in space group *Pnma* #62, which we call "Structure I". (Adiwidjaja, 2003) found that structure, but also noted that it could be described by a doubling of the unit cell into a superstructure of type I, which we call "Structure II" and present on this page. It is unclear to us which structure is correct, so we present both. However, we do note that if we allow some uncertainty in the atoms of this structure, AFLOW-SYM and FINDSYM show it to be identical to Structure I.
- On heating, α-KNO_3 transforms into β-KNO_3 at 128 °C. When heated above 200°C and then cooled, the β phase transforms into the metastable ferroelectric γ-KNO_3 phase, which remains metastable at room temperature.

Base-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$y_1 b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(4*a*)	K I
$\mathbf{B}_2$	=	$y_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$-y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	(4*a*)	K I
$\mathbf{B}_3$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*a*)	K II
$\mathbf{B}_4$	=	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$-y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	(4*a*)	K II
$\mathbf{B}_5$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*a*)	N I
$\mathbf{B}_6$	=	$y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$-y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	(4*a*)	N I
$\mathbf{B}_7$	=	$-y_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(4*a*)	N II
$\mathbf{B}_8$	=	$y_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$-y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	(4*a*)	N II
$\mathbf{B}_9$	=	$-y_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_{10}$	=	$y_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$-y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_{11}$	=	$-y_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(4*a*)	O II
$\mathbf{B}_{12}$	=	$y_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$-y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	(4*a*)	O II
$\mathbf{B}_{13}$	=	$(x_7 - y_7)\,\mathbf{a}_1 + (x_7 + y_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(8*b*)	K III
$\mathbf{B}_{14}$	=	$(-x_7 + y_7)\,\mathbf{a}_1 + (-x_7 - y_7)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right) c\,\hat{\mathbf{z}}$	(8*b*)	K III
$\mathbf{B}_{15}$	=	$(x_7 + y_7)\,\mathbf{a}_1 + (x_7 - y_7)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right) c\,\hat{\mathbf{z}}$	(8*b*)	K III
$\mathbf{B}_{16}$	=	$(-x_7 - y_7)\,\mathbf{a}_1 + (-x_7 + y_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(8*b*)	K III
$\mathbf{B}_{17}$	=	$(x_8 - y_8)\,\mathbf{a}_1 + (x_8 + y_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(8*b*)	N III
$\mathbf{B}_{18}$	=	$(-x_8 + y_8)\,\mathbf{a}_1 + (-x_8 - y_8)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right) c\,\hat{\mathbf{z}}$	(8*b*)	N III
$\mathbf{B}_{19}$	=	$(x_8 + y_8)\,\mathbf{a}_1 + (x_8 - y_8)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right) c\,\hat{\mathbf{z}}$	(8*b*)	N III
$\mathbf{B}_{20}$	=	$(-x_8 - y_8)\,\mathbf{a}_1 + (-x_8 + y_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(8*b*)	N III
$\mathbf{B}_{21}$	=	$(x_9 - y_9)\,\mathbf{a}_1 + (x_9 + y_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(8*b*)	O III
$\mathbf{B}_{22}$	=	$(-x_9 + y_9)\,\mathbf{a}_1 + (-x_9 - y_9)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right) c\,\hat{\mathbf{z}}$	(8*b*)	O III
$\mathbf{B}_{23}$	=	$(x_9 + y_9)\,\mathbf{a}_1 + (x_9 - y_9)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right) c\,\hat{\mathbf{z}}$	(8*b*)	O III
$\mathbf{B}_{24}$	=	$(-x_9 - y_9)\,\mathbf{a}_1 + (-x_9 + y_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(8*b*)	O III
$\mathbf{B}_{25}$	=	$(x_{10} - y_{10})\,\mathbf{a}_1 + (x_{10} + y_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$x_{10} a\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}}$	(8*b*)	O IV
$\mathbf{B}_{26}$	=	$(-x_{10} + y_{10})\,\mathbf{a}_1 + (-x_{10} - y_{10})\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{10}\right)\mathbf{a}_3$	=	$-x_{10} a\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{10}\right) c\,\hat{\mathbf{z}}$	(8*b*)	O IV
$\mathbf{B}_{27}$	=	$(x_{10} + y_{10})\,\mathbf{a}_1 + (x_{10} - y_{10})\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{10}\right)\mathbf{a}_3$	=	$x_{10} a\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{10}\right) c\,\hat{\mathbf{z}}$	(8*b*)	O IV
$\mathbf{B}_{28}$	=	$(-x_{10} - y_{10})\,\mathbf{a}_1 + (-x_{10} + y_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$-x_{10} a\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}}$	(8*b*)	O IV
$\mathbf{B}_{29}$	=	$(x_{11} - y_{11})\,\mathbf{a}_1 + (x_{11} + y_{11})\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$x_{11} a\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + z_{11} c\,\hat{\mathbf{z}}$	(8*b*)	O V

$$\mathbf{B}_{30} = (-x_{11}+y_{11})\,\mathbf{a}_1 + (-x_{11}-y_{11})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{11}\right)\mathbf{a}_3 = -x_{11}a\,\hat{\mathbf{x}} - y_{11}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{11}\right)c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O V}$$

$$\mathbf{B}_{31} = (x_{11}+y_{11})\,\mathbf{a}_1 + (x_{11}-y_{11})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{11}\right)\mathbf{a}_3 = x_{11}a\,\hat{\mathbf{x}} - y_{11}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{11}\right)c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O V}$$

$$\mathbf{B}_{32} = (-x_{11}-y_{11})\,\mathbf{a}_1 + (-x_{11}+y_{11})\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 = -x_{11}a\,\hat{\mathbf{x}} + y_{11}b\,\hat{\mathbf{y}} + z_{11}c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O V}$$

$$\mathbf{B}_{33} = (x_{12}-y_{12})\,\mathbf{a}_1 + (x_{12}+y_{12})\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 = x_{12}a\,\hat{\mathbf{x}} + y_{12}b\,\hat{\mathbf{y}} + z_{12}c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O VI}$$

$$\mathbf{B}_{34} = (-x_{12}+y_{12})\,\mathbf{a}_1 + (-x_{12}-y_{12})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{12}\right)\mathbf{a}_3 = -x_{12}a\,\hat{\mathbf{x}} - y_{12}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{12}\right)c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O VI}$$

$$\mathbf{B}_{35} = (x_{12}+y_{12})\,\mathbf{a}_1 + (x_{12}-y_{12})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{12}\right)\mathbf{a}_3 = x_{12}a\,\hat{\mathbf{x}} - y_{12}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{12}\right)c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O VI}$$

$$\mathbf{B}_{36} = (-x_{12}-y_{12})\,\mathbf{a}_1 + (-x_{12}+y_{12})\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 = -x_{12}a\,\hat{\mathbf{x}} + y_{12}b\,\hat{\mathbf{y}} + z_{12}c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O VI}$$

$$\mathbf{B}_{37} = (x_{13}-y_{13})\,\mathbf{a}_1 + (x_{13}+y_{13})\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3 = x_{13}a\,\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} + z_{13}c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O VII}$$

$$\mathbf{B}_{38} = (-x_{13}+y_{13})\,\mathbf{a}_1 + (-x_{13}-y_{13})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{13}\right)\mathbf{a}_3 = -x_{13}a\,\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{13}\right)c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O VII}$$

$$\mathbf{B}_{39} = (x_{13}+y_{13})\,\mathbf{a}_1 + (x_{13}-y_{13})\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{13}\right)\mathbf{a}_3 = x_{13}a\,\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{13}\right)c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O VII}$$

$$\mathbf{B}_{40} = (-x_{13}-y_{13})\,\mathbf{a}_1 + (-x_{13}+y_{13})\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3 = -x_{13}a\,\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} + z_{13}c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O VII}$$

References:

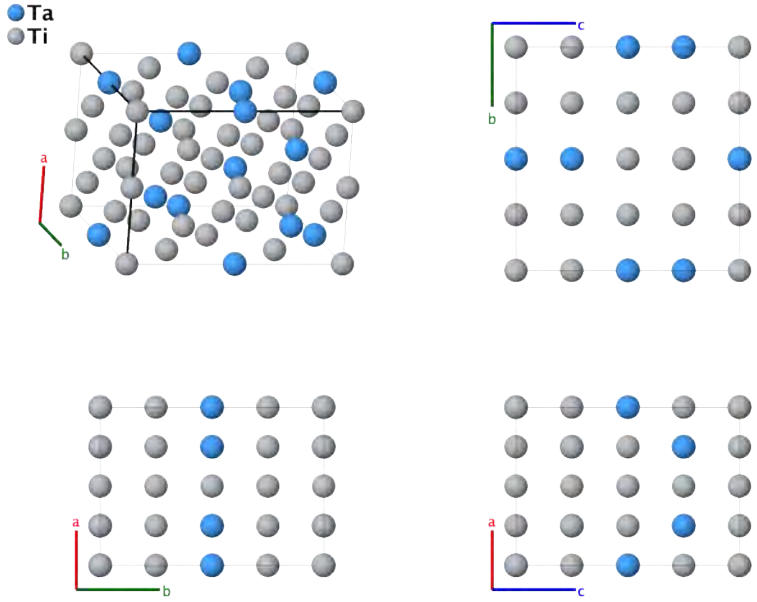
- G. Adiwidjaja and D. Pohl, *Superstructure of α-phase potassium nitrate*, Acta Crystallogr. C **59**, i139–i140 (2003), doi:10.1107/S0108270103025277.
- J. K. Nimmo and B. W. Lucas, *A neutron diffraction determination of the crystal structure of α-phase potassium nitrate at 25°C and 100°C*, J. Phys. C: Solid State Phys. **6**, 201–211 (1973), doi:10.1088/0022-3719/6/2/001.

Geometry files:

- CIF: pp. 1596
- POSCAR: pp. 1596

Ta_3Ti_{13} (BCC SQS-16) Structure: A3B13_oC32_38_ac_a2bcdef

http://aflow.org/prototype-encyclopedia/A3B13_oC32_38_ac_a2bcdef

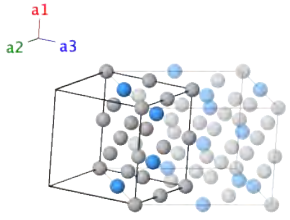


Prototype	:	Ta_3Ti_{13}
AFLOW prototype label	:	A3B13_oC32_38_ac_a2bcdef
***Strukturbericht* designation**	:	None
Pearson symbol	:	oC32
Space group number	:	38
Space group symbol	:	*Amm*2
AFLOW prototype command	:	`aflow --proto=A3B13_oC32_38_ac_a2bcdef --params=`$a, b/a, c/a, z_1, z_2, z_3, z_4, x_5, z_5, x_6, z_6, y_7, z_7, y_8, z_8, x_9, y_9, z_9$

- This is a special quasirandom structure with 16 atoms per unit cell (SQS-16) for the β-phase (high-temperature austenite) bcc substitutional Ti-Ta alloy (Chakraborty, 2016). This protoype contains 18.75% Ta. Prototypes are listed for other Ta-Ti concentrations: 12.5% Ta (AB7_hR16_166_c_c2h), 25% Ta (AB3_mC32_8_4a_4a4b), 31.25% Ta (A5B11_mP16_6_2abc_2a3b3c), and 37.5% Ta (A3B5_oC32_38_abce_abcdf).

Base-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}\,b\,\hat{\mathbf{y}} - \tfrac{1}{2}\,c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}\,b\,\hat{\mathbf{y}} + \tfrac{1}{2}\,c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-z_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$z_1 c\,\hat{\mathbf{z}}$	(2*a*)	Ta I
$\mathbf{B}_2$	=	$-z_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$z_2 c\,\hat{\mathbf{z}}$	(2*a*)	Ti I
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_1 - z_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + z_3 c\,\hat{\mathbf{z}}$	(2*b*)	Ti II
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 - z_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + z_4 c\,\hat{\mathbf{z}}$	(2*b*)	Ti III
$\mathbf{B}_5$	=	$x_5\,\mathbf{a}_1 - z_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + z_5 c\,\hat{\mathbf{z}}$	(4*c*)	Ta II
$\mathbf{B}_6$	=	$-x_5\,\mathbf{a}_1 - z_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + z_5 c\,\hat{\mathbf{z}}$	(4*c*)	Ta II
$\mathbf{B}_7$	=	$x_6\,\mathbf{a}_1 - z_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + z_6 c\,\hat{\mathbf{z}}$	(4*c*)	Ti IV
$\mathbf{B}_8$	=	$-x_6\,\mathbf{a}_1 - z_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + z_6 c\,\hat{\mathbf{z}}$	(4*c*)	Ti IV
$\mathbf{B}_9$	=	$(y_7 - z_7)\,\mathbf{a}_2 + (y_7 + z_7)\,\mathbf{a}_3$	=	$y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(4*d*)	Ti V
$\mathbf{B}_{10}$	=	$(-y_7 - z_7)\,\mathbf{a}_2 + (-y_7 + z_7)\,\mathbf{a}_3$	=	$-y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(4*d*)	Ti V
$\mathbf{B}_{11}$	=	$\frac{1}{2}\,\mathbf{a}_1 + (y_8 - z_8)\,\mathbf{a}_2 + (y_8 + z_8)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(4*e*)	Ti VI
$\mathbf{B}_{12}$	=	$\frac{1}{2}\,\mathbf{a}_1 + (-y_8 - z_8)\,\mathbf{a}_2 + (-y_8 + z_8)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(4*e*)	Ti VI
$\mathbf{B}_{13}$	=	$x_9\,\mathbf{a}_1 + (y_9 - z_9)\,\mathbf{a}_2 + (y_9 + z_9)\,\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(8*f*)	Ti VII
$\mathbf{B}_{14}$	=	$-x_9\,\mathbf{a}_1 + (-y_9 - z_9)\,\mathbf{a}_2 + (-y_9 + z_9)\,\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(8*f*)	Ti VII
$\mathbf{B}_{15}$	=	$x_9\,\mathbf{a}_1 + (-y_9 - z_9)\,\mathbf{a}_2 + (-y_9 + z_9)\,\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(8*f*)	Ti VII
$\mathbf{B}_{16}$	=	$-x_9\,\mathbf{a}_1 + (y_9 - z_9)\,\mathbf{a}_2 + (y_9 + z_9)\,\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(8*f*)	Ti VII

References:

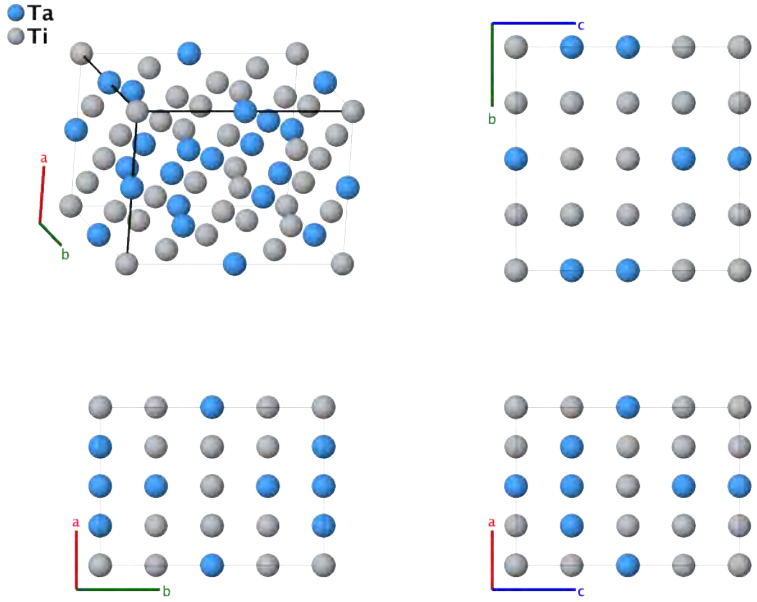
- T. Chakraborty, J. Rogal, and R. Drautz, *Unraveling the composition dependence of the martensitic transformation temperature: A first-principles study of Ti-Ta alloys*, Phys. Rev. B **94**, 224104 (2016), doi:10.1103/PhysRevB.94.224104.

Geometry files:

- CIF: pp. 1596
- POSCAR: pp. 1597

Ta_3Ti_5 (BCC SQS-16) Structure: A3B5_oC32_38_abce_abcdf

http://aflow.org/prototype-encyclopedia/A3B5_oC32_38_abce_abcdf

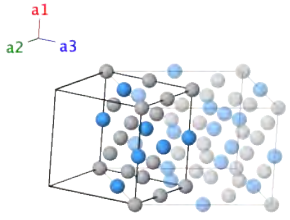


Prototype	:	Ta_3Ti_5
AFLOW prototype label	:	A3B5_oC32_38_abce_abcdf
***Strukturbericht* designation**	:	None
Pearson symbol	:	oC32
Space group number	:	38
Space group symbol	:	*Amm*2
AFLOW prototype command	:	`aflow --proto=A3B5_oC32_38_abce_abcdf --params=`$a, b/a, c/a, z_1, z_2, z_3, z_4, x_5, z_5, x_6, z_6, y_7, z_7, y_8, z_8, x_9, y_9, z_9$

- This is a special quasirandom structure with 16 atoms per unit cell (SQS-16) for the β-phase (high-temperature austenite) bcc substitutional Ti-Ta alloy (Chakraborty, 2016). This protoype contains 37.5% Ta. Prototypes are listed for other Ta-Ti concentrations: 12.5% Ta (AB7_hR16_166_c_c2h), 18.75% Ta (A3B13_oC32_38_ac_a2bcdef), 25% Ta (AB3_mC32_8_4a_4a4b), 31.25% Ta (A5B11_mP16_6_2abc_2a3b3c).

Base-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}\,b\,\hat{\mathbf{y}} - \tfrac{1}{2}\,c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}\,b\,\hat{\mathbf{y}} + \tfrac{1}{2}\,c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-z_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$z_1 c\,\hat{\mathbf{z}}$	($2a$)	Ta I
$\mathbf{B}_2$	=	$-z_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$z_2 c\,\hat{\mathbf{z}}$	($2a$)	Ti I
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_1 - z_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + z_3 c\,\hat{\mathbf{z}}$	($2b$)	Ta II
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 - z_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + z_4 c\,\hat{\mathbf{z}}$	($2b$)	Ti II
$\mathbf{B}_5$	=	$x_5\,\mathbf{a}_1 - z_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + z_5 c\,\hat{\mathbf{z}}$	($4c$)	Ta III
$\mathbf{B}_6$	=	$-x_5\,\mathbf{a}_1 - z_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + z_5 c\,\hat{\mathbf{z}}$	($4c$)	Ta III
$\mathbf{B}_7$	=	$x_6\,\mathbf{a}_1 - z_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + z_6 c\,\hat{\mathbf{z}}$	($4c$)	Ti III
$\mathbf{B}_8$	=	$-x_6\,\mathbf{a}_1 - z_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + z_6 c\,\hat{\mathbf{z}}$	($4c$)	Ti III
$\mathbf{B}_9$	=	$(y_7 - z_7)\,\mathbf{a}_2 + (y_7 + z_7)\,\mathbf{a}_3$	=	$y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	($4d$)	Ti IV
$\mathbf{B}_{10}$	=	$(-y_7 - z_7)\,\mathbf{a}_2 + (-y_7 + z_7)\,\mathbf{a}_3$	=	$-y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	($4d$)	Ti IV
$\mathbf{B}_{11}$	=	$\frac{1}{2}\,\mathbf{a}_1 + (y_8 - z_8)\,\mathbf{a}_2 + (y_8 + z_8)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	($4e$)	Ta IV
$\mathbf{B}_{12}$	=	$\frac{1}{2}\,\mathbf{a}_1 + (-y_8 - z_8)\,\mathbf{a}_2 + (-y_8 + z_8)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	($4e$)	Ta IV
$\mathbf{B}_{13}$	=	$x_9\,\mathbf{a}_1 + (y_9 - z_9)\,\mathbf{a}_2 + (y_9 + z_9)\,\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	($8f$)	Ti V
$\mathbf{B}_{14}$	=	$-x_9\,\mathbf{a}_1 + (-y_9 - z_9)\,\mathbf{a}_2 + (-y_9 + z_9)\,\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	($8f$)	Ti V
$\mathbf{B}_{15}$	=	$x_9\,\mathbf{a}_1 + (-y_9 - z_9)\,\mathbf{a}_2 + (-y_9 + z_9)\,\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	($8f$)	Ti V
$\mathbf{B}_{16}$	=	$-x_9\,\mathbf{a}_1 + (y_9 - z_9)\,\mathbf{a}_2 + (y_9 + z_9)\,\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	($8f$)	Ti V

References:

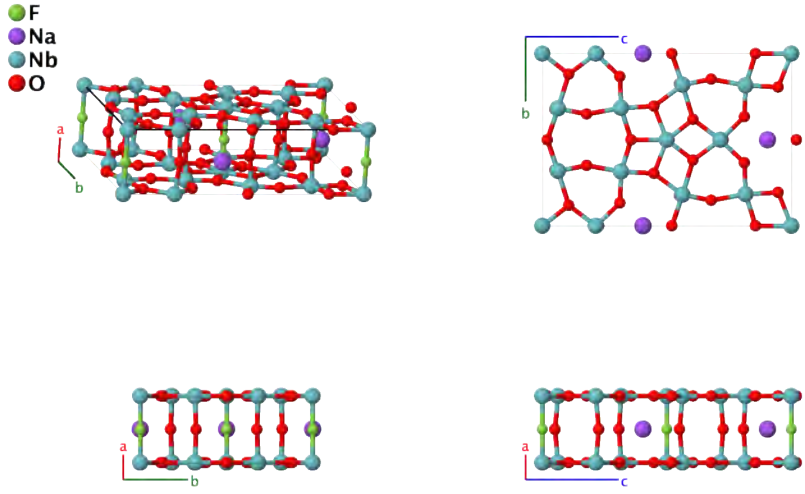
- T. Chakraborty, J. Rogal, and R. Drautz, *Unraveling the composition dependence of the martensitic transformation temperature: A first-principles study of Ti-Ta alloys*, Phys. Rev. B **94**, 224104 (2016), doi:10.1103/PhysRevB.94.224104.

Geometry files:

- CIF: pp. 1597
- POSCAR: pp. 1597

$NaNb_6O_{15}F$ Structure: ABC6D15_oC46_38_b_b_2a2d_2ab4d2e

http://aflow.org/prototype-encyclopedia/ABC6D15_oC46_38_b_b_2a2d_2ab4d2e



Prototype	:	$FNaNb_6O_{15}$
AFLOW prototype label	:	ABC6D15_oC46_38_b_b_2a2d_2ab4d2e
***Strukturbericht* designation**	:	None
Pearson symbol	:	oC46
Space group number	:	38
Space group symbol	:	*Amm*2
AFLOW prototype command	:	`aflow --proto=ABC6D15_oC46_38_b_b_2a2d_2ab4d2e --params=`$a, b/a, c/a, z_1, z_2, z_3, z_4, z_5, z_6, z_7, y_8, z_8, y_9, z_9, y_{10}, z_{10}, y_{11}, z_{11}, y_{12}, z_{12}, y_{13}, z_{13}, y_{14}, z_{14}, y_{15}, z_{15}$

Other compounds with this structure

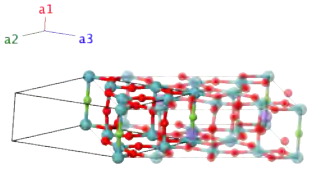
- $NaNb_6O_{15}(OH)$

- The x-ray scattering of an F^- ion is almost identical to that of O^{2-}, and (Andersson, 1965) was not able to distinguish between them. He arbitrarily labeled the (2*b*) site he designated as O(1) as the location of the fluorine ion and we follow this, but in reality we have no idea if the F^- ions are located on this site, are ordered on another site, or are statistically distributed on the oxygen sites. Presumably the same considerations hold for $NaNb_6O_{15}(OH)$.

- Andersson sets $z_4 = 0.159$ as the coordinate of what we label as O-II and he calls O(10), but this gives an unreasonably short distance between the Nb-II and O-II atoms, and the distances between the O-II atom and the other atoms in the structure do not agree with the distances given his paper. If we assume that the first two digits were transposed when printed, so that $z_4 = 0.519$, then we get within 0.1% of Andersson's distances.

Base-centered Orthorhombic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= a\,\hat{\mathbf{x}} \\ \mathbf{a}_2 &= \tfrac{1}{2}b\,\hat{\mathbf{y}} - \tfrac{1}{2}c\,\hat{\mathbf{z}} \\ \mathbf{a}_3 &= \tfrac{1}{2}b\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-z_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$z_1 c\,\hat{\mathbf{z}}$	$(2a)$	Nb I
$\mathbf{B}_2$	=	$-z_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$z_2 c\,\hat{\mathbf{z}}$	$(2a)$	Nb II
$\mathbf{B}_3$	=	$-z_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$z_3 c\,\hat{\mathbf{z}}$	$(2a)$	O I
$\mathbf{B}_4$	=	$-z_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$z_4 c\,\hat{\mathbf{z}}$	$(2a)$	O II
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_1 - z_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + z_5 c\,\hat{\mathbf{z}}$	$(2b)$	F
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_1 - z_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + z_6 c\,\hat{\mathbf{z}}$	$(2b)$	Na
$\mathbf{B}_7$	=	$\frac{1}{2}\,\mathbf{a}_1 - z_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + z_7 c\,\hat{\mathbf{z}}$	$(2b)$	O III
$\mathbf{B}_8$	=	$(y_8 - z_8)\,\mathbf{a}_2 + (y_8 + z_8)\,\mathbf{a}_3$	=	$y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	$(4d)$	Nb III
$\mathbf{B}_9$	=	$(-y_8 - z_8)\,\mathbf{a}_2 + (-y_8 + z_8)\,\mathbf{a}_3$	=	$-y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	$(4d)$	Nb III
$\mathbf{B}_{10}$	=	$(y_9 - z_9)\,\mathbf{a}_2 + (y_9 + z_9)\,\mathbf{a}_3$	=	$y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	$(4d)$	Nb IV
$\mathbf{B}_{11}$	=	$(-y_9 - z_9)\,\mathbf{a}_2 + (-y_9 + z_9)\,\mathbf{a}_3$	=	$-y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	$(4d)$	Nb IV
$\mathbf{B}_{12}$	=	$(y_{10} - z_{10})\,\mathbf{a}_2 + (y_{10} + z_{10})\,\mathbf{a}_3$	=	$y_{10} b\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}}$	$(4d)$	O IV
$\mathbf{B}_{13}$	=	$(-y_{10} - z_{10})\,\mathbf{a}_2 + (-y_{10} + z_{10})\,\mathbf{a}_3$	=	$-y_{10} b\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}}$	$(4d)$	O IV
$\mathbf{B}_{14}$	=	$(y_{11} - z_{11})\,\mathbf{a}_2 + (y_{11} + z_{11})\,\mathbf{a}_3$	=	$y_{11} b\,\hat{\mathbf{y}} + z_{11} c\,\hat{\mathbf{z}}$	$(4d)$	O V
$\mathbf{B}_{15}$	=	$(-y_{11} - z_{11})\,\mathbf{a}_2 + (-y_{11} + z_{11})\,\mathbf{a}_3$	=	$-y_{11} b\,\hat{\mathbf{y}} + z_{11} c\,\hat{\mathbf{z}}$	$(4d)$	O V
$\mathbf{B}_{16}$	=	$(y_{12} - z_{12})\,\mathbf{a}_2 + (y_{12} + z_{12})\,\mathbf{a}_3$	=	$y_{12} b\,\hat{\mathbf{y}} + z_{12} c\,\hat{\mathbf{z}}$	$(4d)$	O VI
$\mathbf{B}_{17}$	=	$(-y_{12} - z_{12})\,\mathbf{a}_2 + (-y_{12} + z_{12})\,\mathbf{a}_3$	=	$-y_{12} b\,\hat{\mathbf{y}} + z_{12} c\,\hat{\mathbf{z}}$	$(4d)$	O VI
$\mathbf{B}_{18}$	=	$(y_{13} - z_{13})\,\mathbf{a}_2 + (y_{13} + z_{13})\,\mathbf{a}_3$	=	$y_{13} b\,\hat{\mathbf{y}} + z_{13} c\,\hat{\mathbf{z}}$	$(4d)$	O VII
$\mathbf{B}_{19}$	=	$(-y_{13} - z_{13})\,\mathbf{a}_2 + (-y_{13} + z_{13})\,\mathbf{a}_3$	=	$-y_{13} b\,\hat{\mathbf{y}} + z_{13} c\,\hat{\mathbf{z}}$	$(4d)$	O VII
$\mathbf{B}_{20}$	=	$\frac{1}{2}\,\mathbf{a}_1 + (y_{14} - z_{14})\,\mathbf{a}_2 + (y_{14} + z_{14})\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + y_{14} b\,\hat{\mathbf{y}} + z_{14} c\,\hat{\mathbf{z}}$	$(4e)$	O VIII
$\mathbf{B}_{21}$	=	$\frac{1}{2}\,\mathbf{a}_1 + (-y_{14} - z_{14})\,\mathbf{a}_2 + (-y_{14} + z_{14})\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - y_{14} b\,\hat{\mathbf{y}} + z_{14} c\,\hat{\mathbf{z}}$	$(4e)$	O VIII
$\mathbf{B}_{22}$	=	$\frac{1}{2}\,\mathbf{a}_1 + (y_{15} - z_{15})\,\mathbf{a}_2 + (y_{15} + z_{15})\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + y_{15} b\,\hat{\mathbf{y}} + z_{15} c\,\hat{\mathbf{z}}$	$(4e)$	O IX
$\mathbf{B}_{23}$	=	$\frac{1}{2}\,\mathbf{a}_1 + (-y_{15} - z_{15})\,\mathbf{a}_2 + (-y_{15} + z_{15})\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - y_{15} b\,\hat{\mathbf{y}} + z_{15} c\,\hat{\mathbf{z}}$	$(4e)$	O IX

References:

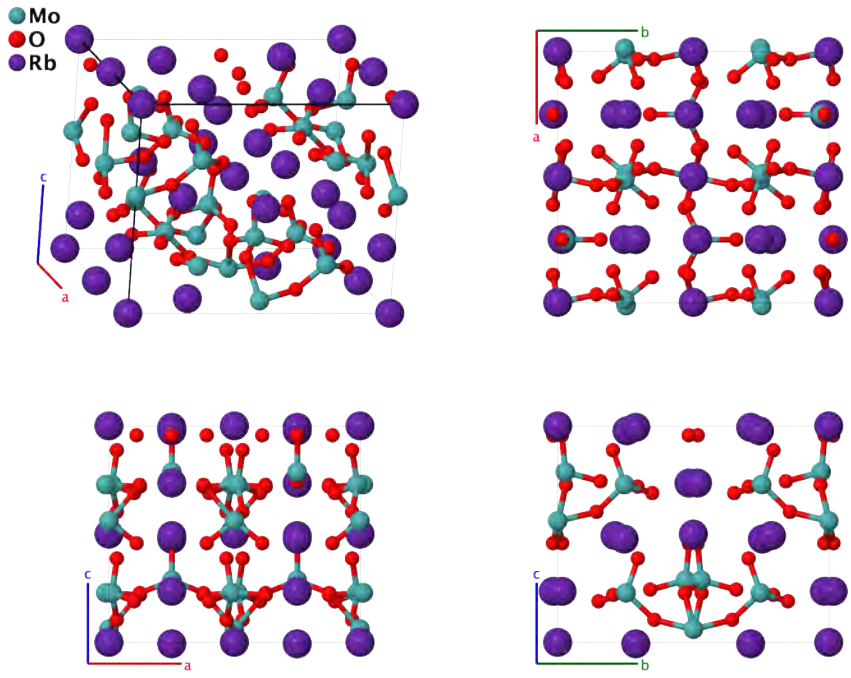
- S. Andersson, *The Crystal Structure of $NaNb_6O_{15}F$ and $NaNb_6O_{15}OH$*, Acta Chem. Scand. **19**, 2285–2290 (1965), doi:10.3891/acta.chem.scand.19-2285.

Geometry files:

- CIF: pp. 1597
- POSCAR: pp. 1598

$Rb_2Mo_2O_7$ Structure: A2B7C2_oC88_40_abc_2b6c_a3b

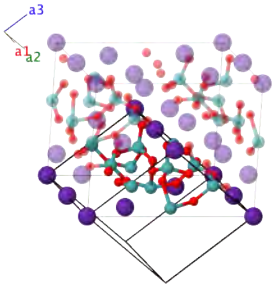
http://aflow.org/prototype-encyclopedia/A2B7C2_oC88_40_abc_2b6c_a3b



Prototype	:	$Mo_2O_7Rb_2$
AFLOW prototype label	:	A2B7C2_oC88_40_abc_2b6c_a3b
***Strukturbericht* designation**	:	None
Pearson symbol	:	oC88
Space group number	:	40
Space group symbol	:	*Ama*2
AFLOW prototype command	:	`aflow --proto=A2B7C2_oC88_40_abc_2b6c_a3b --params=`$a, b/a, c/a, z_1, z_2, y_3, z_3, y_4, z_4, y_5, z_5, y_6, z_6, y_7, z_7, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}$

Base-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}\,b\,\hat{\mathbf{y}} - \tfrac{1}{2}\,c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}\,b\,\hat{\mathbf{y}} + \tfrac{1}{2}\,c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-z_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$z_1 c\,\hat{\mathbf{z}}$	($4a$)	Mo I

$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 - z_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + z_1 c\,\hat{\mathbf{z}}$	(4*a*)	Mo I
$\mathbf{B}_3$	=	$-z_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$z_2 c\,\hat{\mathbf{z}}$	(4*a*)	Rb I
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 - z_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + z_2 c\,\hat{\mathbf{z}}$	(4*a*)	Rb I
$\mathbf{B}_5$	=	$\frac{1}{4}\,\mathbf{a}_1 + (y_3 - z_3)\,\mathbf{a}_2 + (y_3 + z_3)\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*b*)	Mo II
$\mathbf{B}_6$	=	$\frac{3}{4}\,\mathbf{a}_1 + (-y_3 - z_3)\,\mathbf{a}_2 + (-y_3 + z_3)\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*b*)	Mo II
$\mathbf{B}_7$	=	$\frac{1}{4}\,\mathbf{a}_1 + (y_4 - z_4)\,\mathbf{a}_2 + (y_4 + z_4)\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(4*b*)	O I
$\mathbf{B}_8$	=	$\frac{3}{4}\,\mathbf{a}_1 + (-y_4 - z_4)\,\mathbf{a}_2 + (-y_4 + z_4)\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(4*b*)	O I
$\mathbf{B}_9$	=	$\frac{1}{4}\,\mathbf{a}_1 + (y_5 - z_5)\,\mathbf{a}_2 + (y_5 + z_5)\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(4*b*)	O II
$\mathbf{B}_{10}$	=	$\frac{3}{4}\,\mathbf{a}_1 + (-y_5 - z_5)\,\mathbf{a}_2 + (-y_5 + z_5)\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(4*b*)	O II
$\mathbf{B}_{11}$	=	$\frac{1}{4}\,\mathbf{a}_1 + (y_6 - z_6)\,\mathbf{a}_2 + (y_6 + z_6)\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(4*b*)	Rb II
$\mathbf{B}_{12}$	=	$\frac{3}{4}\,\mathbf{a}_1 + (-y_6 - z_6)\,\mathbf{a}_2 + (-y_6 + z_6)\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(4*b*)	Rb II
$\mathbf{B}_{13}$	=	$\frac{1}{4}\,\mathbf{a}_1 + (y_7 - z_7)\,\mathbf{a}_2 + (y_7 + z_7)\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(4*b*)	Rb III
$\mathbf{B}_{14}$	=	$\frac{3}{4}\,\mathbf{a}_1 + (-y_7 - z_7)\,\mathbf{a}_2 + (-y_7 + z_7)\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(4*b*)	Rb III
$\mathbf{B}_{15}$	=	$\frac{1}{4}\,\mathbf{a}_1 + (y_8 - z_8)\,\mathbf{a}_2 + (y_8 + z_8)\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(4*b*)	Rb IV
$\mathbf{B}_{16}$	=	$\frac{3}{4}\,\mathbf{a}_1 + (-y_8 - z_8)\,\mathbf{a}_2 + (-y_8 + z_8)\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(4*b*)	Rb IV
$\mathbf{B}_{17}$	=	$x_9\,\mathbf{a}_1 + (y_9 - z_9)\,\mathbf{a}_2 + (y_9 + z_9)\,\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(8*c*)	Mo III
$\mathbf{B}_{18}$	=	$-x_9\,\mathbf{a}_1 + (-y_9 - z_9)\,\mathbf{a}_2 + (-y_9 + z_9)\,\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(8*c*)	Mo III
$\mathbf{B}_{19}$	=	$\left(\frac{1}{2} + x_9\right)\mathbf{a}_1 + (-y_9 - z_9)\,\mathbf{a}_2 + (-y_9 + z_9)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_9\right)a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(8*c*)	Mo III
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} - x_9\right)\mathbf{a}_1 + (y_9 - z_9)\,\mathbf{a}_2 + (y_9 + z_9)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_9\right)a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(8*c*)	Mo III
$\mathbf{B}_{21}$	=	$x_{10}\,\mathbf{a}_1 + (y_{10} - z_{10})\,\mathbf{a}_2 + (y_{10} + z_{10})\,\mathbf{a}_3$	=	$x_{10}a\,\hat{\mathbf{x}} + y_{10}b\,\hat{\mathbf{y}} + z_{10}c\,\hat{\mathbf{z}}$	(8*c*)	O III
$\mathbf{B}_{22}$	=	$-x_{10}\,\mathbf{a}_1 + (-y_{10} - z_{10})\,\mathbf{a}_2 + (-y_{10} + z_{10})\,\mathbf{a}_3$	=	$-x_{10}a\,\hat{\mathbf{x}} - y_{10}b\,\hat{\mathbf{y}} + z_{10}c\,\hat{\mathbf{z}}$	(8*c*)	O III
$\mathbf{B}_{23}$	=	$\left(\frac{1}{2} + x_{10}\right)\mathbf{a}_1 + (-y_{10} - z_{10})\,\mathbf{a}_2 + (-y_{10} + z_{10})\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{10}\right)a\,\hat{\mathbf{x}} - y_{10}b\,\hat{\mathbf{y}} + z_{10}c\,\hat{\mathbf{z}}$	(8*c*)	O III
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} - x_{10}\right)\mathbf{a}_1 + (y_{10} - z_{10})\,\mathbf{a}_2 + (y_{10} + z_{10})\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{10}\right)a\,\hat{\mathbf{x}} + y_{10}b\,\hat{\mathbf{y}} + z_{10}c\,\hat{\mathbf{z}}$	(8*c*)	O III
$\mathbf{B}_{25}$	=	$x_{11}\,\mathbf{a}_1 + (y_{11} - z_{11})\,\mathbf{a}_2 + (y_{11} + z_{11})\,\mathbf{a}_3$	=	$x_{11}a\,\hat{\mathbf{x}} + y_{11}b\,\hat{\mathbf{y}} + z_{11}c\,\hat{\mathbf{z}}$	(8*c*)	O IV
$\mathbf{B}_{26}$	=	$-x_{11}\,\mathbf{a}_1 + (-y_{11} - z_{11})\,\mathbf{a}_2 + (-y_{11} + z_{11})\,\mathbf{a}_3$	=	$-x_{11}a\,\hat{\mathbf{x}} - y_{11}b\,\hat{\mathbf{y}} + z_{11}c\,\hat{\mathbf{z}}$	(8*c*)	O IV
$\mathbf{B}_{27}$	=	$\left(\frac{1}{2} + x_{11}\right)\mathbf{a}_1 + (-y_{11} - z_{11})\,\mathbf{a}_2 + (-y_{11} + z_{11})\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{11}\right)a\,\hat{\mathbf{x}} - y_{11}b\,\hat{\mathbf{y}} + z_{11}c\,\hat{\mathbf{z}}$	(8*c*)	O IV
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} - x_{11}\right)\mathbf{a}_1 + (y_{11} - z_{11})\,\mathbf{a}_2 + (y_{11} + z_{11})\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{11}\right)a\,\hat{\mathbf{x}} + y_{11}b\,\hat{\mathbf{y}} + z_{11}c\,\hat{\mathbf{z}}$	(8*c*)	O IV
$\mathbf{B}_{29}$	=	$x_{12}\,\mathbf{a}_1 + (y_{12} - z_{12})\,\mathbf{a}_2 + (y_{12} + z_{12})\,\mathbf{a}_3$	=	$x_{12}a\,\hat{\mathbf{x}} + y_{12}b\,\hat{\mathbf{y}} + z_{12}c\,\hat{\mathbf{z}}$	(8*c*)	O V
$\mathbf{B}_{30}$	=	$-x_{12}\,\mathbf{a}_1 + (-y_{12} - z_{12})\,\mathbf{a}_2 + (-y_{12} + z_{12})\,\mathbf{a}_3$	=	$-x_{12}a\,\hat{\mathbf{x}} - y_{12}b\,\hat{\mathbf{y}} + z_{12}c\,\hat{\mathbf{z}}$	(8*c*)	O V
$\mathbf{B}_{31}$	=	$\left(\frac{1}{2} + x_{12}\right)\mathbf{a}_1 + (-y_{12} - z_{12})\,\mathbf{a}_2 + (-y_{12} + z_{12})\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{12}\right)a\,\hat{\mathbf{x}} - y_{12}b\,\hat{\mathbf{y}} + z_{12}c\,\hat{\mathbf{z}}$	(8*c*)	O V
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2} - x_{12}\right)\mathbf{a}_1 + (y_{12} - z_{12})\,\mathbf{a}_2 + (y_{12} + z_{12})\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{12}\right)a\,\hat{\mathbf{x}} + y_{12}b\,\hat{\mathbf{y}} + z_{12}c\,\hat{\mathbf{z}}$	(8*c*)	O V
$\mathbf{B}_{33}$	=	$x_{13}\,\mathbf{a}_1 + (y_{13} - z_{13})\,\mathbf{a}_2 + (y_{13} + z_{13})\,\mathbf{a}_3$	=	$x_{13}a\,\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} + z_{13}c\,\hat{\mathbf{z}}$	(8*c*)	O VI

$\mathbf{B}_{34}$	=	$-x_{13}\,\mathbf{a}_1 + (-y_{13} - z_{13})\,\mathbf{a}_2 + (-y_{13} + z_{13})\,\mathbf{a}_3$	=	$-x_{13}a\,\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} + z_{13}c\,\hat{\mathbf{z}}$	(8*c*)	O VI
$\mathbf{B}_{35}$	=	$\left(\frac{1}{2} + x_{13}\right)\mathbf{a}_1 + (-y_{13} - z_{13})\,\mathbf{a}_2 + (-y_{13} + z_{13})\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{13}\right)a\,\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} + z_{13}c\,\hat{\mathbf{z}}$	(8*c*)	O VI
$\mathbf{B}_{36}$	=	$\left(\frac{1}{2} - x_{13}\right)\mathbf{a}_1 + (y_{13} - z_{13})\,\mathbf{a}_2 + (y_{13} + z_{13})\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{13}\right)a\,\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} + z_{13}c\,\hat{\mathbf{z}}$	(8*c*)	O VI
$\mathbf{B}_{37}$	=	$x_{14}\,\mathbf{a}_1 + (y_{14} - z_{14})\,\mathbf{a}_2 + (y_{14} + z_{14})\,\mathbf{a}_3$	=	$x_{14}a\,\hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} + z_{14}c\,\hat{\mathbf{z}}$	(8*c*)	O VII
$\mathbf{B}_{38}$	=	$-x_{14}\,\mathbf{a}_1 + (-y_{14} - z_{14})\,\mathbf{a}_2 + (-y_{14} + z_{14})\,\mathbf{a}_3$	=	$-x_{14}a\,\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} + z_{14}c\,\hat{\mathbf{z}}$	(8*c*)	O VII
$\mathbf{B}_{39}$	=	$\left(\frac{1}{2} + x_{14}\right)\mathbf{a}_1 + (-y_{14} - z_{14})\,\mathbf{a}_2 + (-y_{14} + z_{14})\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{14}\right)a\,\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} + z_{14}c\,\hat{\mathbf{z}}$	(8*c*)	O VII
$\mathbf{B}_{40}$	=	$\left(\frac{1}{2} - x_{14}\right)\mathbf{a}_1 + (y_{14} - z_{14})\,\mathbf{a}_2 + (y_{14} + z_{14})\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{14}\right)a\,\hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} + z_{14}c\,\hat{\mathbf{z}}$	(8*c*)	O VII
$\mathbf{B}_{41}$	=	$x_{15}\,\mathbf{a}_1 + (y_{15} - z_{15})\,\mathbf{a}_2 + (y_{15} + z_{15})\,\mathbf{a}_3$	=	$x_{15}a\,\hat{\mathbf{x}} + y_{15}b\,\hat{\mathbf{y}} + z_{15}c\,\hat{\mathbf{z}}$	(8*c*)	O VIII
$\mathbf{B}_{42}$	=	$-x_{15}\,\mathbf{a}_1 + (-y_{15} - z_{15})\,\mathbf{a}_2 + (-y_{15} + z_{15})\,\mathbf{a}_3$	=	$-x_{15}a\,\hat{\mathbf{x}} - y_{15}b\,\hat{\mathbf{y}} + z_{15}c\,\hat{\mathbf{z}}$	(8*c*)	O VIII
$\mathbf{B}_{43}$	=	$\left(\frac{1}{2} + x_{15}\right)\mathbf{a}_1 + (-y_{15} - z_{15})\,\mathbf{a}_2 + (-y_{15} + z_{15})\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{15}\right)a\,\hat{\mathbf{x}} - y_{15}b\,\hat{\mathbf{y}} + z_{15}c\,\hat{\mathbf{z}}$	(8*c*)	O VIII
$\mathbf{B}_{44}$	=	$\left(\frac{1}{2} - x_{15}\right)\mathbf{a}_1 + (y_{15} - z_{15})\,\mathbf{a}_2 + (y_{15} + z_{15})\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{15}\right)a\,\hat{\mathbf{x}} + y_{15}b\,\hat{\mathbf{y}} + z_{15}c\,\hat{\mathbf{z}}$	(8*c*)	O VIII

References:

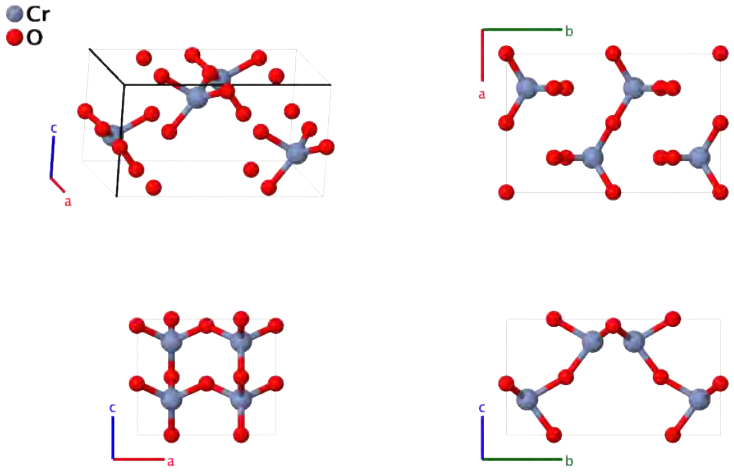
- Z. A. Solodovnikova and S. F. Solodovnikov, *Rubidium dimolybdate, $Rb_2Mo_2O_7$, and caesium dimolybdate, $Cs_2Mo_2O_7$*, Acta Crystallogr. C **62**, i53–i56 (2006), doi:10.1107/S0108270106014880.

Geometry files:

- CIF: pp. 1598
- POSCAR: pp. 1598

Orthorhombic CrO_3 Structure: AB3_oC16_40_b_a2b

http://aflow.org/prototype-encyclopedia/AB3_oC16_40_b_a2b

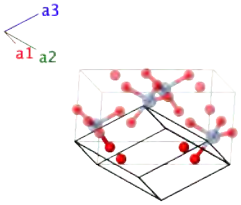


Prototype	:	CrO_3
AFLOW prototype label	:	AB3_oC16_40_b_a2b
***Strukturbericht* designation**	:	None
Pearson symbol	:	oC16
Space group number	:	40
Space group symbol	:	*Ama*2
AFLOW prototype command	:	`aflow --proto=AB3_oC16_40_b_a2b --params=`$a, b/a, c/a, z_1, y_2, z_2, y_3, z_3, y_4, z_4$

- This is a more modern version of the $D0_7$ CrO_3 structure.
- (Byström, 1950) give this structure in the $C2cm$ setting of space group #40. We have shifted it to the standard *Ama*2 setting.

Base-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}\,b\,\hat{\mathbf{y}} - \tfrac{1}{2}\,c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}\,b\,\hat{\mathbf{y}} + \tfrac{1}{2}\,c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-z_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$z_1 c\,\hat{\mathbf{z}}$	$(4a)$	O I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 - z_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + z_1 c\,\hat{\mathbf{z}}$	$(4a)$	O I
$\mathbf{B}_3$	=	$\frac{1}{4}\,\mathbf{a}_1 + (y_2 - z_2)\,\mathbf{a}_2 + (y_2 + z_2)\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(4b)$	Cr

$\mathbf{B}_4$	=	$\frac{3}{4}\,\mathbf{a}_1 + (-y_2 - z_2)\,\mathbf{a}_2 + (-y_2 + z_2)\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*b*)	Cr
$\mathbf{B}_5$	=	$\frac{1}{4}\,\mathbf{a}_1 + (y_3 - z_3)\,\mathbf{a}_2 + (y_3 + z_3)\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*b*)	O II
$\mathbf{B}_6$	=	$\frac{3}{4}\,\mathbf{a}_1 + (-y_3 - z_3)\,\mathbf{a}_2 + (-y_3 + z_3)\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*b*)	O II
$\mathbf{B}_7$	=	$\frac{1}{4}\,\mathbf{a}_1 + (y_4 - z_4)\,\mathbf{a}_2 + (y_4 + z_4)\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(4*b*)	O III
$\mathbf{B}_8$	=	$\frac{3}{4}\,\mathbf{a}_1 + (-y_4 - z_4)\,\mathbf{a}_2 + (-y_4 + z_4)\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(4*b*)	O III

References:

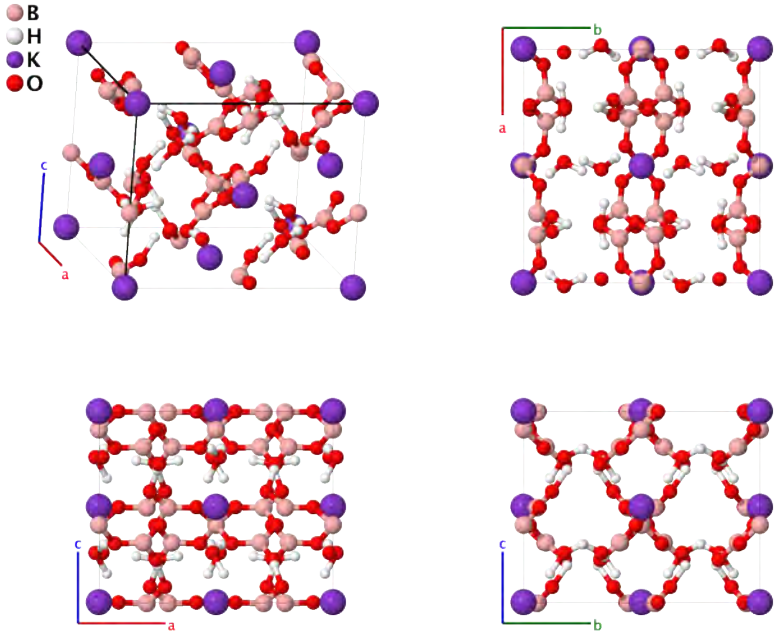
- A. Byström and K.-A. Wilhelmi, *The Crystal Structure of Chromium Trioxide*, Acta Chem. Scand. **4**, 1131–1141 (1950), doi:10.3891/acta.chem.scand.04-1131.
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturebericht Band II, 1928-1932* (Akademsiche Verlagsgesellschaft M. B. H, Leipzig, 1937).

Geometry files:

- CIF: pp. 1599
- POSCAR: pp. 1599

Santite ($KB_5O_8{\cdot}4H_2O$, $K3_5$) Structure: A5B8CD12_oC104_41_a2b_4b_a_6b

http://aflow.org/prototype-encyclopedia/A5B8CD12_oC104_41_a2b_4b_a_6b

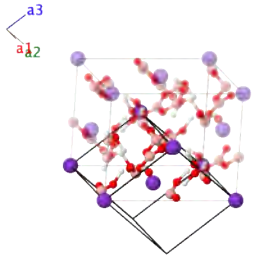


Prototype	:	$B_5H_8KO_{12}$
AFLOW prototype label	:	A5B8CD12_oC104_41_a2b_4b_a_6b
***Strukturbericht* designation**	:	$K3_5$
Pearson symbol	:	oC104
Space group number	:	41
Space group symbol	:	*Aba*2
AFLOW prototype command	:	`aflow --proto=A5B8CD12_oC104_41_a2b_4b_a_6b --params=`$a, b/a, c/a, z_1, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}$

- (Zachariasen, 1938) originally determined the structure of this compound, without knowing the positions of the hydrogens, but believed the structure to be correctly written as $KH_2(H_3O)_2B_5O_{10}$. (Gottfried, 1940) gave this the *Strukturbericht* designation $K3_5$. (Zachariasen, 1963) did a refinement of the structure, including the hydrogen positions, noting that the structure should properly be written $K[B_5O_6(OH)_4]{\cdot}2H_2O$. As neither the positions of the heavy atoms nor the space group have changed, we retain the $K3_5$ designation, but we give the prototype as $KB_5O_8{\cdot}4H_2O$, which seems to be the standard formula even though it is not structurally correct.

Base-centered Orthorhombic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= a\,\hat{\mathbf{x}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,b\,\hat{\mathbf{y}} - \tfrac{1}{2}\,c\,\hat{\mathbf{z}}\\ \mathbf{a}_3 &= \tfrac{1}{2}\,b\,\hat{\mathbf{y}} + \tfrac{1}{2}\,c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-z_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$z_1 c\,\hat{\mathbf{z}}$	(4*a*)	B I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \left(\frac{1}{2} - z_1\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(4*a*)	B I
$\mathbf{B}_3$	=	$-z_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$z_2 c\,\hat{\mathbf{z}}$	(4*a*)	K
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \left(\frac{1}{2} - z_2\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*a*)	K
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + (y_3 - z_3)\,\mathbf{a}_2 + (y_3 + z_3)\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(8*b*)	B II
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 + (-y_3 - z_3)\,\mathbf{a}_2 + (-y_3 + z_3)\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(8*b*)	B II
$\mathbf{B}_7$	=	$\left(\frac{1}{2} + x_3\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - y_3 - z_3\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - y_3 + z_3\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(8*b*)	B II
$\mathbf{B}_8$	=	$\left(\frac{1}{2} - x_3\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + y_3 - z_3\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + y_3 + z_3\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(8*b*)	B II
$\mathbf{B}_9$	=	$x_4\,\mathbf{a}_1 + (y_4 - z_4)\,\mathbf{a}_2 + (y_4 + z_4)\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*b*)	B III
$\mathbf{B}_{10}$	=	$-x_4\,\mathbf{a}_1 + (-y_4 - z_4)\,\mathbf{a}_2 + (-y_4 + z_4)\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*b*)	B III
$\mathbf{B}_{11}$	=	$\left(\frac{1}{2} + x_4\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - y_4 - z_4\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - y_4 + z_4\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*b*)	B III
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} - x_4\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4 - z_4\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + y_4 + z_4\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*b*)	B III
$\mathbf{B}_{13}$	=	$x_5\,\mathbf{a}_1 + (y_5 - z_5)\,\mathbf{a}_2 + (y_5 + z_5)\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(8*b*)	H I
$\mathbf{B}_{14}$	=	$-x_5\,\mathbf{a}_1 + (-y_5 - z_5)\,\mathbf{a}_2 + (-y_5 + z_5)\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(8*b*)	H I
$\mathbf{B}_{15}$	=	$\left(\frac{1}{2} + x_5\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - y_5 - z_5\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - y_5 + z_5\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right)b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(8*b*)	H I
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} - x_5\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + y_5 - z_5\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + y_5 + z_5\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right)b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(8*b*)	H I
$\mathbf{B}_{17}$	=	$x_6\,\mathbf{a}_1 + (y_6 - z_6)\,\mathbf{a}_2 + (y_6 + z_6)\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(8*b*)	H II
$\mathbf{B}_{18}$	=	$-x_6\,\mathbf{a}_1 + (-y_6 - z_6)\,\mathbf{a}_2 + (-y_6 + z_6)\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(8*b*)	H II
$\mathbf{B}_{19}$	=	$\left(\frac{1}{2} + x_6\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - y_6 - z_6\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - y_6 + z_6\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right)b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(8*b*)	H II

$$\mathbf{B}_{20} = \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_6 - z_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + y_6 + z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_6\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_6\right)b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}} \quad (8b) \quad \text{H II}$$

$$\mathbf{B}_{21} = x_7\mathbf{a}_1 + (y_7 - z_7)\mathbf{a}_2 + (y_7 + z_7)\mathbf{a}_3 = x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}} \quad (8b) \quad \text{H III}$$

$$\mathbf{B}_{22} = -x_7\mathbf{a}_1 + (-y_7 - z_7)\mathbf{a}_2 + (-y_7 + z_7)\mathbf{a}_3 = -x_7 a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}} \quad (8b) \quad \text{H III}$$

$$\mathbf{B}_{23} = \left(\tfrac{1}{2} + x_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_7 - z_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - y_7 + z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_7\right)b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}} \quad (8b) \quad \text{H III}$$

$$\mathbf{B}_{24} = \left(\tfrac{1}{2} - x_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_7 - z_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + y_7 + z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_7\right)b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}} \quad (8b) \quad \text{H III}$$

$$\mathbf{B}_{25} = x_8\mathbf{a}_1 + (y_8 - z_8)\mathbf{a}_2 + (y_8 + z_8)\mathbf{a}_3 = x_8 a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}} \quad (8b) \quad \text{H IV}$$

$$\mathbf{B}_{26} = -x_8\mathbf{a}_1 + (-y_8 - z_8)\mathbf{a}_2 + (-y_8 + z_8)\mathbf{a}_3 = -x_8 a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}} \quad (8b) \quad \text{H IV}$$

$$\mathbf{B}_{27} = \left(\tfrac{1}{2} + x_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_8 - z_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - y_8 + z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_8\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_8\right)b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}} \quad (8b) \quad \text{H IV}$$

$$\mathbf{B}_{28} = \left(\tfrac{1}{2} - x_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_8 - z_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + y_8 + z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_8\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_8\right)b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}} \quad (8b) \quad \text{H IV}$$

$$\mathbf{B}_{29} = x_9\mathbf{a}_1 + (y_9 - z_9)\mathbf{a}_2 + (y_9 + z_9)\mathbf{a}_3 = x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O I}$$

$$\mathbf{B}_{30} = -x_9\mathbf{a}_1 + (-y_9 - z_9)\mathbf{a}_2 + (-y_9 + z_9)\mathbf{a}_3 = -x_9 a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O I}$$

$$\mathbf{B}_{31} = \left(\tfrac{1}{2} + x_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_9 - z_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - y_9 + z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_9\right)b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O I}$$

$$\mathbf{B}_{32} = \left(\tfrac{1}{2} - x_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_9 - z_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + y_9 + z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_9\right)b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O I}$$

$$\mathbf{B}_{33} = x_{10}\mathbf{a}_1 + (y_{10} - z_{10})\mathbf{a}_2 + (y_{10} + z_{10})\mathbf{a}_3 = x_{10} a\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O II}$$

$$\mathbf{B}_{34} = -x_{10}\mathbf{a}_1 + (-y_{10} - z_{10})\mathbf{a}_2 + (-y_{10} + z_{10})\mathbf{a}_3 = -x_{10} a\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O II}$$

$$\mathbf{B}_{35} = \left(\tfrac{1}{2} + x_{10}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{10} - z_{10}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - y_{10} + z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{10}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{10}\right)b\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O II}$$

$$\mathbf{B}_{36} = \left(\tfrac{1}{2} - x_{10}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{10} - z_{10}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + y_{10} + z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{10}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{10}\right)b\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O II}$$

$$\mathbf{B}_{37} = x_{11}\mathbf{a}_1 + (y_{11} - z_{11})\mathbf{a}_2 + (y_{11} + z_{11})\mathbf{a}_3 = x_{11} a\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + z_{11} c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O III}$$

$$\mathbf{B}_{38} = -x_{11}\mathbf{a}_1 + (-y_{11} - z_{11})\mathbf{a}_2 + (-y_{11} + z_{11})\mathbf{a}_3 = -x_{11} a\,\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} + z_{11} c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O III}$$

$$\mathbf{B}_{39} = \left(\tfrac{1}{2} + x_{11}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{11} - z_{11}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - y_{11} + z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{11}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{11}\right)b\,\hat{\mathbf{y}} + z_{11} c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O III}$$

$$\mathbf{B}_{40} = \left(\tfrac{1}{2} - x_{11}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{11} - z_{11}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + y_{11} + z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{11}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{11}\right)b\,\hat{\mathbf{y}} + z_{11} c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O III}$$

$$\mathbf{B}_{41} = x_{12}\mathbf{a}_1 + (y_{12} - z_{12})\mathbf{a}_2 + (y_{12} + z_{12})\mathbf{a}_3 = x_{12} a\,\hat{\mathbf{x}} + y_{12} b\,\hat{\mathbf{y}} + z_{12} c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O IV}$$

$$\mathbf{B}_{42} = -x_{12}\mathbf{a}_1 + (-y_{12} - z_{12})\mathbf{a}_2 + (-y_{12} + z_{12})\mathbf{a}_3 = -x_{12} a\,\hat{\mathbf{x}} - y_{12} b\,\hat{\mathbf{y}} + z_{12} c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O IV}$$

$$\mathbf{B}_{43} = \left(\tfrac{1}{2}+x_{12}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{12}-z_{12}\right)\mathbf{a}_2+\left(\tfrac{1}{2}-y_{12}+z_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{12}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{12}\right)b\,\hat{\mathbf{y}}+z_{12}c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O IV}$$

$$\mathbf{B}_{44} = \left(\tfrac{1}{2}-x_{12}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{12}-z_{12}\right)\mathbf{a}_2+\left(\tfrac{1}{2}+y_{12}+z_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{12}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{12}\right)b\,\hat{\mathbf{y}}+z_{12}c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O IV}$$

$$\mathbf{B}_{45} = x_{13}\,\mathbf{a}_1+(y_{13}-z_{13})\,\mathbf{a}_2+(y_{13}+z_{13})\,\mathbf{a}_3 = x_{13}a\,\hat{\mathbf{x}}+y_{13}b\,\hat{\mathbf{y}}+z_{13}c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O V}$$

$$\mathbf{B}_{46} = -x_{13}\,\mathbf{a}_1+(-y_{13}-z_{13})\,\mathbf{a}_2+(-y_{13}+z_{13})\,\mathbf{a}_3 = -x_{13}a\,\hat{\mathbf{x}}-y_{13}b\,\hat{\mathbf{y}}+z_{13}c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O V}$$

$$\mathbf{B}_{47} = \left(\tfrac{1}{2}+x_{13}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{13}-z_{13}\right)\mathbf{a}_2+\left(\tfrac{1}{2}-y_{13}+z_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{13}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{13}\right)b\,\hat{\mathbf{y}}+z_{13}c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O V}$$

$$\mathbf{B}_{48} = \left(\tfrac{1}{2}-x_{13}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{13}-z_{13}\right)\mathbf{a}_2+\left(\tfrac{1}{2}+y_{13}+z_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{13}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{13}\right)b\,\hat{\mathbf{y}}+z_{13}c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O V}$$

$$\mathbf{B}_{49} = x_{14}\,\mathbf{a}_1+(y_{14}-z_{14})\,\mathbf{a}_2+(y_{14}+z_{14})\,\mathbf{a}_3 = x_{14}a\,\hat{\mathbf{x}}+y_{14}b\,\hat{\mathbf{y}}+z_{14}c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O VI}$$

$$\mathbf{B}_{50} = -x_{14}\,\mathbf{a}_1+(-y_{14}-z_{14})\,\mathbf{a}_2+(-y_{14}+z_{14})\,\mathbf{a}_3 = -x_{14}a\,\hat{\mathbf{x}}-y_{14}b\,\hat{\mathbf{y}}+z_{14}c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O VI}$$

$$\mathbf{B}_{51} = \left(\tfrac{1}{2}+x_{14}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{14}-z_{14}\right)\mathbf{a}_2+\left(\tfrac{1}{2}-y_{14}+z_{14}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{14}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{14}\right)b\,\hat{\mathbf{y}}+z_{14}c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O VI}$$

$$\mathbf{B}_{52} = \left(\tfrac{1}{2}-x_{14}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{14}-z_{14}\right)\mathbf{a}_2+\left(\tfrac{1}{2}+y_{14}+z_{14}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{14}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{14}\right)b\,\hat{\mathbf{y}}+z_{14}c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O VI}$$

References:

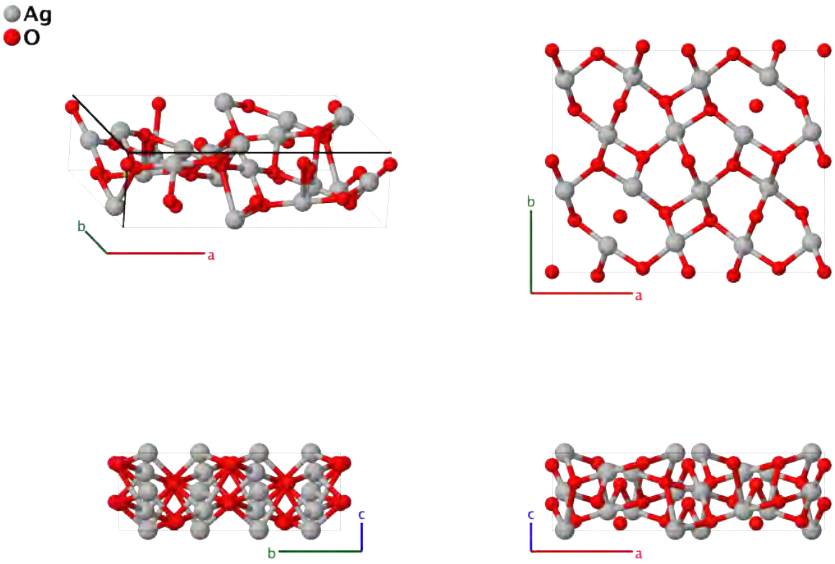
- W. H. Zachariasen and H. A. Plettinger, *Refinement of the structure of potassium pentaborate tetrahydrate*, Acta Cryst. **16**, 376–379 (1963), doi:10.1107/S0365110X63001006.
- W. H. Zachariasen, *The Crystal Structure of Potassium Acid Dihydronium Pentaborate $KH_2(H_3O)_2B_5O_{10}$, (Potassium Pentaborate Tetrahydrate)*, Zeitschrift für Kristallographie - Crystalline Materials **98**, 266–274 (1938), doi:10.1524/zkri.1938.98.1.266.
- C. Gottfried, ed., *Strukturbericht Band V 1937* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1940).

Geometry files:

- CIF: pp. 1599
- POSCAR: pp. 1600

Ag_2O_3 Structure: A2B3_oF40_43_b_ab

http://aflow.org/prototype-encyclopedia/A2B3_oF40_43_b_ab

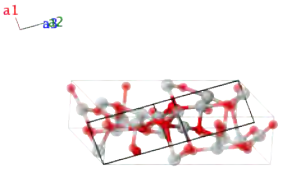


Prototype	:	Ag_2O_3
AFLOW prototype label	:	A2B3_oF40_43_b_ab
***Strukturbericht* designation**	:	None
Pearson symbol	:	oF40
Space group number	:	43
Space group symbol	:	$Fdd2$
AFLOW prototype command	:	`aflow --proto=A2B3_oF40_43_b_ab --params=`$a, b/a, c/a, z_1, x_2, y_2, z_2, x_3, y_3, z_3$

- This structure is a distortion of the $D5_5$ Ag_2O_3 structure (A3B2_cP10_224_d_b), although (Standke, 1986) does not seem to be aware of the earlier work. This is most likely closer to the correct structure for Ag_2O_3 than the $D5_5$ structure is.

Face-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2} b\,\hat{\mathbf{y}} + \frac{1}{2} c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2} a\,\hat{\mathbf{x}} + \frac{1}{2} c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2} a\,\hat{\mathbf{x}} + \frac{1}{2} b\,\hat{\mathbf{y}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$\mathbf{B}_1$	=	$z_1\,\mathbf{a}_1 + z_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$z_1 c\,\hat{\mathbf{z}}$	(8*a*)	O I
$\mathbf{B}_2$	=	$\left(\frac{1}{4}+z_1\right)\mathbf{a}_1 + \left(\frac{1}{4}+z_1\right)\mathbf{a}_2 + \left(\frac{1}{4}-z_1\right)\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{4}+z_1\right)c\,\hat{\mathbf{z}}$	(8*a*)	O I
$\mathbf{B}_3$	=	$(-x_2+y_2+z_2)\,\mathbf{a}_1 + (x_2-y_2+z_2)\,\mathbf{a}_2 + (x_2+y_2-z_2)\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(16*b*)	Ag
$\mathbf{B}_4$	=	$(x_2-y_2+z_2)\,\mathbf{a}_1 + (-x_2+y_2+z_2)\,\mathbf{a}_2 + (-x_2-y_2-z_2)\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(16*b*)	Ag
$\mathbf{B}_5$	=	$\left(\frac{1}{4}-x_2-y_2+z_2\right)\mathbf{a}_1 + \left(\frac{1}{4}+x_2+y_2+z_2\right)\mathbf{a}_2 + \left(\frac{1}{4}+x_2-y_2-z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}+x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4}-y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{4}+z_2\right)c\,\hat{\mathbf{z}}$	(16*b*)	Ag
$\mathbf{B}_6$	=	$\left(\frac{1}{4}+x_2+y_2+z_2\right)\mathbf{a}_1 + \left(\frac{1}{4}-x_2-y_2+z_2\right)\mathbf{a}_2 + \left(\frac{1}{4}-x_2+y_2-z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}-x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4}+y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{4}+z_2\right)c\,\hat{\mathbf{z}}$	(16*b*)	Ag
$\mathbf{B}_7$	=	$(-x_3+y_3+z_3)\,\mathbf{a}_1 + (x_3-y_3+z_3)\,\mathbf{a}_2 + (x_3+y_3-z_3)\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(16*b*)	O II
$\mathbf{B}_8$	=	$(x_3-y_3+z_3)\,\mathbf{a}_1 + (-x_3+y_3+z_3)\,\mathbf{a}_2 + (-x_3-y_3-z_3)\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(16*b*)	O II
$\mathbf{B}_9$	=	$\left(\frac{1}{4}-x_3-y_3+z_3\right)\mathbf{a}_1 + \left(\frac{1}{4}+x_3+y_3+z_3\right)\mathbf{a}_2 + \left(\frac{1}{4}+x_3-y_3-z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}+x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4}-y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{4}+z_3\right)c\,\hat{\mathbf{z}}$	(16*b*)	O II
$\mathbf{B}_{10}$	=	$\left(\frac{1}{4}+x_3+y_3+z_3\right)\mathbf{a}_1 + \left(\frac{1}{4}-x_3-y_3+z_3\right)\mathbf{a}_2 + \left(\frac{1}{4}-x_3+y_3-z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}-x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4}+y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{4}+z_3\right)c\,\hat{\mathbf{z}}$	(16*b*)	O II

References:

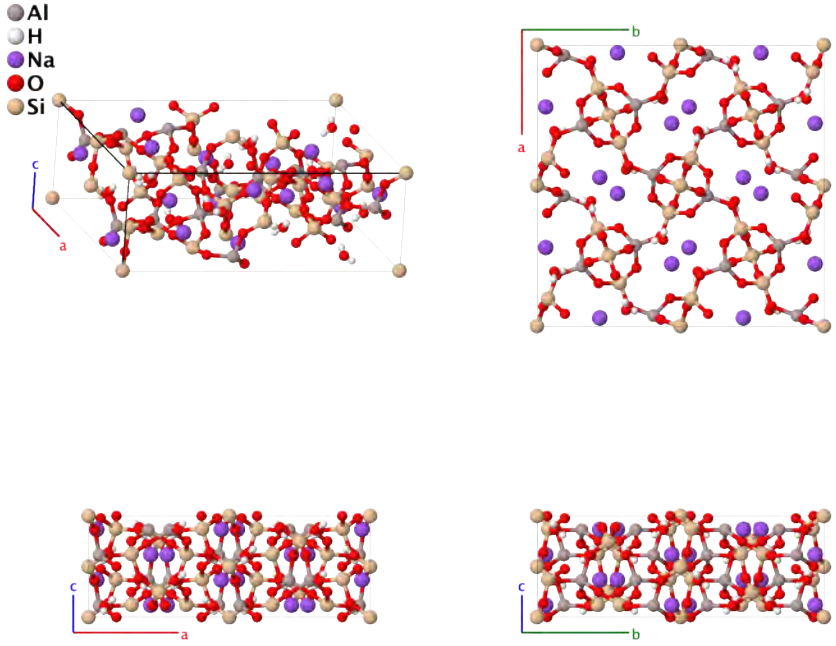
- B. Standke and M. Jansen, *Darstellung und Kristallstruktur von Ag_2O_3*, Z. Anorg. Allg. Chem. **535**, 39–46 (1986), doi:10.1002/zaac.19865350406.

Geometry files:

- CIF: pp. 1600
- POSCAR: pp. 1600

Natrolite ($Na_2Al_2Si_3O_{10}{\cdot}2H_2O$, $S6_{10}$) Structure: A2B4C2D12E3_oF184_43_b_2b_b_6b_ab

http://aflow.org/prototype-encyclopedia/A2B4C2D12E3_oF184_43_b_2b_b_6b_ab

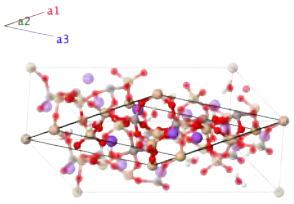


Prototype	:	$Al_2H_4Na_2O_{12}Si_3$
AFLOW prototype label	:	A2B4C2D12E3_oF184_43_b_2b_b_6b_ab
***Strukturbericht* designation**	:	$S6_{10}$
Pearson symbol	:	oF184
Space group number	:	43
Space group symbol	:	$Fdd2$
AFLOW prototype command	:	`aflow --proto=A2B4C2D12E3_oF184_43_b_2b_b_6b_ab --params=`$a, b/a, c/a, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}$

- We use the data from the sample that (Kirfel, 1984) call Crystal II. The origin has been arbitrarily set so that $z_1 = 0$.

Face-centered Orthorhombic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,b\,\hat{\mathbf{y}} + \tfrac{1}{2}\,c\,\hat{\mathbf{z}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,c\,\hat{\mathbf{z}}\\ \mathbf{a}_3 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,b\,\hat{\mathbf{y}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$z_1\,\mathbf{a}_1 + z_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$z_1 c\,\hat{\mathbf{z}}$	(8*a*)	Si I
$\mathbf{B}_2$	=	$\left(\frac{1}{4}+z_1\right)\mathbf{a}_1 + \left(\frac{1}{4}+z_1\right)\mathbf{a}_2 + \left(\frac{1}{4}-z_1\right)\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{4}+z_1\right)c\,\hat{\mathbf{z}}$	(8*a*)	Si I
$\mathbf{B}_3$	=	$(-x_2+y_2+z_2)\,\mathbf{a}_1 + (x_2-y_2+z_2)\,\mathbf{a}_2 + (x_2+y_2-z_2)\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(16*b*)	Al
$\mathbf{B}_4$	=	$(x_2-y_2+z_2)\,\mathbf{a}_1 + (-x_2+y_2+z_2)\,\mathbf{a}_2 + (-x_2-y_2-z_2)\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(16*b*)	Al
$\mathbf{B}_5$	=	$\left(\frac{1}{4}-x_2-y_2+z_2\right)\mathbf{a}_1 + \left(\frac{1}{4}+x_2+y_2+z_2\right)\mathbf{a}_2 + \left(\frac{1}{4}+x_2-y_2-z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}+x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4}-y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{4}+z_2\right)c\,\hat{\mathbf{z}}$	(16*b*)	Al
$\mathbf{B}_6$	=	$\left(\frac{1}{4}+x_2+y_2+z_2\right)\mathbf{a}_1 + \left(\frac{1}{4}-x_2-y_2+z_2\right)\mathbf{a}_2 + \left(\frac{1}{4}-x_2+y_2-z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}-x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4}+y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{4}+z_2\right)c\,\hat{\mathbf{z}}$	(16*b*)	Al
$\mathbf{B}_7$	=	$(-x_3+y_3+z_3)\,\mathbf{a}_1 + (x_3-y_3+z_3)\,\mathbf{a}_2 + (x_3+y_3-z_3)\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(16*b*)	H I
$\mathbf{B}_8$	=	$(x_3-y_3+z_3)\,\mathbf{a}_1 + (-x_3+y_3+z_3)\,\mathbf{a}_2 + (-x_3-y_3-z_3)\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(16*b*)	H I
$\mathbf{B}_9$	=	$\left(\frac{1}{4}-x_3-y_3+z_3\right)\mathbf{a}_1 + \left(\frac{1}{4}+x_3+y_3+z_3\right)\mathbf{a}_2 + \left(\frac{1}{4}+x_3-y_3-z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}+x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4}-y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{4}+z_3\right)c\,\hat{\mathbf{z}}$	(16*b*)	H I
$\mathbf{B}_{10}$	=	$\left(\frac{1}{4}+x_3+y_3+z_3\right)\mathbf{a}_1 + \left(\frac{1}{4}-x_3-y_3+z_3\right)\mathbf{a}_2 + \left(\frac{1}{4}-x_3+y_3-z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}-x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4}+y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{4}+z_3\right)c\,\hat{\mathbf{z}}$	(16*b*)	H I
$\mathbf{B}_{11}$	=	$(-x_4+y_4+z_4)\,\mathbf{a}_1 + (x_4-y_4+z_4)\,\mathbf{a}_2 + (x_4+y_4-z_4)\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(16*b*)	H II
$\mathbf{B}_{12}$	=	$(x_4-y_4+z_4)\,\mathbf{a}_1 + (-x_4+y_4+z_4)\,\mathbf{a}_2 + (-x_4-y_4-z_4)\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(16*b*)	H II

$\mathbf{B}_{13}$	=	$\left(\frac{1}{4} - x_4 - y_4 + z_4\right)\mathbf{a}_1 + \left(\frac{1}{4} + x_4 + y_4 + z_4\right)\mathbf{a}_2 + \left(\frac{1}{4} + x_4 - y_4 - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_4\right)c\,\hat{\mathbf{z}}$	(16*b*)	H II
$\mathbf{B}_{14}$	=	$\left(\frac{1}{4} + x_4 + y_4 + z_4\right)\mathbf{a}_1 + \left(\frac{1}{4} - x_4 - y_4 + z_4\right)\mathbf{a}_2 + \left(\frac{1}{4} - x_4 + y_4 - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_4\right)c\,\hat{\mathbf{z}}$	(16*b*)	H II
$\mathbf{B}_{15}$	=	$(-x_5 + y_5 + z_5)\,\mathbf{a}_1 + (x_5 - y_5 + z_5)\,\mathbf{a}_2 + (x_5 + y_5 - z_5)\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(16*b*)	Na
$\mathbf{B}_{16}$	=	$(x_5 - y_5 + z_5)\,\mathbf{a}_1 + (-x_5 + y_5 + z_5)\,\mathbf{a}_2 + (-x_5 - y_5 - z_5)\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(16*b*)	Na
$\mathbf{B}_{17}$	=	$\left(\frac{1}{4} - x_5 - y_5 + z_5\right)\mathbf{a}_1 + \left(\frac{1}{4} + x_5 + y_5 + z_5\right)\mathbf{a}_2 + \left(\frac{1}{4} + x_5 - y_5 - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_5\right)c\,\hat{\mathbf{z}}$	(16*b*)	Na
$\mathbf{B}_{18}$	=	$\left(\frac{1}{4} + x_5 + y_5 + z_5\right)\mathbf{a}_1 + \left(\frac{1}{4} - x_5 - y_5 + z_5\right)\mathbf{a}_2 + \left(\frac{1}{4} - x_5 + y_5 - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_5\right)c\,\hat{\mathbf{z}}$	(16*b*)	Na
$\mathbf{B}_{19}$	=	$(-x_6 + y_6 + z_6)\,\mathbf{a}_1 + (x_6 - y_6 + z_6)\,\mathbf{a}_2 + (x_6 + y_6 - z_6)\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(16*b*)	O I
$\mathbf{B}_{20}$	=	$(x_6 - y_6 + z_6)\,\mathbf{a}_1 + (-x_6 + y_6 + z_6)\,\mathbf{a}_2 + (-x_6 - y_6 - z_6)\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(16*b*)	O I
$\mathbf{B}_{21}$	=	$\left(\frac{1}{4} - x_6 - y_6 + z_6\right)\mathbf{a}_1 + \left(\frac{1}{4} + x_6 + y_6 + z_6\right)\mathbf{a}_2 + \left(\frac{1}{4} + x_6 - y_6 - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - y_6\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_6\right)c\,\hat{\mathbf{z}}$	(16*b*)	O I
$\mathbf{B}_{22}$	=	$\left(\frac{1}{4} + x_6 + y_6 + z_6\right)\mathbf{a}_1 + \left(\frac{1}{4} - x_6 - y_6 + z_6\right)\mathbf{a}_2 + \left(\frac{1}{4} - x_6 + y_6 - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + y_6\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_6\right)c\,\hat{\mathbf{z}}$	(16*b*)	O I
$\mathbf{B}_{23}$	=	$(-x_7 + y_7 + z_7)\,\mathbf{a}_1 + (x_7 - y_7 + z_7)\,\mathbf{a}_2 + (x_7 + y_7 - z_7)\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(16*b*)	O II
$\mathbf{B}_{24}$	=	$(x_7 - y_7 + z_7)\,\mathbf{a}_1 + (-x_7 + y_7 + z_7)\,\mathbf{a}_2 + (-x_7 - y_7 - z_7)\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(16*b*)	O II
$\mathbf{B}_{25}$	=	$\left(\frac{1}{4} - x_7 - y_7 + z_7\right)\mathbf{a}_1 + \left(\frac{1}{4} + x_7 + y_7 + z_7\right)\mathbf{a}_2 + \left(\frac{1}{4} + x_7 - y_7 - z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - y_7\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_7\right)c\,\hat{\mathbf{z}}$	(16*b*)	O II
$\mathbf{B}_{26}$	=	$\left(\frac{1}{4} + x_7 + y_7 + z_7\right)\mathbf{a}_1 + \left(\frac{1}{4} - x_7 - y_7 + z_7\right)\mathbf{a}_2 + \left(\frac{1}{4} - x_7 + y_7 - z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + y_7\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_7\right)c\,\hat{\mathbf{z}}$	(16*b*)	O II
$\mathbf{B}_{27}$	=	$(-x_8 + y_8 + z_8)\,\mathbf{a}_1 + (x_8 - y_8 + z_8)\,\mathbf{a}_2 + (x_8 + y_8 - z_8)\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(16*b*)	O III

$$\mathbf{B}_{28} = (x_8 - y_8 + z_8)\,\mathbf{a}_1 + (-x_8 + y_8 + z_8)\,\mathbf{a}_2 + (-x_8 - y_8 - z_8)\,\mathbf{a}_3 = -x_8 a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}} \quad (16b) \quad \text{O III}$$

$$\mathbf{B}_{29} = \left(\tfrac{1}{4} - x_8 - y_8 + z_8\right)\mathbf{a}_1 + \left(\tfrac{1}{4} + x_8 + y_8 + z_8\right)\mathbf{a}_2 + \left(\tfrac{1}{4} + x_8 - y_8 - z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + x_8\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - y_8\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_8\right)c\,\hat{\mathbf{z}} \quad (16b) \quad \text{O III}$$

$$\mathbf{B}_{30} = \left(\tfrac{1}{4} + x_8 + y_8 + z_8\right)\mathbf{a}_1 + \left(\tfrac{1}{4} - x_8 - y_8 + z_8\right)\mathbf{a}_2 + \left(\tfrac{1}{4} - x_8 + y_8 - z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - x_8\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + y_8\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_8\right)c\,\hat{\mathbf{z}} \quad (16b) \quad \text{O III}$$

$$\mathbf{B}_{31} = (-x_9 + y_9 + z_9)\,\mathbf{a}_1 + (x_9 - y_9 + z_9)\,\mathbf{a}_2 + (x_9 + y_9 - z_9)\,\mathbf{a}_3 = x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}} \quad (16b) \quad \text{O IV}$$

$$\mathbf{B}_{32} = (x_9 - y_9 + z_9)\,\mathbf{a}_1 + (-x_9 + y_9 + z_9)\,\mathbf{a}_2 + (-x_9 - y_9 - z_9)\,\mathbf{a}_3 = -x_9 a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}} \quad (16b) \quad \text{O IV}$$

$$\mathbf{B}_{33} = \left(\tfrac{1}{4} - x_9 - y_9 + z_9\right)\mathbf{a}_1 + \left(\tfrac{1}{4} + x_9 + y_9 + z_9\right)\mathbf{a}_2 + \left(\tfrac{1}{4} + x_9 - y_9 - z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + x_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - y_9\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_9\right)c\,\hat{\mathbf{z}} \quad (16b) \quad \text{O IV}$$

$$\mathbf{B}_{34} = \left(\tfrac{1}{4} + x_9 + y_9 + z_9\right)\mathbf{a}_1 + \left(\tfrac{1}{4} - x_9 - y_9 + z_9\right)\mathbf{a}_2 + \left(\tfrac{1}{4} - x_9 + y_9 - z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - x_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + y_9\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_9\right)c\,\hat{\mathbf{z}} \quad (16b) \quad \text{O IV}$$

$$\mathbf{B}_{35} = (-x_{10} + y_{10} + z_{10})\,\mathbf{a}_1 + (x_{10} - y_{10} + z_{10})\,\mathbf{a}_2 + (x_{10} + y_{10} - z_{10})\,\mathbf{a}_3 = x_{10} a\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}} \quad (16b) \quad \text{O V}$$

$$\mathbf{B}_{36} = (x_{10} - y_{10} + z_{10})\,\mathbf{a}_1 + (-x_{10} + y_{10} + z_{10})\,\mathbf{a}_2 + (-x_{10} - y_{10} - z_{10})\,\mathbf{a}_3 = -x_{10} a\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}} \quad (16b) \quad \text{O V}$$

$$\mathbf{B}_{37} = \left(\tfrac{1}{4} - x_{10} - y_{10} + z_{10}\right)\mathbf{a}_1 + \left(\tfrac{1}{4} + x_{10} + y_{10} + z_{10}\right)\mathbf{a}_2 + \left(\tfrac{1}{4} + x_{10} - y_{10} - z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + x_{10}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - y_{10}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_{10}\right)c\,\hat{\mathbf{z}} \quad (16b) \quad \text{O V}$$

$$\mathbf{B}_{38} = \left(\tfrac{1}{4} + x_{10} + y_{10} + z_{10}\right)\mathbf{a}_1 + \left(\tfrac{1}{4} - x_{10} - y_{10} + z_{10}\right)\mathbf{a}_2 + \left(\tfrac{1}{4} - x_{10} + y_{10} - z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - x_{10}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + y_{10}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_{10}\right)c\,\hat{\mathbf{z}} \quad (16b) \quad \text{O V}$$

$$\mathbf{B}_{39} = (-x_{11} + y_{11} + z_{11})\,\mathbf{a}_1 + (x_{11} - y_{11} + z_{11})\,\mathbf{a}_2 + (x_{11} + y_{11} - z_{11})\,\mathbf{a}_3 = x_{11} a\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + z_{11} c\,\hat{\mathbf{z}} \quad (16b) \quad \text{O VI}$$

$$\mathbf{B}_{40} = (x_{11} - y_{11} + z_{11})\,\mathbf{a}_1 + (-x_{11} + y_{11} + z_{11})\,\mathbf{a}_2 + (-x_{11} - y_{11} - z_{11})\,\mathbf{a}_3 = -x_{11} a\,\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} + z_{11} c\,\hat{\mathbf{z}} \quad (16b) \quad \text{O VI}$$

$$\mathbf{B}_{41} = \left(\tfrac{1}{4} - x_{11} - y_{11} + z_{11}\right)\mathbf{a}_1 + \left(\tfrac{1}{4} + x_{11} + y_{11} + z_{11}\right)\mathbf{a}_2 + \left(\tfrac{1}{4} + x_{11} - y_{11} - z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + x_{11}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - y_{11}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_{11}\right)c\,\hat{\mathbf{z}} \quad (16b) \quad \text{O VI}$$

$$\mathbf{B}_{42} = \left(\tfrac{1}{4} + x_{11} + y_{11} + z_{11}\right)\mathbf{a}_1 + \left(\tfrac{1}{4} - x_{11} - y_{11} + z_{11}\right)\mathbf{a}_2 + \left(\tfrac{1}{4} - x_{11} + y_{11} - z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - x_{11}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + y_{11}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_{11}\right)c\,\hat{\mathbf{z}} \quad (16b) \quad \text{O VI}$$

$\mathbf{B}_{43}$	=	$(-x_{12}+y_{12}+z_{12})\,\mathbf{a}_1+$ $(x_{12}-y_{12}+z_{12})\,\mathbf{a}_2+$ $(x_{12}+y_{12}-z_{12})\,\mathbf{a}_3$	=	$x_{12}a\,\hat{\mathbf{x}}+y_{12}b\,\hat{\mathbf{y}}+z_{12}c\,\hat{\mathbf{z}}$	(16*b*)	Si II
$\mathbf{B}_{44}$	=	$(x_{12}-y_{12}+z_{12})\,\mathbf{a}_1+$ $(-x_{12}+y_{12}+z_{12})\,\mathbf{a}_2+$ $(-x_{12}-y_{12}-z_{12})\,\mathbf{a}_3$	=	$-x_{12}a\,\hat{\mathbf{x}}-y_{12}b\,\hat{\mathbf{y}}+z_{12}c\,\hat{\mathbf{z}}$	(16*b*)	Si II
$\mathbf{B}_{45}$	=	$\left(\frac{1}{4}-x_{12}-y_{12}+z_{12}\right)\mathbf{a}_1+$ $\left(\frac{1}{4}+x_{12}+y_{12}+z_{12}\right)\mathbf{a}_2+$ $\left(\frac{1}{4}+x_{12}-y_{12}-z_{12}\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}+x_{12}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-y_{12}\right)b\,\hat{\mathbf{y}}+$ $\left(\frac{1}{4}+z_{12}\right)c\,\hat{\mathbf{z}}$	(16*b*)	Si II
$\mathbf{B}_{46}$	=	$\left(\frac{1}{4}+x_{12}+y_{12}+z_{12}\right)\mathbf{a}_1+$ $\left(\frac{1}{4}-x_{12}-y_{12}+z_{12}\right)\mathbf{a}_2+$ $\left(\frac{1}{4}-x_{12}+y_{12}-z_{12}\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}-x_{12}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+y_{12}\right)b\,\hat{\mathbf{y}}+$ $\left(\frac{1}{4}+z_{12}\right)c\,\hat{\mathbf{z}}$	(16*b*)	Si II

References:

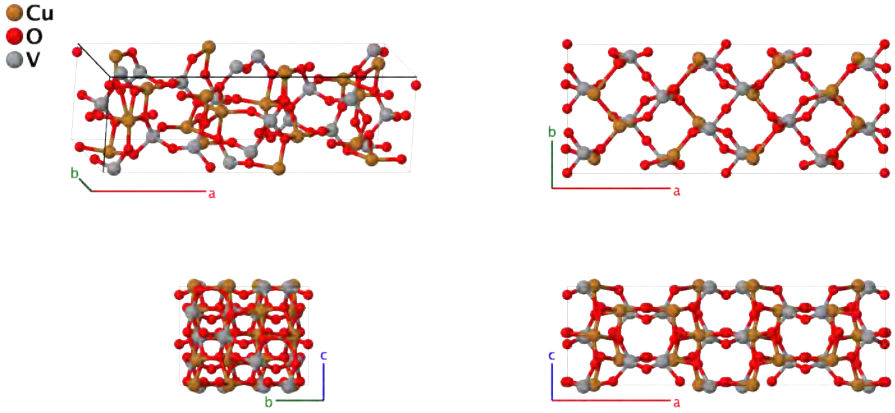
- A. Kirfel, M. Orthen, and G. Will, *Natrolite: refinement of the crystal structure of two samples from Marienberg (Usti nad Labem, CSSR)*, Zeolites **4**, 140–146 (1984), doi:10.1016/0144-2449(84)90052-6.

Geometry files:

- CIF: pp. 1600
- POSCAR: pp. 1601

Blossite (α-$Cu_2V_2O_7$) Structure: A2B7C2_oF88_43_b_a3b_b

http://aflow.org/prototype-encyclopedia/A2B7C2_oF88_43_b_a3b_b

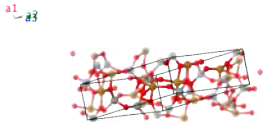


Prototype	:	$Cu_2O_7V_2$
AFLOW prototype label	:	A2B7C2_oF88_43_b_a3b_b
***Strukturbericht* designation**	:	None
Pearson symbol	:	oF88
Space group number	:	43
Space group symbol	:	$Fdd2$
AFLOW prototype command	:	`aflow --proto=A2B7C2_oF88_43_b_a3b_b --params=`$a, b/a, c/a, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6$

- This structure was given the name "blossite" by (Robinson, 1987).
- Space group $Fdd2$ #43 does not fix the $z = 0$ plane. We do this by setting $z_2 = 3/4$ for the copper atom.

Face-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2} b\,\hat{\mathbf{y}} + \frac{1}{2} c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2} a\,\hat{\mathbf{x}} + \frac{1}{2} c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2} a\,\hat{\mathbf{x}} + \frac{1}{2} b\,\hat{\mathbf{y}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$z_1\,\mathbf{a}_1 + z_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$z_1 c\,\hat{\mathbf{z}}$	($8a$)	O I
$\mathbf{B}_2$	=	$\left(\frac{1}{4} + z_1\right)\mathbf{a}_1 + \left(\frac{1}{4} + z_1\right)\mathbf{a}_2 + \left(\frac{1}{4} - z_1\right)\mathbf{a}_3$	=	$\frac{1}{4} a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_1\right) c\,\hat{\mathbf{z}}$	($8a$)	O I
$\mathbf{B}_3$	=	$(-x_2 + y_2 + z_2)\,\mathbf{a}_1 + (x_2 - y_2 + z_2)\,\mathbf{a}_2 + (x_2 + y_2 - z_2)\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	($16b$)	Cu

$\mathbf{B}_4$	=	$(x_2 - y_2 + z_2)\,\mathbf{a}_1 +$ $(-x_2 + y_2 + z_2)\,\mathbf{a}_2 +$ $(-x_2 - y_2 - z_2)\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(16*b*)	Cu
$\mathbf{B}_5$	=	$\left(\frac{1}{4} - x_2 - y_2 + z_2\right)\mathbf{a}_1 +$ $\left(\frac{1}{4} + x_2 + y_2 + z_2\right)\mathbf{a}_2 +$ $\left(\frac{1}{4} + x_2 - y_2 - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - y_2\right)b\,\hat{\mathbf{y}} +$ $\left(\frac{1}{4} + z_2\right)c\,\hat{\mathbf{z}}$	(16*b*)	Cu
$\mathbf{B}_6$	=	$\left(\frac{1}{4} + x_2 + y_2 + z_2\right)\mathbf{a}_1 +$ $\left(\frac{1}{4} - x_2 - y_2 + z_2\right)\mathbf{a}_2 +$ $\left(\frac{1}{4} - x_2 + y_2 - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + y_2\right)b\,\hat{\mathbf{y}} +$ $\left(\frac{1}{4} + z_2\right)c\,\hat{\mathbf{z}}$	(16*b*)	Cu
$\mathbf{B}_7$	=	$(-x_3 + y_3 + z_3)\,\mathbf{a}_1 +$ $(x_3 - y_3 + z_3)\,\mathbf{a}_2 +$ $(x_3 + y_3 - z_3)\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(16*b*)	O II
$\mathbf{B}_8$	=	$(x_3 - y_3 + z_3)\,\mathbf{a}_1 +$ $(-x_3 + y_3 + z_3)\,\mathbf{a}_2 +$ $(-x_3 - y_3 - z_3)\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(16*b*)	O II
$\mathbf{B}_9$	=	$\left(\frac{1}{4} - x_3 - y_3 + z_3\right)\mathbf{a}_1 +$ $\left(\frac{1}{4} + x_3 + y_3 + z_3\right)\mathbf{a}_2 +$ $\left(\frac{1}{4} + x_3 - y_3 - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - y_3\right)b\,\hat{\mathbf{y}} +$ $\left(\frac{1}{4} + z_3\right)c\,\hat{\mathbf{z}}$	(16*b*)	O II
$\mathbf{B}_{10}$	=	$\left(\frac{1}{4} + x_3 + y_3 + z_3\right)\mathbf{a}_1 +$ $\left(\frac{1}{4} - x_3 - y_3 + z_3\right)\mathbf{a}_2 +$ $\left(\frac{1}{4} - x_3 + y_3 - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + y_3\right)b\,\hat{\mathbf{y}} +$ $\left(\frac{1}{4} + z_3\right)c\,\hat{\mathbf{z}}$	(16*b*)	O II
$\mathbf{B}_{11}$	=	$(-x_4 + y_4 + z_4)\,\mathbf{a}_1 +$ $(x_4 - y_4 + z_4)\,\mathbf{a}_2 +$ $(x_4 + y_4 - z_4)\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(16*b*)	O III
$\mathbf{B}_{12}$	=	$(x_4 - y_4 + z_4)\,\mathbf{a}_1 +$ $(-x_4 + y_4 + z_4)\,\mathbf{a}_2 +$ $(-x_4 - y_4 - z_4)\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(16*b*)	O III
$\mathbf{B}_{13}$	=	$\left(\frac{1}{4} - x_4 - y_4 + z_4\right)\mathbf{a}_1 +$ $\left(\frac{1}{4} + x_4 + y_4 + z_4\right)\mathbf{a}_2 +$ $\left(\frac{1}{4} + x_4 - y_4 - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - y_4\right)b\,\hat{\mathbf{y}} +$ $\left(\frac{1}{4} + z_4\right)c\,\hat{\mathbf{z}}$	(16*b*)	O III
$\mathbf{B}_{14}$	=	$\left(\frac{1}{4} + x_4 + y_4 + z_4\right)\mathbf{a}_1 +$ $\left(\frac{1}{4} - x_4 - y_4 + z_4\right)\mathbf{a}_2 +$ $\left(\frac{1}{4} - x_4 + y_4 - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + y_4\right)b\,\hat{\mathbf{y}} +$ $\left(\frac{1}{4} + z_4\right)c\,\hat{\mathbf{z}}$	(16*b*)	O III
$\mathbf{B}_{15}$	=	$(-x_5 + y_5 + z_5)\,\mathbf{a}_1 +$ $(x_5 - y_5 + z_5)\,\mathbf{a}_2 +$ $(x_5 + y_5 - z_5)\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(16*b*)	O IV
$\mathbf{B}_{16}$	=	$(x_5 - y_5 + z_5)\,\mathbf{a}_1 +$ $(-x_5 + y_5 + z_5)\,\mathbf{a}_2 +$ $(-x_5 - y_5 - z_5)\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(16*b*)	O IV
$\mathbf{B}_{17}$	=	$\left(\frac{1}{4} - x_5 - y_5 + z_5\right)\mathbf{a}_1 +$ $\left(\frac{1}{4} + x_5 + y_5 + z_5\right)\mathbf{a}_2 +$ $\left(\frac{1}{4} + x_5 - y_5 - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - y_5\right)b\,\hat{\mathbf{y}} +$ $\left(\frac{1}{4} + z_5\right)c\,\hat{\mathbf{z}}$	(16*b*)	O IV
$\mathbf{B}_{18}$	=	$\left(\frac{1}{4} + x_5 + y_5 + z_5\right)\mathbf{a}_1 +$ $\left(\frac{1}{4} - x_5 - y_5 + z_5\right)\mathbf{a}_2 +$ $\left(\frac{1}{4} - x_5 + y_5 - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + y_5\right)b\,\hat{\mathbf{y}} +$ $\left(\frac{1}{4} + z_5\right)c\,\hat{\mathbf{z}}$	(16*b*)	O IV

$\mathbf{B}_{19}$	=	$(-x_6 + y_6 + z_6)\,\mathbf{a}_1 +$ $(x_6 - y_6 + z_6)\,\mathbf{a}_2 +$ $(x_6 + y_6 - z_6)\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(16*b*)	V
$\mathbf{B}_{20}$	=	$(x_6 - y_6 + z_6)\,\mathbf{a}_1 +$ $(-x_6 + y_6 + z_6)\,\mathbf{a}_2 +$ $(-x_6 - y_6 - z_6)\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(16*b*)	V
$\mathbf{B}_{21}$	=	$\left(\frac{1}{4} - x_6 - y_6 + z_6\right)\mathbf{a}_1 +$ $\left(\frac{1}{4} + x_6 + y_6 + z_6\right)\mathbf{a}_2 +$ $\left(\frac{1}{4} + x_6 - y_6 - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - y_6\right)b\,\hat{\mathbf{y}} +$ $\left(\frac{1}{4} + z_6\right)c\,\hat{\mathbf{z}}$	(16*b*)	V
$\mathbf{B}_{22}$	=	$\left(\frac{1}{4} + x_6 + y_6 + z_6\right)\mathbf{a}_1 +$ $\left(\frac{1}{4} - x_6 - y_6 + z_6\right)\mathbf{a}_2 +$ $\left(\frac{1}{4} - x_6 + y_6 - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + y_6\right)b\,\hat{\mathbf{y}} +$ $\left(\frac{1}{4} + z_6\right)c\,\hat{\mathbf{z}}$	(16*b*)	V

References:

- C. Calvo and R. Faggiani, *α Cupric Divanadate*, Acta Crystallogr. Sect. B Struct. Sci. **31**, 603–605 (1975), doi:10.1107/S0567740875003354.
- P. D. Robinson, J. M. Hughes, and M. L. Malinconico, *Blossite, α-$Cu_2^{2+}V_2^{5+}O_7$, an new fumarolic sublimate from Izalco volcano, El Salvador*, Am. Mineral. **72**, 397–400 (1987).

Found in:

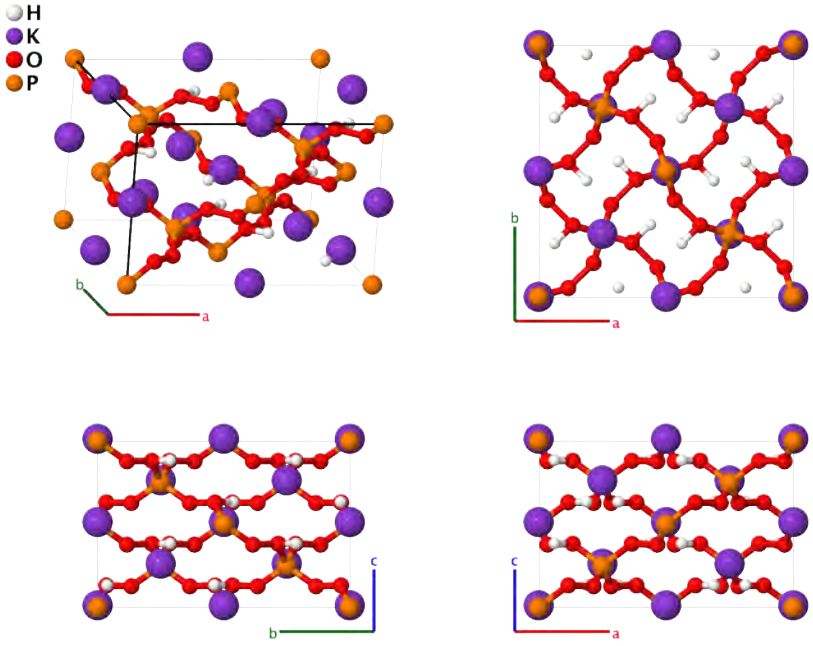
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1601
- POSCAR: pp. 1602

Archerite (KH_2PO_4) Structure: A2BC4D_oF64_43_b_a_2b_a

http://aflow.org/prototype-encyclopedia/A2BC4D_oF64_43_b_a_2b_a

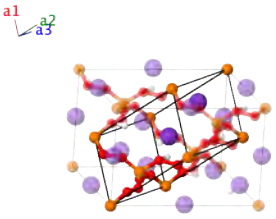


Prototype	:	H_2KO_4P
AFLOW prototype label	:	A2BC4D_oF64_43_b_a_2b_a
***Strukturbericht* designation**	:	None
Pearson symbol	:	oF64
Space group number	:	43
Space group symbol	:	$Fdd2$
AFLOW prototype command	:	`aflow --proto=A2BC4D_oF64_43_b_a_2b_a --params=`$a, b/a, c/a, z_1, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5$

- This structure is stable below 121 K, and the data was taken at 113 K. The high temperature structure is tetragonal with disordered hydrogen (Levy, 1954).
- The origin of the z-axis is not restricted in space group $Fdd2$ #43. Here it is fixed by putting the phosphorus atom at the origin.

Face-centered Orthorhombic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,b\,\hat{\mathbf{y}} + \tfrac{1}{2}\,c\,\hat{\mathbf{z}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,c\,\hat{\mathbf{z}}\\ \mathbf{a}_3 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,b\,\hat{\mathbf{y}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$z_1\,\mathbf{a}_1 + z_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$z_1 c\,\hat{\mathbf{z}}$	$(8a)$	K
$\mathbf{B}_2$	=	$\left(\frac{1}{4}+z_1\right)\mathbf{a}_1 + \left(\frac{1}{4}+z_1\right)\mathbf{a}_2 + \left(\frac{1}{4}-z_1\right)\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{4}+z_1\right)c\,\hat{\mathbf{z}}$	$(8a)$	K
$\mathbf{B}_3$	=	$z_2\,\mathbf{a}_1 + z_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$z_2 c\,\hat{\mathbf{z}}$	$(8a)$	P
$\mathbf{B}_4$	=	$\left(\frac{1}{4}+z_2\right)\mathbf{a}_1 + \left(\frac{1}{4}+z_2\right)\mathbf{a}_2 + \left(\frac{1}{4}-z_2\right)\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{4}+z_2\right)c\,\hat{\mathbf{z}}$	$(8a)$	P
$\mathbf{B}_5$	=	$(-x_3+y_3+z_3)\,\mathbf{a}_1 + (x_3-y_3+z_3)\,\mathbf{a}_2 + (x_3+y_3-z_3)\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(16b)$	H
$\mathbf{B}_6$	=	$(x_3-y_3+z_3)\,\mathbf{a}_1 + (-x_3+y_3+z_3)\,\mathbf{a}_2 + (-x_3-y_3-z_3)\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(16b)$	H
$\mathbf{B}_7$	=	$\left(\frac{1}{4}-x_3-y_3+z_3\right)\mathbf{a}_1 + \left(\frac{1}{4}+x_3+y_3+z_3\right)\mathbf{a}_2 + \left(\frac{1}{4}+x_3-y_3-z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}+x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4}-y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{4}+z_3\right)c\,\hat{\mathbf{z}}$	$(16b)$	H
$\mathbf{B}_8$	=	$\left(\frac{1}{4}+x_3+y_3+z_3\right)\mathbf{a}_1 + \left(\frac{1}{4}-x_3-y_3+z_3\right)\mathbf{a}_2 + \left(\frac{1}{4}-x_3+y_3-z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}-x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4}+y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{4}+z_3\right)c\,\hat{\mathbf{z}}$	$(16b)$	H
$\mathbf{B}_9$	=	$(-x_4+y_4+z_4)\,\mathbf{a}_1 + (x_4-y_4+z_4)\,\mathbf{a}_2 + (x_4+y_4-z_4)\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(16b)$	O I
$\mathbf{B}_{10}$	=	$(x_4-y_4+z_4)\,\mathbf{a}_1 + (-x_4+y_4+z_4)\,\mathbf{a}_2 + (-x_4-y_4-z_4)\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(16b)$	O I
$\mathbf{B}_{11}$	=	$\left(\frac{1}{4}-x_4-y_4+z_4\right)\mathbf{a}_1 + \left(\frac{1}{4}+x_4+y_4+z_4\right)\mathbf{a}_2 + \left(\frac{1}{4}+x_4-y_4-z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}+x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4}-y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{4}+z_4\right)c\,\hat{\mathbf{z}}$	$(16b)$	O I
$\mathbf{B}_{12}$	=	$\left(\frac{1}{4}+x_4+y_4+z_4\right)\mathbf{a}_1 + \left(\frac{1}{4}-x_4-y_4+z_4\right)\mathbf{a}_2 + \left(\frac{1}{4}-x_4+y_4-z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}-x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4}+y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{4}+z_4\right)c\,\hat{\mathbf{z}}$	$(16b)$	O I
$\mathbf{B}_{13}$	=	$(-x_5+y_5+z_5)\,\mathbf{a}_1 + (x_5-y_5+z_5)\,\mathbf{a}_2 + (x_5+y_5-z_5)\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(16b)$	O II

$\mathbf{B}_{14}$	=	$(x_5 - y_5 + z_5)\,\mathbf{a}_1 +$ $(-x_5 + y_5 + z_5)\,\mathbf{a}_2 +$ $(-x_5 - y_5 - z_5)\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(16*b*)	O II
$\mathbf{B}_{15}$	=	$\left(\frac{1}{4} - x_5 - y_5 + z_5\right)\mathbf{a}_1 +$ $\left(\frac{1}{4} + x_5 + y_5 + z_5\right)\mathbf{a}_2 +$ $\left(\frac{1}{4} + x_5 - y_5 - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - y_5\right)b\,\hat{\mathbf{y}} +$ $\left(\frac{1}{4} + z_5\right)c\,\hat{\mathbf{z}}$	(16*b*)	O II
$\mathbf{B}_{16}$	=	$\left(\frac{1}{4} + x_5 + y_5 + z_5\right)\mathbf{a}_1 +$ $\left(\frac{1}{4} - x_5 - y_5 + z_5\right)\mathbf{a}_2 +$ $\left(\frac{1}{4} - x_5 + y_5 - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + y_5\right)b\,\hat{\mathbf{y}} +$ $\left(\frac{1}{4} + z_5\right)c\,\hat{\mathbf{z}}$	(16*b*)	O II

References:

- H. A. Levy, S. W. Peterson, and S. H. Simonsen, *Neutron Diffraction Study of the Ferroelectric Modification of Potassium Dihydrogen Phosphate*, Phys. Rev. **93**, 1120–1121 (1954), doi:10.1103/PhysRev.93.1120.

Found in:

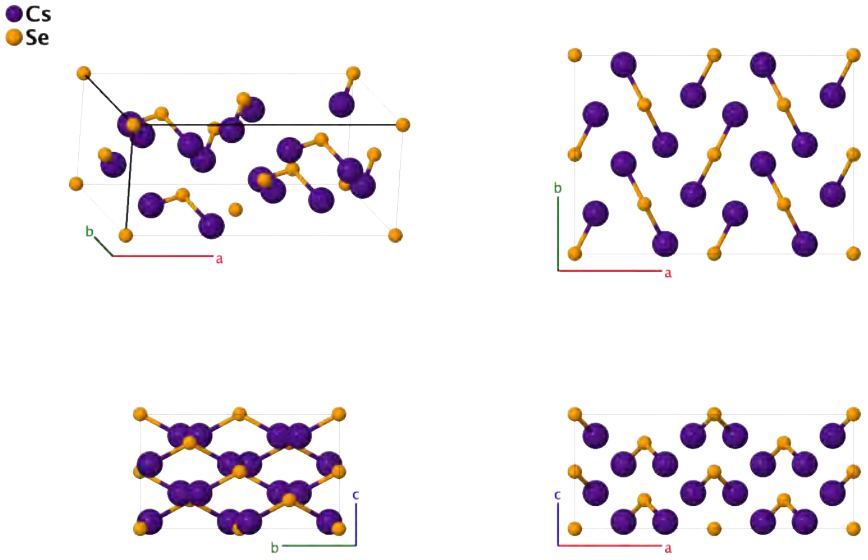
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1602
- POSCAR: pp. 1602

Cs_2Se Structure: A2B_oF24_43_b_a

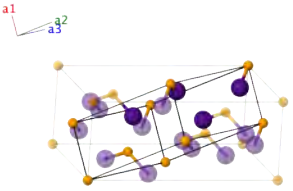
http://aflow.org/prototype-encyclopedia/A2B_oF24_43_b_a



Prototype	:	Cs_2Se
AFLOW prototype label	:	A2B_oF24_43_b_a
***Strukturbericht* designation**	:	None
Pearson symbol	:	oF24
Space group number	:	43
Space group symbol	:	$Fdd2$
AFLOW prototype command	:	`aflow --proto=A2B_oF24_43_b_a --params=`$a, b/a, c/a, z_1, x_2, y_2, z_2$

Face-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,b\,\hat{\mathbf{y}} + \frac{1}{2}\,c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{1}{2}\,c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{1}{2}\,b\,\hat{\mathbf{y}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$z_1\,\mathbf{a}_1 + z_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$z_1 c\,\hat{\mathbf{z}}$	$(8a)$	Se
$\mathbf{B}_2$	=	$\left(\frac{1}{4}+z_1\right)\mathbf{a}_1 + \left(\frac{1}{4}+z_1\right)\mathbf{a}_2 + \left(\frac{1}{4}-z_1\right)\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{4}+z_1\right)c\,\hat{\mathbf{z}}$	$(8a)$	Se
$\mathbf{B}_3$	=	$(-x_2+y_2+z_2)\,\mathbf{a}_1 + (x_2-y_2+z_2)\,\mathbf{a}_2 + (x_2+y_2-z_2)\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(16b)$	Cs

$\mathbf{B}_4$	=	$(x_2 - y_2 + z_2)\,\mathbf{a}_1 +$ $(-x_2 + y_2 + z_2)\,\mathbf{a}_2 +$ $(-x_2 - y_2 - z_2)\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(16*b*)	Cs
$\mathbf{B}_5$	=	$\left(\frac{1}{4} - x_2 - y_2 + z_2\right)\mathbf{a}_1 +$ $\left(\frac{1}{4} + x_2 + y_2 + z_2\right)\mathbf{a}_2 +$ $\left(\frac{1}{4} + x_2 - y_2 - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - y_2\right)b\,\hat{\mathbf{y}} +$ $\left(\frac{1}{4} + z_2\right)c\,\hat{\mathbf{z}}$	(16*b*)	Cs
$\mathbf{B}_6$	=	$\left(\frac{1}{4} + x_2 + y_2 + z_2\right)\mathbf{a}_1 +$ $\left(\frac{1}{4} - x_2 - y_2 + z_2\right)\mathbf{a}_2 +$ $\left(\frac{1}{4} - x_2 + y_2 - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + y_2\right)b\,\hat{\mathbf{y}} +$ $\left(\frac{1}{4} + z_2\right)c\,\hat{\mathbf{z}}$	(16*b*)	Cs

References:

- P. Böttcher, *Zur Kenntnis von Cs_2Se*, J. Less-Common Met. **76**, 271–277 (1980), doi:10.1016/0022-5088(80)90029-6.

Found in:

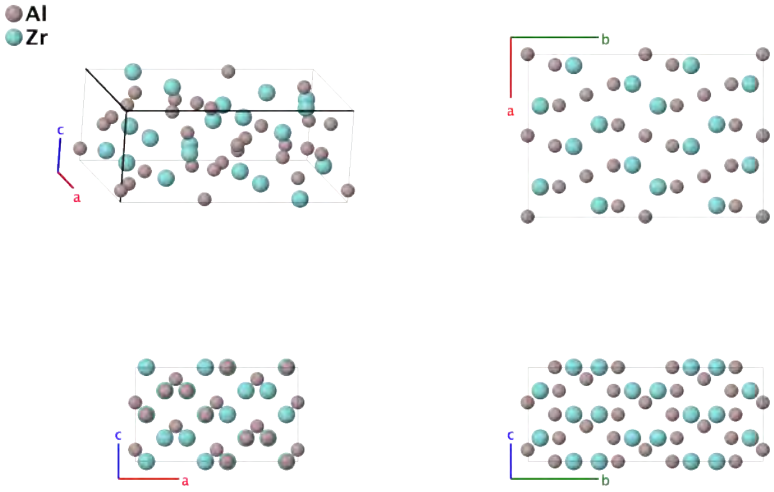
- P. Villars and L. D. Calvert, eds., *Pearson's Handbook of Crystallographic Data* (ASM International, Materials Park OH, 1991), vol. III, p. 2781.

Geometry files:

- CIF: pp. 1602
- POSCAR: pp. 1603

Zr_2Al_3 Structure: A3B2_oF40_43_ab_b

http://aflow.org/prototype-encyclopedia/A3B2_oF40_43_ab_b



Prototype	:	Al_3Zr_2
AFLOW prototype label	:	A3B2_oF40_43_ab_b
***Strukturbericht* designation**	:	None
Pearson symbol	:	oF40
Space group number	:	43
Space group symbol	:	$Fdd2$
AFLOW prototype command	:	`aflow --proto=A3B2_oF40_43_ab_b --params=`$a, b/a, c/a, z_1, x_2, y_2, z_2, x_3, y_3, z_3$

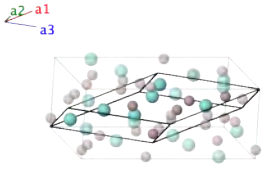
Other compounds with this structure

- Hf_2Al_3

- The $z = 0$ plane is not fixed in space group $Fdd2$ #43. Here it is arbitrarily chosen so that $z_3 = 0$ for the zirconium atom.

Face-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2} b\,\hat{\mathbf{y}} + \frac{1}{2} c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2} a\,\hat{\mathbf{x}} + \frac{1}{2} c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2} a\,\hat{\mathbf{x}} + \frac{1}{2} b\,\hat{\mathbf{y}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$z_1\,\mathbf{a}_1 + z_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$z_1 c\,\hat{\mathbf{z}}$	$(8a)$	Al I

$\mathbf{B}_2$	=	$\left(\frac{1}{4}+z_1\right)\mathbf{a}_1+\left(\frac{1}{4}+z_1\right)\mathbf{a}_2+\left(\frac{1}{4}-z_1\right)\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{4}+z_1\right)c\,\hat{\mathbf{z}}$	(8*a*)	Al I
$\mathbf{B}_3$	=	$(-x_2+y_2+z_2)\,\mathbf{a}_1+(x_2-y_2+z_2)\,\mathbf{a}_2+(x_2+y_2-z_2)\,\mathbf{a}_3$	=	$x_2a\,\hat{\mathbf{x}}+y_2b\,\hat{\mathbf{y}}+z_2c\,\hat{\mathbf{z}}$	(16*b*)	Al II
$\mathbf{B}_4$	=	$(x_2-y_2+z_2)\,\mathbf{a}_1+(-x_2+y_2+z_2)\,\mathbf{a}_2+(-x_2-y_2-z_2)\,\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}}-y_2b\,\hat{\mathbf{y}}+z_2c\,\hat{\mathbf{z}}$	(16*b*)	Al II
$\mathbf{B}_5$	=	$\left(\frac{1}{4}-x_2-y_2+z_2\right)\mathbf{a}_1+\left(\frac{1}{4}+x_2+y_2+z_2\right)\mathbf{a}_2+\left(\frac{1}{4}+x_2-y_2-z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}+x_2\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-y_2\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{4}+z_2\right)c\,\hat{\mathbf{z}}$	(16*b*)	Al II
$\mathbf{B}_6$	=	$\left(\frac{1}{4}+x_2+y_2+z_2\right)\mathbf{a}_1+\left(\frac{1}{4}-x_2-y_2+z_2\right)\mathbf{a}_2+\left(\frac{1}{4}-x_2+y_2-z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}-x_2\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+y_2\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{4}+z_2\right)c\,\hat{\mathbf{z}}$	(16*b*)	Al II
$\mathbf{B}_7$	=	$(-x_3+y_3+z_3)\,\mathbf{a}_1+(x_3-y_3+z_3)\,\mathbf{a}_2+(x_3+y_3-z_3)\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}}+y_3b\,\hat{\mathbf{y}}+z_3c\,\hat{\mathbf{z}}$	(16*b*)	Zr
$\mathbf{B}_8$	=	$(x_3-y_3+z_3)\,\mathbf{a}_1+(-x_3+y_3+z_3)\,\mathbf{a}_2+(-x_3-y_3-z_3)\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}}-y_3b\,\hat{\mathbf{y}}+z_3c\,\hat{\mathbf{z}}$	(16*b*)	Zr
$\mathbf{B}_9$	=	$\left(\frac{1}{4}-x_3-y_3+z_3\right)\mathbf{a}_1+\left(\frac{1}{4}+x_3+y_3+z_3\right)\mathbf{a}_2+\left(\frac{1}{4}+x_3-y_3-z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}+x_3\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-y_3\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{4}+z_3\right)c\,\hat{\mathbf{z}}$	(16*b*)	Zr
$\mathbf{B}_{10}$	=	$\left(\frac{1}{4}+x_3+y_3+z_3\right)\mathbf{a}_1+\left(\frac{1}{4}-x_3-y_3+z_3\right)\mathbf{a}_2+\left(\frac{1}{4}-x_3+y_3-z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}-x_3\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+y_3\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{4}+z_3\right)c\,\hat{\mathbf{z}}$	(16*b*)	Zr

References:

- T. J. Renouf and C. A. Beevers, *The Crystal Structure of* Zr_2Al_3, Acta Cryst. **14**, 469–472 (1961), doi:10.1107/S0365110X61001510.

Found in:

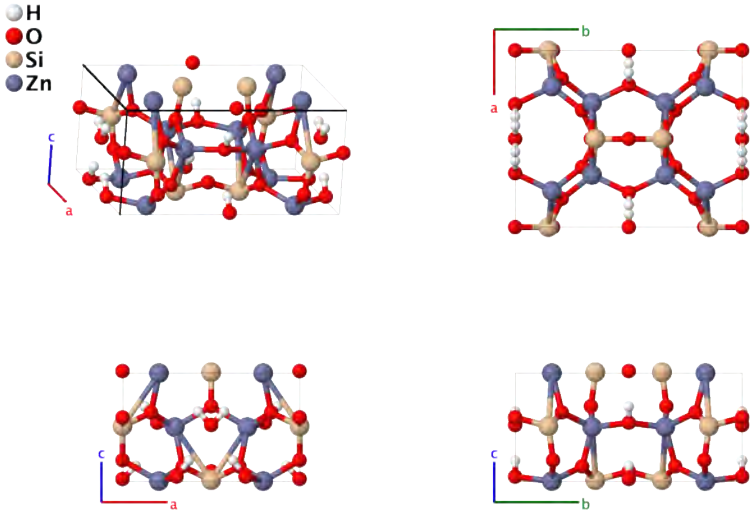
- L.-E. Edshammar, *Crystal Structure Investigations on the Zr-Al and Hf-Al Systems*, Acta Chem. Scand. **16**, 20–30 (1962), doi:10.3891/acta.chem.scand.16-0020.

Geometry files:

- CIF: pp. 1603
- POSCAR: pp. 1603

Hemimorphite ($Zn_4Si_2O_7(OH)_2{\cdot}H_2O$, $S\,2_2$) Structure: A2B5CD2_oI40_44_2c_abcde_d_e

http://aflow.org/prototype-encyclopedia/A2B5CD2_oI40_44_2c_abcde_d_e

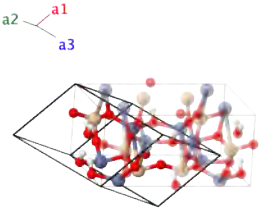


Prototype	:	$H_2O_5SiZn_2$
AFLOW prototype label	:	A2B5CD2_oI40_44_2c_abcde_d_e
***Strukturbericht* designation**	:	$S\,2_2$
Pearson symbol	:	oI40
Space group number	:	44
Space group symbol	:	*Imm*2
AFLOW prototype command	:	`aflow --proto=A2B5CD2_oI40_44_2c_abcde_d_e --params=`$a, b/a, c/a, z_1, z_2, x_3, z_3, x_4, z_4, x_5, z_5, y_6, z_6, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9$

- The original (Ito, 1932) determination of this structure did not locate the positions of the hydrogen atoms. (Hill, 1977) were able to do this, so we use the updated structure as the prototype.
- (Hill, 1977) gives the z coordinates of the atoms on the ($2c$) sites as 0.0190, 0.0643, and 0.0410, respectively, but this gives unrealistic H-O distances. Examination of the figures and distance tables shows that we should take $z_2 = 0.190$, $z_3 = 0.643$, and $z_4 = 0.041$, a conclusion also reached by (Downs, 2003).

Body-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = -\tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,b\,\hat{\mathbf{y}} + \tfrac{1}{2}\,c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2}\,b\,\hat{\mathbf{y}} + \tfrac{1}{2}\,c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,b\,\hat{\mathbf{y}} - \tfrac{1}{2}\,c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$z_1\,\mathbf{a}_1 + z_1\,\mathbf{a}_2$	=	$z_1 c\,\hat{\mathbf{z}}$	(2*a*)	O I
$\mathbf{B}_2$	=	$\left(\frac{1}{2} + z_2\right)\mathbf{a}_1 + z_2\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(2*b*)	O II
$\mathbf{B}_3$	=	$z_3\,\mathbf{a}_1 + (x_3 + z_3)\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + z_3 c\,\hat{\mathbf{z}}$	(4*c*)	H I
$\mathbf{B}_4$	=	$z_3\,\mathbf{a}_1 + (-x_3 + z_3)\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + z_3 c\,\hat{\mathbf{z}}$	(4*c*)	H I
$\mathbf{B}_5$	=	$z_4\,\mathbf{a}_1 + (x_4 + z_4)\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + z_4 c\,\hat{\mathbf{z}}$	(4*c*)	H II
$\mathbf{B}_6$	=	$z_4\,\mathbf{a}_1 + (-x_4 + z_4)\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + z_4 c\,\hat{\mathbf{z}}$	(4*c*)	H II
$\mathbf{B}_7$	=	$z_5\,\mathbf{a}_1 + (x_5 + z_5)\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + z_5 c\,\hat{\mathbf{z}}$	(4*c*)	O III
$\mathbf{B}_8$	=	$z_5\,\mathbf{a}_1 + (-x_5 + z_5)\,\mathbf{a}_2 - x_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + z_5 c\,\hat{\mathbf{z}}$	(4*c*)	O III
$\mathbf{B}_9$	=	$(y_6 + z_6)\,\mathbf{a}_1 + z_6\,\mathbf{a}_2 + y_6\,\mathbf{a}_3$	=	$y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(4*d*)	O IV
$\mathbf{B}_{10}$	=	$(-y_6 + z_6)\,\mathbf{a}_1 + z_6\,\mathbf{a}_2 - y_6\,\mathbf{a}_3$	=	$-y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(4*d*)	O IV
$\mathbf{B}_{11}$	=	$(y_7 + z_7)\,\mathbf{a}_1 + z_7\,\mathbf{a}_2 + y_7\,\mathbf{a}_3$	=	$y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(4*d*)	Si
$\mathbf{B}_{12}$	=	$(-y_7 + z_7)\,\mathbf{a}_1 + z_7\,\mathbf{a}_2 - y_7\,\mathbf{a}_3$	=	$-y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(4*d*)	Si
$\mathbf{B}_{13}$	=	$(y_8 + z_8)\,\mathbf{a}_1 + (x_8 + z_8)\,\mathbf{a}_2 + (x_8 + y_8)\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(8*e*)	O V
$\mathbf{B}_{14}$	=	$(-y_8 + z_8)\,\mathbf{a}_1 + (-x_8 + z_8)\,\mathbf{a}_2 + (-x_8 - y_8)\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(8*e*)	O V
$\mathbf{B}_{15}$	=	$(-y_8 + z_8)\,\mathbf{a}_1 + (x_8 + z_8)\,\mathbf{a}_2 + (x_8 - y_8)\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(8*e*)	O V
$\mathbf{B}_{16}$	=	$(y_8 + z_8)\,\mathbf{a}_1 + (-x_8 + z_8)\,\mathbf{a}_2 + (-x_8 + y_8)\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(8*e*)	O V
$\mathbf{B}_{17}$	=	$(y_9 + z_9)\,\mathbf{a}_1 + (x_9 + z_9)\,\mathbf{a}_2 + (x_9 + y_9)\,\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(8*e*)	Zn
$\mathbf{B}_{18}$	=	$(-y_9 + z_9)\,\mathbf{a}_1 + (-x_9 + z_9)\,\mathbf{a}_2 + (-x_9 - y_9)\,\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(8*e*)	Zn
$\mathbf{B}_{19}$	=	$(-y_9 + z_9)\,\mathbf{a}_1 + (x_9 + z_9)\,\mathbf{a}_2 + (x_9 - y_9)\,\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(8*e*)	Zn
$\mathbf{B}_{20}$	=	$(y_9 + z_9)\,\mathbf{a}_1 + (-x_9 + z_9)\,\mathbf{a}_2 + (-x_9 + y_9)\,\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(8*e*)	Zn

References:

- R. J. Hill, G. V. Gibbs, J. R. Craig, F. K. Ross, and J. M. Williams, *A neutron-diffraction study of hemimorphite*, Zeitschrift für Kristallographie - Crystalline Materials **146**, 241–259 (1977), doi:10.1524/zkri.1978.146.16.241.
- T. Ito and J. West, *The Structure of Hemimorphite ($H_2Zn_2SiO_5$)*, Zeitschrift für Kristallographie - Crystalline Materials **83**, 1–8 (1932), doi:10.1524/zkri.1932.83.1.1.

Found in:

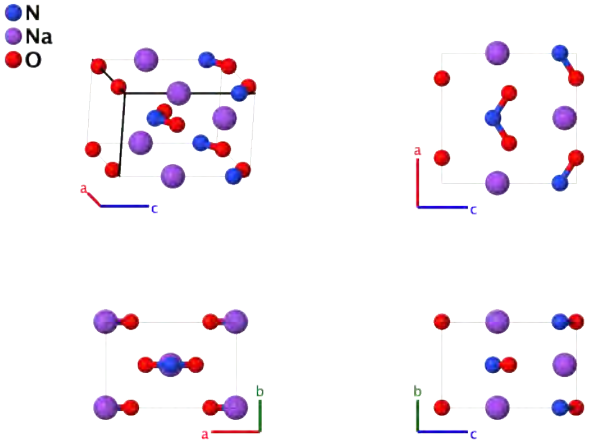
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1603
- POSCAR: pp. 1604

Ferroelectric $NaNO_2$ ($F5_5$) Structure: ABC2_oI8_44_a_a_c

http://aflow.org/prototype-encyclopedia/ABC2_oI8_44_a_a_c



Prototype	:	$NaNO_2$
AFLOW prototype label	:	ABC2_oI8_44_a_a_c
***Strukturbericht* designation**	:	$F5_5$
Pearson symbol	:	oI8
Space group number	:	44
Space group symbol	:	*Imm*2
AFLOW prototype command	:	`aflow --proto=ABC2_oI8_44_a_a_c --params=`$a, b/a, c/a, z_1, z_2, x_3, z_3$

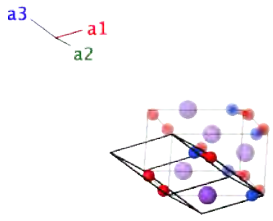
Other compounds with this structure

- $AgNO_2$ ($F5_{12}$)

- This is the low-temperature phase of $NaNO_2$. It is ferroelectric below 158°C.
- (Kay, 1961) gives the structure in the *Im*2*m* setting of space group #44. We have used FINDSYM to put it in the standard *Imm*2 setting.
- This structure is very similar to silver nitrite, $AgNO_2$, $F5_{12}$. We list the two structures as they are listed separately in the *Strukturbericht* volumes.
- The origin of the z coordinate in *Imm*2, or the y coordinate in *Im*2*m*, is arbitrary, and z_3 is set to zero.

Body-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = -\tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,b\,\hat{\mathbf{y}} + \tfrac{1}{2}\,c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2}\,b\,\hat{\mathbf{y}} + \tfrac{1}{2}\,c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,b\,\hat{\mathbf{y}} - \tfrac{1}{2}\,c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$z_1\,\mathbf{a}_1 + z_1\,\mathbf{a}_2$	=	$z_1 c\,\hat{\mathbf{z}}$	$(2a)$	N
$\mathbf{B}_2$	=	$z_2\,\mathbf{a}_1 + z_2\,\mathbf{a}_2$	=	$z_2 c\,\hat{\mathbf{z}}$	$(2a)$	Na
$\mathbf{B}_3$	=	$z_3\,\mathbf{a}_1 + (x_3 + z_3)\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + z_3 c\,\hat{\mathbf{z}}$	$(4c)$	O
$\mathbf{B}_4$	=	$z_3\,\mathbf{a}_1 + (-x_3 + z_3)\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + z_3 c\,\hat{\mathbf{z}}$	$(4c)$	O

References:

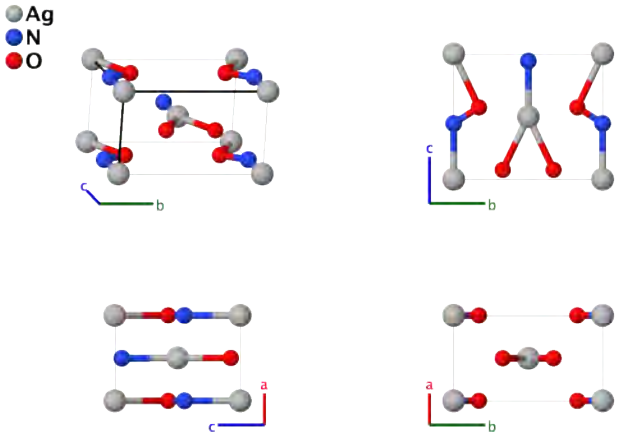
- M. I. Kay and B. C. Frazier, *A neutron diffraction refinement of the low temperature phase of $NaNO_2$*, Acta Cryst. **14**, 56–57 (1961), doi:10.1107/S0365110X61000103.

Geometry files:

- CIF: pp. 1604
- POSCAR: pp. 1604

$AgNO_2$ ($F5_{12}$) Structure: ABC2_oI8_44_a_a_d

http://aflow.org/prototype-encyclopedia/ABC2_oI8_44_a_a_d



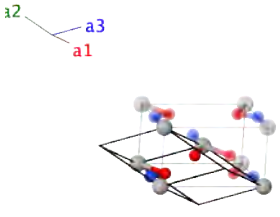
Prototype	:	$AgNO_2$
AFLOW prototype label	:	ABC2_oI8_44_a_a_d
***Strukturbericht* designation**	:	$F5_{12}$
Pearson symbol	:	oI8
Space group number	:	44
Space group symbol	:	*Imm*2
AFLOW prototype command	:	`aflow --proto=ABC2_oI8_44_a_a_d --params=`$a, b/a, c/a, z_1, z_2, y_3, z_3$

Other compounds with this structure

- $NaNO_2$ ($F5_5$)

- This structure is very similar to sodium nitrite, $NbNO_2$, $F5_5$. We list the two structures as they are listed separately in the *Strukturbericht* volumes.
- The origin of the z coordinate in *Imm*2 is arbitrary, and we set z_1 to zero.

Body-centered Orthorhombic primitive vectors:



$$\mathbf{a}_1 = -\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} - \frac{1}{2}c\,\hat{\mathbf{z}}$$

Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$z_1\,\mathbf{a}_1 + z_1\,\mathbf{a}_2$	=	$z_1 c\,\hat{\mathbf{z}}$	($2a$)	Ag

$\mathbf{B}_2$	=	$z_2\,\mathbf{a}_1 + z_2\,\mathbf{a}_2$	=	$z_2 c\,\hat{\mathbf{z}}$	($2a$)	N
$\mathbf{B}_3$	=	$(y_3 + z_3)\,\mathbf{a}_1 + z_3\,\mathbf{a}_2 + y_3\,\mathbf{a}_3$	=	$y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	($4d$)	O
$\mathbf{B}_4$	=	$(-y_3 + z_3)\,\mathbf{a}_1 + z_3\,\mathbf{a}_2 - y_3\,\mathbf{a}_3$	=	$-y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	($4d$)	O

References:

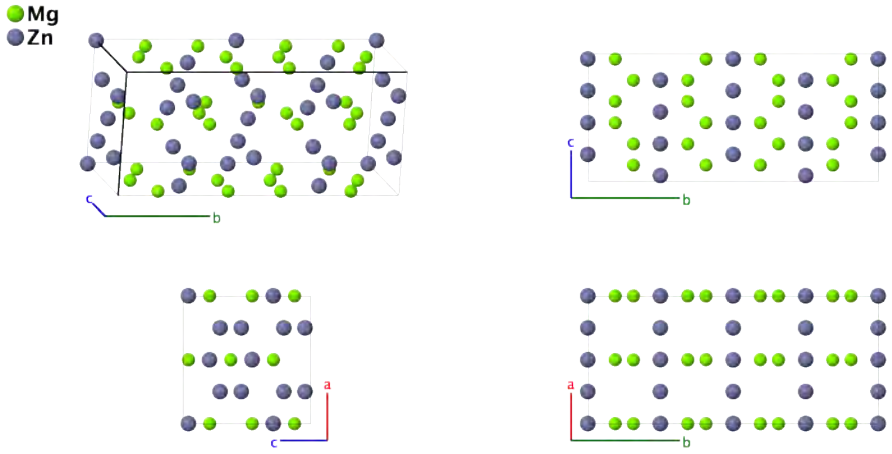
- S. Ohba and Y. Saito, *Structure of silver(I) nitrite, a redetermination*, Acta Crystallogr. Sect. B Struct. Sci. **37**, 1911–1913 (1981), doi:10.1107/S0567740881007565.

Geometry files:

- CIF: pp. 1604
- POSCAR: pp. 1605

*B*30 (MgZn?) Structure: AB_oI48_44_6d_ab2cde

http://aflow.org/prototype-encyclopedia/AB_oI48_44_6d_ab2cde

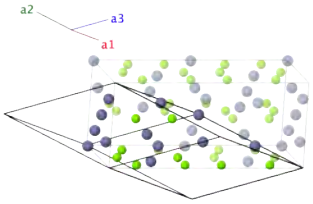


Prototype	:	MgZn
AFLOW prototype label	:	AB_oI48_44_6d_ab2cde
***Strukturbericht* designation**	:	*B*30
Pearson symbol	:	oI48
Space group number	:	44
Space group symbol	:	*Imm*2
AFLOW prototype command	:	`aflow --proto=AB_oI48_44_6d_ab2cde --params=`$a, b/a, c/a, z_1, z_2, x_3, z_3, x_4, z_4, y_5, z_5, y_6, z_6, y_7, z_7, y_8, z_8, y_9, z_9, y_{10}, z_{10}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}$

- It is rather a mystery why (Hermann, 1937) gave this the *Strukturbericht* designation *B*30, as the structure presented in the literature contradicts itself. (Tarschish, 1933) derived this structure from the hexagonal Laves structure $MgZn_2$ (*C*14) by doubling the unit cell in all directions to obtain a 96 atom unit cell, replacing 16 of the zinc atoms in this structure by magnesium, and shifting the z-coordinates of these atoms by $\pm c/16$. He then states that the space group remains $P6_3/mmc$ #194. (McKeehan, 1935) pointed out that this is impossible, as the converted Mg atoms only have a two-fold rotation axis about the z-axis. He assigned the structure to space group *Pmm*2 #25. (Hermann, 1937) referenced both papers, giving the space group as $P6_3/mmc$ but listing the atomic coordinates enumerated by McKeehan. In fact, the McKeehan structure has space group *Imm*2 #44, with 48 atoms in the conventional cell, half of the original, and 24 atoms in the primitive cell. This was noted, without reference, by (Parthé, 1993), which is the only comprehensive list of *Strukturbericht* symbols to include the *B*30 structure. We have reproduced this *Imm*2 structure from McKeehan's data. The true structure of MgZn is unclear, although it is seen in the Mg-Zn binary phase diagram (Massalski, 1990) over a small range of compositions, a complete crystallographic study has never been published. It is possible that the actual structure is off-stoichiometry. There is some evidence of a $Mg_{12}Zn_{13}$ structure (Mezbahul-Islam, 2014), and $Mg_{21}Zn_{25}$ has been determined (Cerný, 2002) to have the $Zr_{21}Re_{25}$ structure. There are similar problems with the $D2_2$ $MgZn_5$ structure, which we discuss on that page.

Body-centered Orthorhombic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= -\tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,b\,\hat{\mathbf{y}} + \tfrac{1}{2}\,c\,\hat{\mathbf{z}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2}\,b\,\hat{\mathbf{y}} + \tfrac{1}{2}\,c\,\hat{\mathbf{z}}\\ \mathbf{a}_3 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,b\,\hat{\mathbf{y}} - \tfrac{1}{2}\,c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$z_1\,\mathbf{a}_1 + z_1\,\mathbf{a}_2$	=	$z_1 c\,\hat{\mathbf{z}}$	$(2a)$	Zn I
$\mathbf{B}_2$	=	$\left(\tfrac{1}{2} + z_2\right)\mathbf{a}_1 + z_2\,\mathbf{a}_2 + \tfrac{1}{2}\,\mathbf{a}_3$	=	$\tfrac{1}{2}b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(2b)$	Zn II
$\mathbf{B}_3$	=	$z_3\,\mathbf{a}_1 + (x_3 + z_3)\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + z_3 c\,\hat{\mathbf{z}}$	$(4c)$	Zn III
$\mathbf{B}_4$	=	$z_3\,\mathbf{a}_1 + (-x_3 + z_3)\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + z_3 c\,\hat{\mathbf{z}}$	$(4c)$	Zn III
$\mathbf{B}_5$	=	$z_4\,\mathbf{a}_1 + (x_4 + z_4)\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + z_4 c\,\hat{\mathbf{z}}$	$(4c)$	Zn IV
$\mathbf{B}_6$	=	$z_4\,\mathbf{a}_1 + (-x_4 + z_4)\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + z_4 c\,\hat{\mathbf{z}}$	$(4c)$	Zn IV
$\mathbf{B}_7$	=	$(y_5 + z_5)\,\mathbf{a}_1 + z_5\,\mathbf{a}_2 + y_5\,\mathbf{a}_3$	=	$y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(4d)$	Mg I
$\mathbf{B}_8$	=	$(-y_5 + z_5)\,\mathbf{a}_1 + z_5\,\mathbf{a}_2 - y_5\,\mathbf{a}_3$	=	$-y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(4d)$	Mg I
$\mathbf{B}_9$	=	$(y_6 + z_6)\,\mathbf{a}_1 + z_6\,\mathbf{a}_2 + y_6\,\mathbf{a}_3$	=	$y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(4d)$	Mg II
$\mathbf{B}_{10}$	=	$(-y_6 + z_6)\,\mathbf{a}_1 + z_6\,\mathbf{a}_2 - y_6\,\mathbf{a}_3$	=	$-y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(4d)$	Mg II
$\mathbf{B}_{11}$	=	$(y_7 + z_7)\,\mathbf{a}_1 + z_7\,\mathbf{a}_2 + y_7\,\mathbf{a}_3$	=	$y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	$(4d)$	Mg III
$\mathbf{B}_{12}$	=	$(-y_7 + z_7)\,\mathbf{a}_1 + z_7\,\mathbf{a}_2 - y_7\,\mathbf{a}_3$	=	$-y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	$(4d)$	Mg III
$\mathbf{B}_{13}$	=	$(y_8 + z_8)\,\mathbf{a}_1 + z_8\,\mathbf{a}_2 + y_8\,\mathbf{a}_3$	=	$y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	$(4d)$	Mg IV
$\mathbf{B}_{14}$	=	$(-y_8 + z_8)\,\mathbf{a}_1 + z_8\,\mathbf{a}_2 - y_8\,\mathbf{a}_3$	=	$-y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	$(4d)$	Mg IV
$\mathbf{B}_{15}$	=	$(y_9 + z_9)\,\mathbf{a}_1 + z_9\,\mathbf{a}_2 + y_9\,\mathbf{a}_3$	=	$y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	$(4d)$	Mg V
$\mathbf{B}_{16}$	=	$(-y_9 + z_9)\,\mathbf{a}_1 + z_9\,\mathbf{a}_2 - y_9\,\mathbf{a}_3$	=	$-y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	$(4d)$	Mg V
$\mathbf{B}_{17}$	=	$(y_{10} + z_{10})\,\mathbf{a}_1 + z_{10}\,\mathbf{a}_2 + y_{10}\,\mathbf{a}_3$	=	$y_{10} b\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}}$	$(4d)$	Mg VI
$\mathbf{B}_{18}$	=	$(-y_{10} + z_{10})\,\mathbf{a}_1 + z_{10}\,\mathbf{a}_2 - y_{10}\,\mathbf{a}_3$	=	$-y_{10} b\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}}$	$(4d)$	Mg VI
$\mathbf{B}_{19}$	=	$(y_{11} + z_{11})\,\mathbf{a}_1 + z_{11}\,\mathbf{a}_2 + y_{11}\,\mathbf{a}_3$	=	$y_{11} b\,\hat{\mathbf{y}} + z_{11} c\,\hat{\mathbf{z}}$	$(4d)$	Zn V
$\mathbf{B}_{20}$	=	$(-y_{11} + z_{11})\,\mathbf{a}_1 + z_{11}\,\mathbf{a}_2 - y_{11}\,\mathbf{a}_3$	=	$-y_{11} b\,\hat{\mathbf{y}} + z_{11} c\,\hat{\mathbf{z}}$	$(4d)$	Zn V
$\mathbf{B}_{21}$	=	$(y_{12} + z_{12})\,\mathbf{a}_1 + (x_{12} + z_{12})\,\mathbf{a}_2 + (x_{12} + y_{12})\,\mathbf{a}_3$	=	$x_{12} a\,\hat{\mathbf{x}} + y_{12} b\,\hat{\mathbf{y}} + z_{12} c\,\hat{\mathbf{z}}$	$(8e)$	Zn VI
$\mathbf{B}_{22}$	=	$(-y_{12} + z_{12})\,\mathbf{a}_1 + (-x_{12} + z_{12})\,\mathbf{a}_2 + (-x_{12} - y_{12})\,\mathbf{a}_3$	=	$-x_{12} a\,\hat{\mathbf{x}} - y_{12} b\,\hat{\mathbf{y}} + z_{12} c\,\hat{\mathbf{z}}$	$(8e)$	Zn VI
$\mathbf{B}_{23}$	=	$(-y_{12} + z_{12})\,\mathbf{a}_1 + (x_{12} + z_{12})\,\mathbf{a}_2 + (x_{12} - y_{12})\,\mathbf{a}_3$	=	$x_{12} a\,\hat{\mathbf{x}} - y_{12} b\,\hat{\mathbf{y}} + z_{12} c\,\hat{\mathbf{z}}$	$(8e)$	Zn VI
$\mathbf{B}_{24}$	=	$(y_{12} + z_{12})\,\mathbf{a}_1 + (-x_{12} + z_{12})\,\mathbf{a}_2 + (-x_{12} + y_{12})\,\mathbf{a}_3$	=	$-x_{12} a\,\hat{\mathbf{x}} + y_{12} b\,\hat{\mathbf{y}} + z_{12} c\,\hat{\mathbf{z}}$	$(8e)$	Zn VI

References:

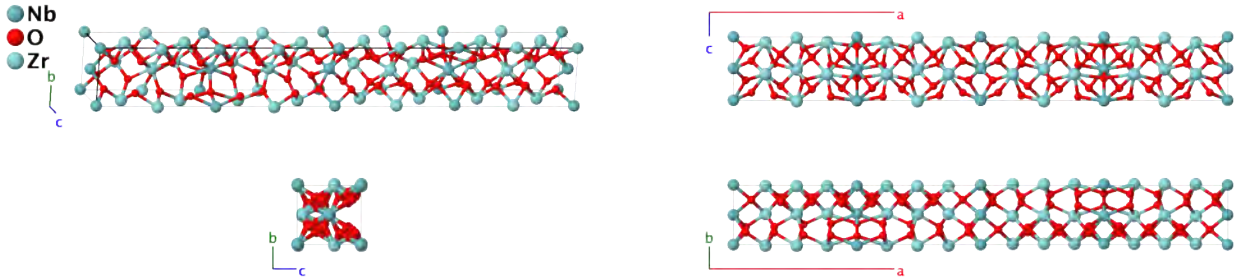
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

- L. Tarschisch, *Röntgenographische Untersuchung der Verbindungen MgZn und $MgZn_5$*, Zeitschrift für Kristallographie - Crystalline Materials **86**, 423–438 (1933), doi:10.1524/zkri.1933.86.1.423.
- L. W. McKeehan, *Note on MgZn and $MgZn_5$*, Zeitschrift für Kristallographie - Crystalline Materials **91**, 501–503 (1935), doi:10.1524/zkri.1935.91.1.501.
- E. Parthé, L. Gelato, B. Chabot, M. Penso, K. Cenzual, and R. Gladyshevskii, in *Standardized Data and Crystal Chemical Characterization of Inorganic Structure Types* (Springer-Verlag, Berlin, Heidelberg, 1993), *Gmelin Handbook of Inorganic and Organometallic Chemistry*, vol. 2, chap. Crystal Chemical Characterization of Inorganic Structure Types, 8 edn., doi:10.1007/978-3-662-02909-1_3.
- T. B. Massalski, H. Okamoto, P. R. Subramanian, and L. Kacprzak, eds., *Binary Alloy Phase Diagrams*, vol. 3 (ASM International, Materials Park, Ohio, USA, 1990), 2^{nd} edn. Hf-Re to Zn-Zr.
- M. Mezbahul-Islam, A. O. Mostafa, and M. Medraj, *Essential Magnesium Alloys Binary Phase Diagrams and Their Thermochemical Data*, J. Mater. **2014**, 704283 (2014), doi:10.1155/2014/704283.
- R. Cerný and G. Renaudin, *The intermetallic compound $Mg_{21}Zn_{25}$*, Acta Crystallogr. C **58**, i154–i155 (2002), doi:10.1107/S0108270102018103.

Geometry files:
- CIF: pp. 1605
- POSCAR: pp. 1605

$Nb_2Zr_6O_{17}$ Structure: A2B17C6_oI100_46_ab_b8c_3c

http://aflow.org/prototype-encyclopedia/A2B17C6_oI100_46_ab_b8c_3c



Prototype	:	$Nb_2O_{17}Zr_6$
AFLOW prototype label	:	A2B17C6_oI100_46_ab_b8c_3c
***Strukturbericht* designation**	:	None
Pearson symbol	:	oI100
Space group number	:	46
Space group symbol	:	*Ima*2
AFLOW prototype command	:	`aflow --proto=A2B17C6_oI100_46_ab_b8c_3c --params=`$a, b/a, c/a, z_1, y_2, z_2, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}$

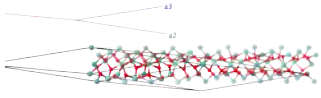
Other compounds with this structure

- $Nb_2Hf_6O_{17}$, $Nb_2Zr_6O_{17}$, $Ta_2Hf_6O_{17}$, and $Ta_2Zr_6O_{17}$

- Both (Galy, 1973) and (McCormack, 2019) state that the metallic atom sites are disordered, that is, for the prototype each metallic site has the average composition, $Nb_{0.25}Zr_{0.75}$. We place the niobium atoms on the (4*a*) and (4*b*) sites, and the zirconium on the (8*c*) sites so that the different symmetries are displayed.
- (McCormack, 2019) notes that the metallic composition of these compounds can deviate from the stoichiometry shown here.

Body-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = -\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}b\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}a\,\hat{\mathbf{x}} - \tfrac{1}{2}b\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}b\,\hat{\mathbf{y}} - \tfrac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$z_1\,\mathbf{a}_1 + z_1\,\mathbf{a}_2$	=	$z_1 c\,\hat{\mathbf{z}}$	(4*a*)	Nb I
$\mathbf{B}_2$	=	$z_1\,\mathbf{a}_1 + \left(\frac{1}{2}+z_1\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + z_1 c\,\hat{\mathbf{z}}$	(4*a*)	Nb I
$\mathbf{B}_3$	=	$(y_2+z_2)\,\mathbf{a}_1 + \left(\frac{1}{4}+z_2\right)\mathbf{a}_2 + \left(\frac{1}{4}+y_2\right)\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*b*)	Nb II
$\mathbf{B}_4$	=	$(-y_2+z_2)\,\mathbf{a}_1 + \left(\frac{3}{4}+z_2\right)\mathbf{a}_2 + \left(\frac{3}{4}-y_2\right)\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*b*)	Nb II
$\mathbf{B}_5$	=	$(y_3+z_3)\,\mathbf{a}_1 + \left(\frac{1}{4}+z_3\right)\mathbf{a}_2 + \left(\frac{1}{4}+y_3\right)\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*b*)	O I
$\mathbf{B}_6$	=	$(-y_3+z_3)\,\mathbf{a}_1 + \left(\frac{3}{4}+z_3\right)\mathbf{a}_2 + \left(\frac{3}{4}-y_3\right)\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*b*)	O I

$$\mathbf{B}_7 = (y_4+z_4)\,\mathbf{a}_1+(x_4+z_4)\,\mathbf{a}_2+(x_4+y_4)\,\mathbf{a}_3 = x_4a\,\hat{\mathbf{x}}+y_4b\,\hat{\mathbf{y}}+z_4c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O II}$$

$$\mathbf{B}_8 = (-y_4+z_4)\,\mathbf{a}_1+(-x_4+z_4)\,\mathbf{a}_2+(-x_4-y_4)\,\mathbf{a}_3 = -x_4a\,\hat{\mathbf{x}}-y_4b\,\hat{\mathbf{y}}+z_4c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O II}$$

$$\mathbf{B}_9 = (-y_4+z_4)\,\mathbf{a}_1+\left(\tfrac{1}{2}+x_4+z_4\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_4-y_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_4\right)a\,\hat{\mathbf{x}}-y_4b\,\hat{\mathbf{y}}+z_4c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O II}$$

$$\mathbf{B}_{10} = (y_4+z_4)\,\mathbf{a}_1+\left(\tfrac{1}{2}-x_4+z_4\right)\mathbf{a}_2+\left(\tfrac{1}{2}-x_4+y_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_4\right)a\,\hat{\mathbf{x}}+y_4b\,\hat{\mathbf{y}}+z_4c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O II}$$

$$\mathbf{B}_{11} = (y_5+z_5)\,\mathbf{a}_1+(x_5+z_5)\,\mathbf{a}_2+(x_5+y_5)\,\mathbf{a}_3 = x_5a\,\hat{\mathbf{x}}+y_5b\,\hat{\mathbf{y}}+z_5c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O III}$$

$$\mathbf{B}_{12} = (-y_5+z_5)\,\mathbf{a}_1+(-x_5+z_5)\,\mathbf{a}_2+(-x_5-y_5)\,\mathbf{a}_3 = -x_5a\,\hat{\mathbf{x}}-y_5b\,\hat{\mathbf{y}}+z_5c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O III}$$

$$\mathbf{B}_{13} = (-y_5+z_5)\,\mathbf{a}_1+\left(\tfrac{1}{2}+x_5+z_5\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_5-y_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_5\right)a\,\hat{\mathbf{x}}-y_5b\,\hat{\mathbf{y}}+z_5c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O III}$$

$$\mathbf{B}_{14} = (y_5+z_5)\,\mathbf{a}_1+\left(\tfrac{1}{2}-x_5+z_5\right)\mathbf{a}_2+\left(\tfrac{1}{2}-x_5+y_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_5\right)a\,\hat{\mathbf{x}}+y_5b\,\hat{\mathbf{y}}+z_5c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O III}$$

$$\mathbf{B}_{15} = (y_6+z_6)\,\mathbf{a}_1+(x_6+z_6)\,\mathbf{a}_2+(x_6+y_6)\,\mathbf{a}_3 = x_6a\,\hat{\mathbf{x}}+y_6b\,\hat{\mathbf{y}}+z_6c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O IV}$$

$$\mathbf{B}_{16} = (-y_6+z_6)\,\mathbf{a}_1+(-x_6+z_6)\,\mathbf{a}_2+(-x_6-y_6)\,\mathbf{a}_3 = -x_6a\,\hat{\mathbf{x}}-y_6b\,\hat{\mathbf{y}}+z_6c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O IV}$$

$$\mathbf{B}_{17} = (-y_6+z_6)\,\mathbf{a}_1+\left(\tfrac{1}{2}+x_6+z_6\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_6-y_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{x}}-y_6b\,\hat{\mathbf{y}}+z_6c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O IV}$$

$$\mathbf{B}_{18} = (y_6+z_6)\,\mathbf{a}_1+\left(\tfrac{1}{2}-x_6+z_6\right)\mathbf{a}_2+\left(\tfrac{1}{2}-x_6+y_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{x}}+y_6b\,\hat{\mathbf{y}}+z_6c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O IV}$$

$$\mathbf{B}_{19} = (y_7+z_7)\,\mathbf{a}_1+(x_7+z_7)\,\mathbf{a}_2+(x_7+y_7)\,\mathbf{a}_3 = x_7a\,\hat{\mathbf{x}}+y_7b\,\hat{\mathbf{y}}+z_7c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O V}$$

$$\mathbf{B}_{20} = (-y_7+z_7)\,\mathbf{a}_1+(-x_7+z_7)\,\mathbf{a}_2+(-x_7-y_7)\,\mathbf{a}_3 = -x_7a\,\hat{\mathbf{x}}-y_7b\,\hat{\mathbf{y}}+z_7c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O V}$$

$$\mathbf{B}_{21} = (-y_7+z_7)\,\mathbf{a}_1+\left(\tfrac{1}{2}+x_7+z_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_7-y_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{x}}-y_7b\,\hat{\mathbf{y}}+z_7c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O V}$$

$$\mathbf{B}_{22} = (y_7+z_7)\,\mathbf{a}_1+\left(\tfrac{1}{2}-x_7+z_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}-x_7+y_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{x}}+y_7b\,\hat{\mathbf{y}}+z_7c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O V}$$

$$\mathbf{B}_{23} = (y_8+z_8)\,\mathbf{a}_1+(x_8+z_8)\,\mathbf{a}_2+(x_8+y_8)\,\mathbf{a}_3 = x_8a\,\hat{\mathbf{x}}+y_8b\,\hat{\mathbf{y}}+z_8c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O VI}$$

$$\mathbf{B}_{24} = (-y_8+z_8)\,\mathbf{a}_1+(-x_8+z_8)\,\mathbf{a}_2+(-x_8-y_8)\,\mathbf{a}_3 = -x_8a\,\hat{\mathbf{x}}-y_8b\,\hat{\mathbf{y}}+z_8c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O VI}$$

$$\mathbf{B}_{25} = (-y_8+z_8)\,\mathbf{a}_1+\left(\tfrac{1}{2}+x_8+z_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_8-y_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{x}}-y_8b\,\hat{\mathbf{y}}+z_8c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O VI}$$

$$\mathbf{B}_{26} = (y_8+z_8)\,\mathbf{a}_1+\left(\tfrac{1}{2}-x_8+z_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}-x_8+y_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{x}}+y_8b\,\hat{\mathbf{y}}+z_8c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O VI}$$

$$\mathbf{B}_{27} = (y_9+z_9)\,\mathbf{a}_1+(x_9+z_9)\,\mathbf{a}_2+(x_9+y_9)\,\mathbf{a}_3 = x_9a\,\hat{\mathbf{x}}+y_9b\,\hat{\mathbf{y}}+z_9c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O VII}$$

$$\mathbf{B}_{28} = (-y_9+z_9)\,\mathbf{a}_1+(-x_9+z_9)\,\mathbf{a}_2+(-x_9-y_9)\,\mathbf{a}_3 = -x_9a\,\hat{\mathbf{x}}-y_9b\,\hat{\mathbf{y}}+z_9c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O VII}$$

$$\mathbf{B}_{29} = (-y_9+z_9)\,\mathbf{a}_1+\left(\tfrac{1}{2}+x_9+z_9\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_9-y_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_9\right)a\,\hat{\mathbf{x}}-y_9b\,\hat{\mathbf{y}}+z_9c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O VII}$$

$$\mathbf{B}_{30} = (y_9+z_9)\,\mathbf{a}_1+\left(\tfrac{1}{2}-x_9+z_9\right)\mathbf{a}_2+\left(\tfrac{1}{2}-x_9+y_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_9\right)a\,\hat{\mathbf{x}}+y_9b\,\hat{\mathbf{y}}+z_9c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O VII}$$

$$\mathbf{B}_{31} = (y_{10}+z_{10})\,\mathbf{a}_1 + (x_{10}+z_{10})\,\mathbf{a}_2 + (x_{10}+y_{10})\,\mathbf{a}_3 = x_{10}a\,\hat{\mathbf{x}} + y_{10}b\,\hat{\mathbf{y}} + z_{10}c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O VIII}$$

$$\mathbf{B}_{32} = (-y_{10}+z_{10})\,\mathbf{a}_1 + (-x_{10}+z_{10})\,\mathbf{a}_2 + (-x_{10}-y_{10})\,\mathbf{a}_3 = -x_{10}a\,\hat{\mathbf{x}} - y_{10}b\,\hat{\mathbf{y}} + z_{10}c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O VIII}$$

$$\mathbf{B}_{33} = (-y_{10}+z_{10})\,\mathbf{a}_1 + \left(\tfrac{1}{2}+x_{10}+z_{10}\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+x_{10}-y_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{10}\right)a\,\hat{\mathbf{x}} - y_{10}b\,\hat{\mathbf{y}} + z_{10}c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O VIII}$$

$$\mathbf{B}_{34} = (y_{10}+z_{10})\,\mathbf{a}_1 + \left(\tfrac{1}{2}-x_{10}+z_{10}\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-x_{10}+y_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{10}\right)a\,\hat{\mathbf{x}} + y_{10}b\,\hat{\mathbf{y}} + z_{10}c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O VIII}$$

$$\mathbf{B}_{35} = (y_{11}+z_{11})\,\mathbf{a}_1 + (x_{11}+z_{11})\,\mathbf{a}_2 + (x_{11}+y_{11})\,\mathbf{a}_3 = x_{11}a\,\hat{\mathbf{x}} + y_{11}b\,\hat{\mathbf{y}} + z_{11}c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O IX}$$

$$\mathbf{B}_{36} = (-y_{11}+z_{11})\,\mathbf{a}_1 + (-x_{11}+z_{11})\,\mathbf{a}_2 + (-x_{11}-y_{11})\,\mathbf{a}_3 = -x_{11}a\,\hat{\mathbf{x}} - y_{11}b\,\hat{\mathbf{y}} + z_{11}c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O IX}$$

$$\mathbf{B}_{37} = (-y_{11}+z_{11})\,\mathbf{a}_1 + \left(\tfrac{1}{2}+x_{11}+z_{11}\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+x_{11}-y_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{11}\right)a\,\hat{\mathbf{x}} - y_{11}b\,\hat{\mathbf{y}} + z_{11}c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O IX}$$

$$\mathbf{B}_{38} = (y_{11}+z_{11})\,\mathbf{a}_1 + \left(\tfrac{1}{2}-x_{11}+z_{11}\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-x_{11}+y_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{11}\right)a\,\hat{\mathbf{x}} + y_{11}b\,\hat{\mathbf{y}} + z_{11}c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O IX}$$

$$\mathbf{B}_{39} = (y_{12}+z_{12})\,\mathbf{a}_1 + (x_{12}+z_{12})\,\mathbf{a}_2 + (x_{12}+y_{12})\,\mathbf{a}_3 = x_{12}a\,\hat{\mathbf{x}} + y_{12}b\,\hat{\mathbf{y}} + z_{12}c\,\hat{\mathbf{z}} \quad (8c) \quad \text{Zr I}$$

$$\mathbf{B}_{40} = (-y_{12}+z_{12})\,\mathbf{a}_1 + (-x_{12}+z_{12})\,\mathbf{a}_2 + (-x_{12}-y_{12})\,\mathbf{a}_3 = -x_{12}a\,\hat{\mathbf{x}} - y_{12}b\,\hat{\mathbf{y}} + z_{12}c\,\hat{\mathbf{z}} \quad (8c) \quad \text{Zr I}$$

$$\mathbf{B}_{41} = (-y_{12}+z_{12})\,\mathbf{a}_1 + \left(\tfrac{1}{2}+x_{12}+z_{12}\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+x_{12}-y_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{12}\right)a\,\hat{\mathbf{x}} - y_{12}b\,\hat{\mathbf{y}} + z_{12}c\,\hat{\mathbf{z}} \quad (8c) \quad \text{Zr I}$$

$$\mathbf{B}_{42} = (y_{12}+z_{12})\,\mathbf{a}_1 + \left(\tfrac{1}{2}-x_{12}+z_{12}\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-x_{12}+y_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{12}\right)a\,\hat{\mathbf{x}} + y_{12}b\,\hat{\mathbf{y}} + z_{12}c\,\hat{\mathbf{z}} \quad (8c) \quad \text{Zr I}$$

$$\mathbf{B}_{43} = (y_{13}+z_{13})\,\mathbf{a}_1 + (x_{13}+z_{13})\,\mathbf{a}_2 + (x_{13}+y_{13})\,\mathbf{a}_3 = x_{13}a\,\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} + z_{13}c\,\hat{\mathbf{z}} \quad (8c) \quad \text{Zr II}$$

$$\mathbf{B}_{44} = (-y_{13}+z_{13})\,\mathbf{a}_1 + (-x_{13}+z_{13})\,\mathbf{a}_2 + (-x_{13}-y_{13})\,\mathbf{a}_3 = -x_{13}a\,\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} + z_{13}c\,\hat{\mathbf{z}} \quad (8c) \quad \text{Zr II}$$

$$\mathbf{B}_{45} = (-y_{13}+z_{13})\,\mathbf{a}_1 + \left(\tfrac{1}{2}+x_{13}+z_{13}\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+x_{13}-y_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{13}\right)a\,\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} + z_{13}c\,\hat{\mathbf{z}} \quad (8c) \quad \text{Zr II}$$

$$\mathbf{B}_{46} = (y_{13}+z_{13})\,\mathbf{a}_1 + \left(\tfrac{1}{2}-x_{13}+z_{13}\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-x_{13}+y_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{13}\right)a\,\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} + z_{13}c\,\hat{\mathbf{z}} \quad (8c) \quad \text{Zr II}$$

$$\mathbf{B}_{47} = (y_{14}+z_{14})\,\mathbf{a}_1 + (x_{14}+z_{14})\,\mathbf{a}_2 + (x_{14}+y_{14})\,\mathbf{a}_3 = x_{14}a\,\hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} + z_{14}c\,\hat{\mathbf{z}} \quad (8c) \quad \text{Zr III}$$

$$\mathbf{B}_{48} = (-y_{14}+z_{14})\,\mathbf{a}_1 + (-x_{14}+z_{14})\,\mathbf{a}_2 + (-x_{14}-y_{14})\,\mathbf{a}_3 = -x_{14}a\,\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} + z_{14}c\,\hat{\mathbf{z}} \quad (8c) \quad \text{Zr III}$$

$$\mathbf{B}_{49} = (-y_{14}+z_{14})\,\mathbf{a}_1 + \left(\tfrac{1}{2}+x_{14}+z_{14}\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+x_{14}-y_{14}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{14}\right)a\,\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} + z_{14}c\,\hat{\mathbf{z}} \quad (8c) \quad \text{Zr III}$$

$$\mathbf{B}_{50} = (y_{14}+z_{14})\,\mathbf{a}_1 + \left(\tfrac{1}{2}-x_{14}+z_{14}\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-x_{14}+y_{14}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{14}\right)a\,\hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} + z_{14}c\,\hat{\mathbf{z}} \quad (8c) \quad \text{Zr III}$$

References:

- J. Galy and R. S. Roth, *The Crystal Structure of $Nb_2Zr_6O_{17}$*, J. Solid State Chem. **7**, 277–285 (1973), doi:10.1016/0022-4596(73)90134-5.

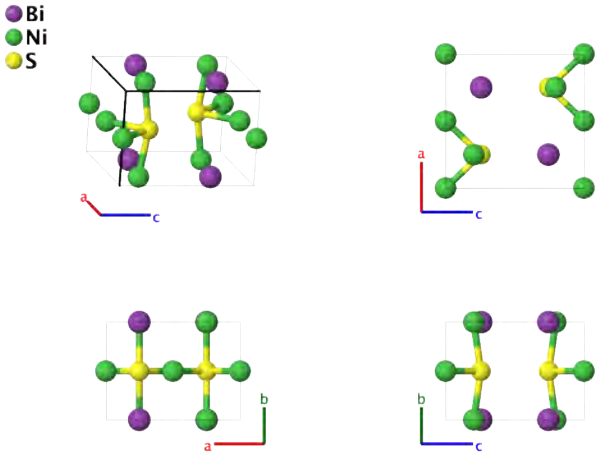
Found in:

- S. J. McCormack and W. M. Kriven, *Crystal structure solution for the $A_6B_2O_{17}$ (A = Zr, Hf; B = Nb, Ta) superstructure*, Acta Crystallogr. Sect. B Struct. Sci. **75**, 227–234 (2019), doi:10.1107/S2052520619001963.

Geometry files:
- CIF: pp. 1605
- POSCAR: pp. 1606

Parkerite ($Ni_3Bi_2S_2$) Structure: AB2C_oP8_51_e_be_f

http://aflow.org/prototype-encyclopedia/AB2C_oP8_51_e_be_f



Prototype	:	$Bi_2Ni_3S_2$
AFLOW prototype label	:	AB2C_oP8_51_e_be_f
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP8
Space group number	:	51
Space group symbol	:	*Pmma*
AFLOW prototype command	:	`aflow --proto=AB2C_oP8_51_e_be_f --params=`$a, b/a, c/a, z_2, z_3, z_4$

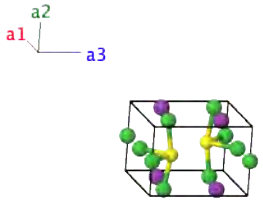
Other compounds with this structure

- $Ni_3(Bi,Pb)_2S_2$

- (Fleet, 1973) states that parkerite is a derivative of the shandite ($Ni_3Pb_2S_2$) structure, and changes to shandite if more than 4% of the bismuth is replaced by lead.
- The Ni-II ($2e$) site is occupied 50% of the time, given the observed stoichiometry.
- Earlier sources give parkerite a monoclinic structure. This may be due to an ordering of the nickel atoms at lower temperature. We follow (Downs, 2003) and use the orthorhombic structure.
- (Fleet, 1973) describes the structure in the *Pmam* setting of space group #51. We used FINDSYM to transform it to the standard *Pmma* setting.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}b\,\hat{\mathbf{y}}$	$(2b)$	Ni I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$	$(2b)$	Ni I
$\mathbf{B}_3$	=	$\frac{1}{4}\,\mathbf{a}_1 + z_2\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + z_2 c\,\hat{\mathbf{z}}$	$(2e)$	Bi
$\mathbf{B}_4$	=	$\frac{3}{4}\,\mathbf{a}_1 - z_2\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} - z_2 c\,\hat{\mathbf{z}}$	$(2e)$	Bi
$\mathbf{B}_5$	=	$\frac{1}{4}\,\mathbf{a}_1 + z_3\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + z_3 c\,\hat{\mathbf{z}}$	$(2e)$	Ni II
$\mathbf{B}_6$	=	$\frac{3}{4}\,\mathbf{a}_1 - z_3\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} - z_3 c\,\hat{\mathbf{z}}$	$(2e)$	Ni II
$\mathbf{B}_7$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(2f)$	S
$\mathbf{B}_8$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	$(2f)$	S

References:

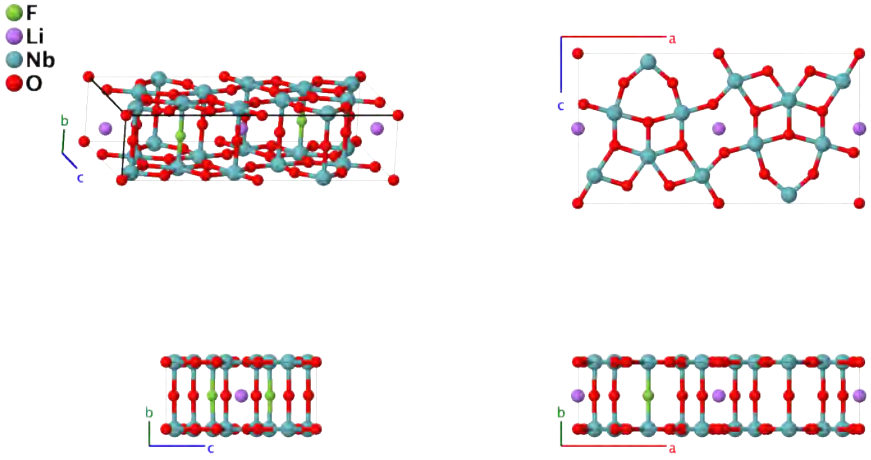
- M. E. Fleet, *The Crystal Structure of Parkerite ($Ni_3Bi_2S_2$)*, Am. Mineral. **58**, 435–439 (1973).
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1606
- POSCAR: pp. 1606

$LiNb_6O_{15}F$ Structure: ABC6D15_oP46_51_f_d_2e2i_aef4i2j

http://aflow.org/prototype-encyclopedia/ABC6D15_oP46_51_f_d_2e2i_aef4i2j

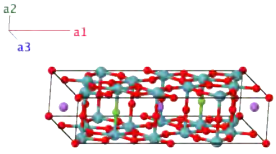


Prototype	:	$FLiNb_6O_{15}$
AFLOW prototype label	:	ABC6D15_oP46_51_f_d_2e2i_aef4i2j
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP46
Space group number	:	51
Space group symbol	:	*Pmma*
AFLOW prototype command	:	`aflow --proto=ABC6D15_oP46_51_f_d_2e2i_aef4i2j --params=`$a, b/a, c/a, z_3, z_4, z_5, z_6, z_7, x_8, z_8, x_9, z_9, x_{10}, z_{10}, x_{11}, z_{11}, x_{12}, z_{12}, x_{13}, z_{13}, x_{14}, z_{14}, x_{15}, z_{15}$

- (Lundberg, 1965) suggests that the lithium atoms are either on the (2*d*) site or are statistically distributed on a (4*j*) site with approximate coordinates (0.08,1/2,0.10). For simplicity we place the atoms on the (2*d*) site.
- The x-ray scattering of a F^- ion is almost identical to that of O^{2-}, and Lundberg was not able to distinguish between them. She arbitrarily designated the (2*f*) site she also called O_4 as the location of the fluorine ion and we follow this, but in reality we have no idea if the F^- ions are located on this site, are ordered on another site, or are statistically distributed on the oxygen sites.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	O I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	$(2a)$	O I
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(2d)$	Li
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(2d)$	Li
$\mathbf{B}_5$	=	$\frac{1}{4}\,\mathbf{a}_1 + z_3\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + z_3c\,\hat{\mathbf{z}}$	$(2e)$	Nb I
$\mathbf{B}_6$	=	$\frac{3}{4}\,\mathbf{a}_1 - z_3\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} - z_3c\,\hat{\mathbf{z}}$	$(2e)$	Nb I
$\mathbf{B}_7$	=	$\frac{1}{4}\,\mathbf{a}_1 + z_4\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + z_4c\,\hat{\mathbf{z}}$	$(2e)$	Nb II
$\mathbf{B}_8$	=	$\frac{3}{4}\,\mathbf{a}_1 - z_4\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} - z_4c\,\hat{\mathbf{z}}$	$(2e)$	Nb II
$\mathbf{B}_9$	=	$\frac{1}{4}\,\mathbf{a}_1 + z_5\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + z_5c\,\hat{\mathbf{z}}$	$(2e)$	O II
$\mathbf{B}_{10}$	=	$\frac{3}{4}\,\mathbf{a}_1 - z_5\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} - z_5c\,\hat{\mathbf{z}}$	$(2e)$	O II
$\mathbf{B}_{11}$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + z_6c\,\hat{\mathbf{z}}$	$(2f)$	F
$\mathbf{B}_{12}$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} - z_6c\,\hat{\mathbf{z}}$	$(2f)$	F
$\mathbf{B}_{13}$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + z_7c\,\hat{\mathbf{z}}$	$(2f)$	O III
$\mathbf{B}_{14}$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} - z_7c\,\hat{\mathbf{z}}$	$(2f)$	O III
$\mathbf{B}_{15}$	=	$x_8\,\mathbf{a}_1 + z_8\,\mathbf{a}_3$	=	$x_8a\,\hat{\mathbf{x}} + z_8c\,\hat{\mathbf{z}}$	$(4i)$	Nb III
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} - x_8\right)\mathbf{a}_1 + z_8\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_8\right)a\,\hat{\mathbf{x}} + z_8c\,\hat{\mathbf{z}}$	$(4i)$	Nb III
$\mathbf{B}_{17}$	=	$-x_8\,\mathbf{a}_1 - z_8\,\mathbf{a}_3$	=	$-x_8a\,\hat{\mathbf{x}} - z_8c\,\hat{\mathbf{z}}$	$(4i)$	Nb III
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} + x_8\right)\mathbf{a}_1 - z_8\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_8\right)a\,\hat{\mathbf{x}} - z_8c\,\hat{\mathbf{z}}$	$(4i)$	Nb III
$\mathbf{B}_{19}$	=	$x_9\,\mathbf{a}_1 + z_9\,\mathbf{a}_3$	=	$x_9a\,\hat{\mathbf{x}} + z_9c\,\hat{\mathbf{z}}$	$(4i)$	Nb IV
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} - x_9\right)\mathbf{a}_1 + z_9\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_9\right)a\,\hat{\mathbf{x}} + z_9c\,\hat{\mathbf{z}}$	$(4i)$	Nb IV
$\mathbf{B}_{21}$	=	$-x_9\,\mathbf{a}_1 - z_9\,\mathbf{a}_3$	=	$-x_9a\,\hat{\mathbf{x}} - z_9c\,\hat{\mathbf{z}}$	$(4i)$	Nb IV
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} + x_9\right)\mathbf{a}_1 - z_9\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_9\right)a\,\hat{\mathbf{x}} - z_9c\,\hat{\mathbf{z}}$	$(4i)$	Nb IV
$\mathbf{B}_{23}$	=	$x_{10}\,\mathbf{a}_1 + z_{10}\,\mathbf{a}_3$	=	$x_{10}a\,\hat{\mathbf{x}} + z_{10}c\,\hat{\mathbf{z}}$	$(4i)$	O IV
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} - x_{10}\right)\mathbf{a}_1 + z_{10}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{10}\right)a\,\hat{\mathbf{x}} + z_{10}c\,\hat{\mathbf{z}}$	$(4i)$	O IV
$\mathbf{B}_{25}$	=	$-x_{10}\,\mathbf{a}_1 - z_{10}\,\mathbf{a}_3$	=	$-x_{10}a\,\hat{\mathbf{x}} - z_{10}c\,\hat{\mathbf{z}}$	$(4i)$	O IV
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} + x_{10}\right)\mathbf{a}_1 - z_{10}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{10}\right)a\,\hat{\mathbf{x}} - z_{10}c\,\hat{\mathbf{z}}$	$(4i)$	O IV
$\mathbf{B}_{27}$	=	$x_{11}\,\mathbf{a}_1 + z_{11}\,\mathbf{a}_3$	=	$x_{11}a\,\hat{\mathbf{x}} + z_{11}c\,\hat{\mathbf{z}}$	$(4i)$	O V
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} - x_{11}\right)\mathbf{a}_1 + z_{11}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{11}\right)a\,\hat{\mathbf{x}} + z_{11}c\,\hat{\mathbf{z}}$	$(4i)$	O V
$\mathbf{B}_{29}$	=	$-x_{11}\,\mathbf{a}_1 - z_{11}\,\mathbf{a}_3$	=	$-x_{11}a\,\hat{\mathbf{x}} - z_{11}c\,\hat{\mathbf{z}}$	$(4i)$	O V
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2} + x_{11}\right)\mathbf{a}_1 - z_{11}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{11}\right)a\,\hat{\mathbf{x}} - z_{11}c\,\hat{\mathbf{z}}$	$(4i)$	O V
$\mathbf{B}_{31}$	=	$x_{12}\,\mathbf{a}_1 + z_{12}\,\mathbf{a}_3$	=	$x_{12}a\,\hat{\mathbf{x}} + z_{12}c\,\hat{\mathbf{z}}$	$(4i)$	O VI
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2} - x_{12}\right)\mathbf{a}_1 + z_{12}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{12}\right)a\,\hat{\mathbf{x}} + z_{12}c\,\hat{\mathbf{z}}$	$(4i)$	O VI
$\mathbf{B}_{33}$	=	$-x_{12}\,\mathbf{a}_1 - z_{12}\,\mathbf{a}_3$	=	$-x_{12}a\,\hat{\mathbf{x}} - z_{12}c\,\hat{\mathbf{z}}$	$(4i)$	O VI
$\mathbf{B}_{34}$	=	$\left(\frac{1}{2} + x_{12}\right)\mathbf{a}_1 - z_{12}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{12}\right)a\,\hat{\mathbf{x}} - z_{12}c\,\hat{\mathbf{z}}$	$(4i)$	O VI
$\mathbf{B}_{35}$	=	$x_{13}\,\mathbf{a}_1 + z_{13}\,\mathbf{a}_3$	=	$x_{13}a\,\hat{\mathbf{x}} + z_{13}c\,\hat{\mathbf{z}}$	$(4i)$	O VII

$\mathbf{B}_{36}$	=	$\left(\frac{1}{2} - x_{13}\right)\mathbf{a}_1 + z_{13}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{13}\right)a\,\hat{\mathbf{x}} + z_{13}c\,\hat{\mathbf{z}}$	$(4i)$	O VII
$\mathbf{B}_{37}$	=	$-x_{13}\,\mathbf{a}_1 - z_{13}\,\mathbf{a}_3$	=	$-x_{13}a\,\hat{\mathbf{x}} - z_{13}c\,\hat{\mathbf{z}}$	$(4i)$	O VII
$\mathbf{B}_{38}$	=	$\left(\frac{1}{2} + x_{13}\right)\mathbf{a}_1 - z_{13}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{13}\right)a\,\hat{\mathbf{x}} - z_{13}c\,\hat{\mathbf{z}}$	$(4i)$	O VII
$\mathbf{B}_{39}$	=	$x_{14}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3$	=	$x_{14}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + z_{14}c\,\hat{\mathbf{z}}$	$(4j)$	O VIII
$\mathbf{B}_{40}$	=	$\left(\frac{1}{2} - x_{14}\right)\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{14}\right)a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + z_{14}c\,\hat{\mathbf{z}}$	$(4j)$	O VIII
$\mathbf{B}_{41}$	=	$-x_{14}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 - z_{14}\,\mathbf{a}_3$	=	$-x_{14}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} - z_{14}c\,\hat{\mathbf{z}}$	$(4j)$	O VIII
$\mathbf{B}_{42}$	=	$\left(\frac{1}{2} + x_{14}\right)\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 - z_{14}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{14}\right)a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} - z_{14}c\,\hat{\mathbf{z}}$	$(4j)$	O VIII
$\mathbf{B}_{43}$	=	$x_{15}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3$	=	$x_{15}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + z_{15}c\,\hat{\mathbf{z}}$	$(4j)$	O IX
$\mathbf{B}_{44}$	=	$\left(\frac{1}{2} - x_{15}\right)\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{15}\right)a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + z_{15}c\,\hat{\mathbf{z}}$	$(4j)$	O IX
$\mathbf{B}_{45}$	=	$-x_{15}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 - z_{15}\,\mathbf{a}_3$	=	$-x_{15}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} - z_{15}c\,\hat{\mathbf{z}}$	$(4j)$	O IX
$\mathbf{B}_{46}$	=	$\left(\frac{1}{2} + x_{15}\right)\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 - z_{15}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{15}\right)a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} - z_{15}c\,\hat{\mathbf{z}}$	$(4j)$	O IX

References:

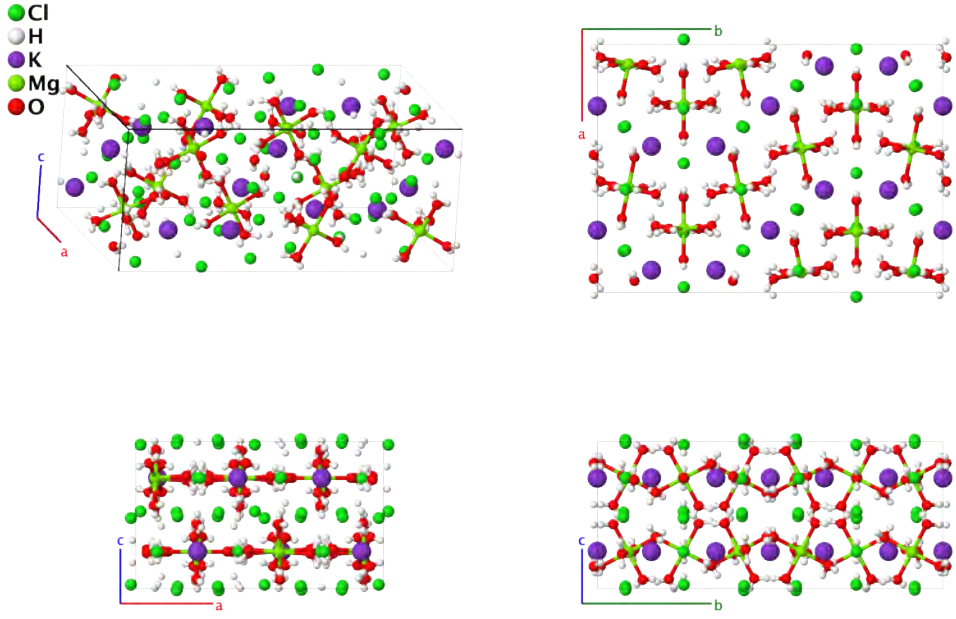
- M. Lundberg, *The Crystal Structure of $LiNb_6O_{15}F$*, Acta Chem. Scand. **19**, 2274–2284 (1965), doi:10.3891/acta.chem.scand.19-2274.

Geometry files:

- CIF: pp. 1606
- POSCAR: pp. 1607

Carnallite [$Mg(H_2O)_6KCl_3$] Structure: A3B12CDE6_oP276_52_d4e_18e_ce_de_2d8e

http://aflow.org/prototype-encyclopedia/A3B12CDE6_oP276_52_d4e_18e_ce_de_2d8e



Prototype	:	$Cl_3H_{12}KMgO_6$
AFLOW prototype label	:	A3B12CDE6_oP276_52_d4e_18e_ce_de_2d8e
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP276
Space group number	:	52
Space group symbol	:	*Pnna*
AFLOW prototype command	:	`aflow --proto=A3B12CDE6_oP276_52_d4e_18e_ce_de_2d8e --params=`$a, b/a, c/a, z_1, x_2, x_3, x_4, x_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}, x_{17}, y_{17}, z_{17}, x_{18}, y_{18}, z_{18}, x_{19}, y_{19}, z_{19}, x_{20}, y_{20}, z_{20}, x_{21}, y_{21}, z_{21}, x_{22}, y_{22}, z_{22}, x_{23}, y_{23}, z_{23}, x_{24}, y_{24}, z_{24}, x_{25}, y_{25}, z_{25}, x_{26}, y_{26}, z_{26}, x_{27}, y_{27}, z_{27}, x_{28}, y_{28}, z_{28}, x_{29}, y_{29}, z_{29}, x_{30}, y_{30}, z_{30}, x_{31}, y_{31}, z_{31}, x_{32}, y_{32}, z_{32}, x_{33}, y_{33}, z_{33}, x_{34}, y_{34}, z_{34}, x_{35}, y_{35}, z_{35}, x_{36}, y_{36}, z_{36}, x_{37}, y_{37}, z_{37}$

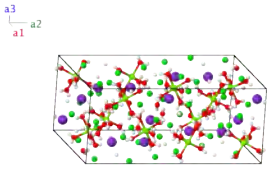
Other compounds with this structure

- $Mg(H_2O)_6K(Cl,Br)_3$

- (Andreß, 1939) determined the crystal structure of brom-carnallite, $Mg(H_2O)_6K(Cl,Br)_3$, finding that it was in space group $P4/n$ #85. (Herrmann, 1939) gave this structure the *Strukturbericht* designation $E2_6$. Later, (Schlemper, 1985), determined that the true space group was *Pnna* #52, and were able to determine the positions of the hydrogen atoms in the water molecules. We present this structure here. For the original structure, see the $E2_6$ page.

Simple Orthorhombic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= a\,\hat{\mathbf{x}}\\ \mathbf{a}_2 &= b\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$\frac{1}{4}\,\mathbf{a}_1 + z_1\,\mathbf{a}_3$	$\frac{1}{4}a\,\hat{\mathbf{x}} + z_1 c\,\hat{\mathbf{z}}$	($4c$)	K I
$\mathbf{B}_2$	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right)c\,\hat{\mathbf{z}}$	($4c$)	K I
$\mathbf{B}_3$	$\frac{3}{4}\,\mathbf{a}_1 - z_1\,\mathbf{a}_3$	$\frac{3}{4}a\,\hat{\mathbf{x}} - z_1 c\,\hat{\mathbf{z}}$	($4c$)	K I
$\mathbf{B}_4$	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\,\hat{\mathbf{z}}$	($4c$)	K I
$\mathbf{B}_5$	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	$x_2 a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($4d$)	Cl I
$\mathbf{B}_6$	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($4d$)	Cl I
$\mathbf{B}_7$	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	$-x_2 a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	($4d$)	Cl I
$\mathbf{B}_8$	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	($4d$)	Cl I
$\mathbf{B}_9$	$x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	$x_3 a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($4d$)	Mg I
$\mathbf{B}_{10}$	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($4d$)	Mg I
$\mathbf{B}_{11}$	$-x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	$-x_3 a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	($4d$)	Mg I
$\mathbf{B}_{12}$	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	($4d$)	Mg I
$\mathbf{B}_{13}$	$x_4\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	$x_4 a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($4d$)	O I
$\mathbf{B}_{14}$	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($4d$)	O I
$\mathbf{B}_{15}$	$-x_4\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	$-x_4 a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	($4d$)	O I
$\mathbf{B}_{16}$	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	($4d$)	O I
$\mathbf{B}_{17}$	$x_5\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	$x_5 a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($4d$)	O II
$\mathbf{B}_{18}$	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($4d$)	O II
$\mathbf{B}_{19}$	$-x_5\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	$-x_5 a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	($4d$)	O II
$\mathbf{B}_{20}$	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	($4d$)	O II
$\mathbf{B}_{21}$	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	($8e$)	Cl II
$\mathbf{B}_{22}$	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 - y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	$\left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	($8e$)	Cl II
$\mathbf{B}_{23}$	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\,\hat{\mathbf{z}}$	($8e$)	Cl II
$\mathbf{B}_{24}$	$x_6\,\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	$x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\,\hat{\mathbf{z}}$	($8e$)	Cl II
$\mathbf{B}_{25}$	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	$-x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	($8e$)	Cl II
$\mathbf{B}_{26}$	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	($8e$)	Cl II

$$\mathbf{B}_{27} = \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_6\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_6\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_6\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Cl II}$$

$$\mathbf{B}_{28} = -x_6\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_6\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_6\right)\mathbf{a}_3 = -x_6a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_6\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_6\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Cl II}$$

$$\mathbf{B}_{29} = x_7\,\mathbf{a}_1+y_7\,\mathbf{a}_2+z_7\,\mathbf{a}_3 = x_7a\,\hat{\mathbf{x}}+y_7b\,\hat{\mathbf{y}}+z_7c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Cl III}$$

$$\mathbf{B}_{30} = \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_1-y_7\,\mathbf{a}_2+z_7\,\mathbf{a}_3 = \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{x}}-y_7b\,\hat{\mathbf{y}}+z_7c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Cl III}$$

$$\mathbf{B}_{31} = \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_7\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_7\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Cl III}$$

$$\mathbf{B}_{32} = x_7\,\mathbf{a}_1+\left(\tfrac{1}{2}-y_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_7\right)\mathbf{a}_3 = x_7a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_7\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_7\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Cl III}$$

$$\mathbf{B}_{33} = -x_7\,\mathbf{a}_1-y_7\,\mathbf{a}_2-z_7\,\mathbf{a}_3 = -x_7a\,\hat{\mathbf{x}}-y_7b\,\hat{\mathbf{y}}-z_7c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Cl III}$$

$$\mathbf{B}_{34} = \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1+y_7\,\mathbf{a}_2-z_7\,\mathbf{a}_3 = \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{x}}+y_7b\,\hat{\mathbf{y}}-z_7c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Cl III}$$

$$\mathbf{B}_{35} = \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_7\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_7\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Cl III}$$

$$\mathbf{B}_{36} = -x_7\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 = -x_7a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_7\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_7\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Cl III}$$

$$\mathbf{B}_{37} = x_8\,\mathbf{a}_1+y_8\,\mathbf{a}_2+z_8\,\mathbf{a}_3 = x_8a\,\hat{\mathbf{x}}+y_8b\,\hat{\mathbf{y}}+z_8c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Cl IV}$$

$$\mathbf{B}_{38} = \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_1-y_8\,\mathbf{a}_2+z_8\,\mathbf{a}_3 = \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{x}}-y_8b\,\hat{\mathbf{y}}+z_8c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Cl IV}$$

$$\mathbf{B}_{39} = \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_8\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_8\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Cl IV}$$

$$\mathbf{B}_{40} = x_8\,\mathbf{a}_1+\left(\tfrac{1}{2}-y_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_8\right)\mathbf{a}_3 = x_8a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_8\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_8\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Cl IV}$$

$$\mathbf{B}_{41} = -x_8\,\mathbf{a}_1-y_8\,\mathbf{a}_2-z_8\,\mathbf{a}_3 = -x_8a\,\hat{\mathbf{x}}-y_8b\,\hat{\mathbf{y}}-z_8c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Cl IV}$$

$$\mathbf{B}_{42} = \left(\tfrac{1}{2}+x_8\right)\mathbf{a}_1+y_8\,\mathbf{a}_2-z_8\,\mathbf{a}_3 = \left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{x}}+y_8b\,\hat{\mathbf{y}}-z_8c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Cl IV}$$

$$\mathbf{B}_{43} = \left(\tfrac{1}{2}+x_8\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_8\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_8\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Cl IV}$$

$$\mathbf{B}_{44} = -x_8\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_8\right)\mathbf{a}_3 = -x_8a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_8\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_8\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Cl IV}$$

$$\mathbf{B}_{45} = x_9\,\mathbf{a}_1+y_9\,\mathbf{a}_2+z_9\,\mathbf{a}_3 = x_9a\,\hat{\mathbf{x}}+y_9b\,\hat{\mathbf{y}}+z_9c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Cl V}$$

$$\mathbf{B}_{46} = \left(\tfrac{1}{2}-x_9\right)\mathbf{a}_1-y_9\,\mathbf{a}_2+z_9\,\mathbf{a}_3 = \left(\tfrac{1}{2}-x_9\right)a\,\hat{\mathbf{x}}-y_9b\,\hat{\mathbf{y}}+z_9c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Cl V}$$

$$\mathbf{B}_{47} = \left(\tfrac{1}{2}-x_9\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_9\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_9\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_9\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_9\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Cl V}$$

$$\mathbf{B}_{48} = x_9\,\mathbf{a}_1+\left(\tfrac{1}{2}-y_9\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_9\right)\mathbf{a}_3 = x_9a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_9\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_9\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Cl V}$$

$$\mathbf{B}_{49} = -x_9\,\mathbf{a}_1-y_9\,\mathbf{a}_2-z_9\,\mathbf{a}_3 = -x_9a\,\hat{\mathbf{x}}-y_9b\,\hat{\mathbf{y}}-z_9c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Cl V}$$

$$\mathbf{B}_{50} = \left(\tfrac{1}{2}+x_9\right)\mathbf{a}_1+y_9\,\mathbf{a}_2-z_9\,\mathbf{a}_3 = \left(\tfrac{1}{2}+x_9\right)a\,\hat{\mathbf{x}}+y_9b\,\hat{\mathbf{y}}-z_9c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Cl V}$$

$$\mathbf{B}_{51} = \left(\tfrac{1}{2}+x_9\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_9\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_9\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_9\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_9\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Cl V}$$

$$\mathbf{B}_{52} = -x_9\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_9\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_9\right)\mathbf{a}_3 = -x_9a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_9\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_9\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Cl V}$$

$$\mathbf{B}_{53} = x_{10}\,\mathbf{a}_1+y_{10}\,\mathbf{a}_2+z_{10}\,\mathbf{a}_3 = x_{10}a\,\hat{\mathbf{x}}+y_{10}b\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H I}$$

$$\mathbf{B}_{54} = \left(\tfrac{1}{2}-x_{10}\right)\mathbf{a}_1-y_{10}\,\mathbf{a}_2+z_{10}\,\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{10}\right)a\,\hat{\mathbf{x}}-y_{10}b\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H I}$$

$$\mathbf{B}_{55} = \left(\tfrac{1}{2}-x_{10}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{10}\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{10}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{10}\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{10}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H I}$$

$$\mathbf{B}_{56} = x_{10}\,\mathbf{a}_1+\left(\tfrac{1}{2}-y_{10}\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_{10}\right)\mathbf{a}_3 = x_{10}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{10}\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{10}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H I}$$

$$\mathbf{B}_{57} = -x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 = -x_{10}a\,\hat{\mathbf{x}} - y_{10}b\,\hat{\mathbf{y}} - z_{10}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H I}$$

$$\mathbf{B}_{58} = \left(\tfrac{1}{2} + x_{10}\right)\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{10}\right)a\,\hat{\mathbf{x}} + y_{10}b\,\hat{\mathbf{y}} - z_{10}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H I}$$

$$\mathbf{B}_{59} = \left(\tfrac{1}{2} + x_{10}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{10}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{10}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{10}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{10}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H I}$$

$$\mathbf{B}_{60} = -x_{10}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{10}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{10}\right)\mathbf{a}_3 = -x_{10}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{10}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{10}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H I}$$

$$\mathbf{B}_{61} = x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 = x_{11}a\,\hat{\mathbf{x}} + y_{11}b\,\hat{\mathbf{y}} + z_{11}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H II}$$

$$\mathbf{B}_{62} = \left(\tfrac{1}{2} - x_{11}\right)\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{11}\right)a\,\hat{\mathbf{x}} - y_{11}b\,\hat{\mathbf{y}} + z_{11}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H II}$$

$$\mathbf{B}_{63} = \left(\tfrac{1}{2} - x_{11}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{11}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{11}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{11}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{11}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H II}$$

$$\mathbf{B}_{64} = x_{11}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{11}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{11}\right)\mathbf{a}_3 = x_{11}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{11}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{11}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H II}$$

$$\mathbf{B}_{65} = -x_{11}\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3 = -x_{11}a\,\hat{\mathbf{x}} - y_{11}b\,\hat{\mathbf{y}} - z_{11}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H II}$$

$$\mathbf{B}_{66} = \left(\tfrac{1}{2} + x_{11}\right)\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{11}\right)a\,\hat{\mathbf{x}} + y_{11}b\,\hat{\mathbf{y}} - z_{11}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H II}$$

$$\mathbf{B}_{67} = \left(\tfrac{1}{2} + x_{11}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{11}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{11}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{11}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{11}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H II}$$

$$\mathbf{B}_{68} = -x_{11}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{11}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{11}\right)\mathbf{a}_3 = -x_{11}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{11}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{11}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H II}$$

$$\mathbf{B}_{69} = x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 = x_{12}a\,\hat{\mathbf{x}} + y_{12}b\,\hat{\mathbf{y}} + z_{12}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H III}$$

$$\mathbf{B}_{70} = \left(\tfrac{1}{2} - x_{12}\right)\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{12}\right)a\,\hat{\mathbf{x}} - y_{12}b\,\hat{\mathbf{y}} + z_{12}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H III}$$

$$\mathbf{B}_{71} = \left(\tfrac{1}{2} - x_{12}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{12}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{12}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{12}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{12}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H III}$$

$$\mathbf{B}_{72} = x_{12}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{12}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{12}\right)\mathbf{a}_3 = x_{12}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{12}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{12}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H III}$$

$$\mathbf{B}_{73} = -x_{12}\,\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3 = -x_{12}a\,\hat{\mathbf{x}} - y_{12}b\,\hat{\mathbf{y}} - z_{12}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H III}$$

$$\mathbf{B}_{74} = \left(\tfrac{1}{2} + x_{12}\right)\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{12}\right)a\,\hat{\mathbf{x}} + y_{12}b\,\hat{\mathbf{y}} - z_{12}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H III}$$

$$\mathbf{B}_{75} = \left(\tfrac{1}{2} + x_{12}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{12}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{12}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{12}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{12}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H III}$$

$$\mathbf{B}_{76} = -x_{12}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{12}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{12}\right)\mathbf{a}_3 = -x_{12}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{12}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{12}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H III}$$

$$\mathbf{B}_{77} = x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3 = x_{13}a\,\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} + z_{13}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H IV}$$

$$\mathbf{B}_{78} = \left(\tfrac{1}{2} - x_{13}\right)\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{13}\right)a\,\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} + z_{13}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H IV}$$

$$\mathbf{B}_{79} = \left(\tfrac{1}{2} - x_{13}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{13}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{13}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{13}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{13}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H IV}$$

$$\mathbf{B}_{80} = x_{13}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{13}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{13}\right)\mathbf{a}_3 = x_{13}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{13}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{13}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H IV}$$

$$\mathbf{B}_{81} = -x_{13}\,\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 - z_{13}\,\mathbf{a}_3 = -x_{13}a\,\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} - z_{13}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H IV}$$

$$\mathbf{B}_{82} = \left(\tfrac{1}{2} + x_{13}\right)\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 - z_{13}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{13}\right)a\,\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} - z_{13}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H IV}$$

$$\mathbf{B}_{83} = \left(\tfrac{1}{2} + x_{13}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{13}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{13}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{13}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{13}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H IV}$$

$$\mathbf{B}_{84} = -x_{13}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{13}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{13}\right)\mathbf{a}_3 = -x_{13}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{13}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{13}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H IV}$$

$$
\begin{array}{lcccccc}
\mathbf{B}_{85} & = & x_{14}\,\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3 & = & x_{14}a\,\hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} + z_{14}c\,\hat{\mathbf{z}} & (8e) & \text{H V} \\
\mathbf{B}_{86} & = & \left(\tfrac{1}{2} - x_{14}\right)\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3 & = & \left(\tfrac{1}{2} - x_{14}\right)a\,\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} + z_{14}c\,\hat{\mathbf{z}} & (8e) & \text{H V} \\
\mathbf{B}_{87} & = & \left(\tfrac{1}{2} - x_{14}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{14}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{14}\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} - x_{14}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{14}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{14}\right)c\,\hat{\mathbf{z}} & (8e) & \text{H V} \\
\mathbf{B}_{88} & = & x_{14}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{14}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{14}\right)\mathbf{a}_3 & = & x_{14}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{14}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{14}\right)c\,\hat{\mathbf{z}} & (8e) & \text{H V} \\
\mathbf{B}_{89} & = & -x_{14}\,\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 - z_{14}\,\mathbf{a}_3 & = & -x_{14}a\,\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} - z_{14}c\,\hat{\mathbf{z}} & (8e) & \text{H V} \\
\mathbf{B}_{90} & = & \left(\tfrac{1}{2} + x_{14}\right)\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 - z_{14}\,\mathbf{a}_3 & = & \left(\tfrac{1}{2} + x_{14}\right)a\,\hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} - z_{14}c\,\hat{\mathbf{z}} & (8e) & \text{H V} \\
\mathbf{B}_{91} & = & \left(\tfrac{1}{2} + x_{14}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{14}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{14}\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} + x_{14}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{14}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{14}\right)c\,\hat{\mathbf{z}} & (8e) & \text{H V} \\
\mathbf{B}_{92} & = & -x_{14}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{14}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{14}\right)\mathbf{a}_3 & = & -x_{14}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{14}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{14}\right)c\,\hat{\mathbf{z}} & (8e) & \text{H V} \\
\mathbf{B}_{93} & = & x_{15}\,\mathbf{a}_1 + y_{15}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 & = & x_{15}a\,\hat{\mathbf{x}} + y_{15}b\,\hat{\mathbf{y}} + z_{15}c\,\hat{\mathbf{z}} & (8e) & \text{H VI} \\
\mathbf{B}_{94} & = & \left(\tfrac{1}{2} - x_{15}\right)\mathbf{a}_1 - y_{15}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 & = & \left(\tfrac{1}{2} - x_{15}\right)a\,\hat{\mathbf{x}} - y_{15}b\,\hat{\mathbf{y}} + z_{15}c\,\hat{\mathbf{z}} & (8e) & \text{H VI} \\
\mathbf{B}_{95} & = & \left(\tfrac{1}{2} - x_{15}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{15}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{15}\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} - x_{15}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{15}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{15}\right)c\,\hat{\mathbf{z}} & (8e) & \text{H VI} \\
\mathbf{B}_{96} & = & x_{15}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{15}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{15}\right)\mathbf{a}_3 & = & x_{15}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{15}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{15}\right)c\,\hat{\mathbf{z}} & (8e) & \text{H VI} \\
\mathbf{B}_{97} & = & -x_{15}\,\mathbf{a}_1 - y_{15}\,\mathbf{a}_2 - z_{15}\,\mathbf{a}_3 & = & -x_{15}a\,\hat{\mathbf{x}} - y_{15}b\,\hat{\mathbf{y}} - z_{15}c\,\hat{\mathbf{z}} & (8e) & \text{H VI} \\
\mathbf{B}_{98} & = & \left(\tfrac{1}{2} + x_{15}\right)\mathbf{a}_1 + y_{15}\,\mathbf{a}_2 - z_{15}\,\mathbf{a}_3 & = & \left(\tfrac{1}{2} + x_{15}\right)a\,\hat{\mathbf{x}} + y_{15}b\,\hat{\mathbf{y}} - z_{15}c\,\hat{\mathbf{z}} & (8e) & \text{H VI} \\
\mathbf{B}_{99} & = & \left(\tfrac{1}{2} + x_{15}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{15}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{15}\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} + x_{15}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{15}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{15}\right)c\,\hat{\mathbf{z}} & (8e) & \text{H VI} \\
\mathbf{B}_{100} & = & -x_{15}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{15}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{15}\right)\mathbf{a}_3 & = & -x_{15}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{15}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{15}\right)c\,\hat{\mathbf{z}} & (8e) & \text{H VI} \\
\mathbf{B}_{101} & = & x_{16}\,\mathbf{a}_1 + y_{16}\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 & = & x_{16}a\,\hat{\mathbf{x}} + y_{16}b\,\hat{\mathbf{y}} + z_{16}c\,\hat{\mathbf{z}} & (8e) & \text{H VII} \\
\mathbf{B}_{102} & = & \left(\tfrac{1}{2} - x_{16}\right)\mathbf{a}_1 - y_{16}\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 & = & \left(\tfrac{1}{2} - x_{16}\right)a\,\hat{\mathbf{x}} - y_{16}b\,\hat{\mathbf{y}} + z_{16}c\,\hat{\mathbf{z}} & (8e) & \text{H VII} \\
\mathbf{B}_{103} & = & \left(\tfrac{1}{2} - x_{16}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{16}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{16}\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} - x_{16}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{16}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{16}\right)c\,\hat{\mathbf{z}} & (8e) & \text{H VII} \\
\mathbf{B}_{104} & = & x_{16}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{16}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{16}\right)\mathbf{a}_3 & = & x_{16}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{16}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{16}\right)c\,\hat{\mathbf{z}} & (8e) & \text{H VII} \\
\mathbf{B}_{105} & = & -x_{16}\,\mathbf{a}_1 - y_{16}\,\mathbf{a}_2 - z_{16}\,\mathbf{a}_3 & = & -x_{16}a\,\hat{\mathbf{x}} - y_{16}b\,\hat{\mathbf{y}} - z_{16}c\,\hat{\mathbf{z}} & (8e) & \text{H VII} \\
\mathbf{B}_{106} & = & \left(\tfrac{1}{2} + x_{16}\right)\mathbf{a}_1 + y_{16}\,\mathbf{a}_2 - z_{16}\,\mathbf{a}_3 & = & \left(\tfrac{1}{2} + x_{16}\right)a\,\hat{\mathbf{x}} + y_{16}b\,\hat{\mathbf{y}} - z_{16}c\,\hat{\mathbf{z}} & (8e) & \text{H VII} \\
\mathbf{B}_{107} & = & \left(\tfrac{1}{2} + x_{16}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{16}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{16}\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} + x_{16}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{16}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{16}\right)c\,\hat{\mathbf{z}} & (8e) & \text{H VII} \\
\mathbf{B}_{108} & = & -x_{16}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{16}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{16}\right)\mathbf{a}_3 & = & -x_{16}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{16}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{16}\right)c\,\hat{\mathbf{z}} & (8e) & \text{H VII} \\
\mathbf{B}_{109} & = & x_{17}\,\mathbf{a}_1 + y_{17}\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 & = & x_{17}a\,\hat{\mathbf{x}} + y_{17}b\,\hat{\mathbf{y}} + z_{17}c\,\hat{\mathbf{z}} & (8e) & \text{H VIII} \\
\mathbf{B}_{110} & = & \left(\tfrac{1}{2} - x_{17}\right)\mathbf{a}_1 - y_{17}\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 & = & \left(\tfrac{1}{2} - x_{17}\right)a\,\hat{\mathbf{x}} - y_{17}b\,\hat{\mathbf{y}} + z_{17}c\,\hat{\mathbf{z}} & (8e) & \text{H VIII} \\
\mathbf{B}_{111} & = & \left(\tfrac{1}{2} - x_{17}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{17}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{17}\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} - x_{17}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{17}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{17}\right)c\,\hat{\mathbf{z}} & (8e) & \text{H VIII} \\
\mathbf{B}_{112} & = & x_{17}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{17}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{17}\right)\mathbf{a}_3 & = & x_{17}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{17}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{17}\right)c\,\hat{\mathbf{z}} & (8e) & \text{H VIII}
\end{array}
$$

$$\mathbf{B}_{113} = -x_{17}\,\mathbf{a}_1 - y_{17}\,\mathbf{a}_2 - z_{17}\,\mathbf{a}_3 = -x_{17}a\,\hat{\mathbf{x}} - y_{17}b\,\hat{\mathbf{y}} - z_{17}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H VIII}$$

$$\mathbf{B}_{114} = \left(\tfrac{1}{2}+x_{17}\right)\mathbf{a}_1 + y_{17}\,\mathbf{a}_2 - z_{17}\,\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{17}\right)a\,\hat{\mathbf{x}} + y_{17}b\,\hat{\mathbf{y}} - z_{17}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H VIII}$$

$$\mathbf{B}_{115} = \left(\tfrac{1}{2}+x_{17}\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_{17}\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{17}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{17}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_{17}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{17}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H VIII}$$

$$\mathbf{B}_{116} = -x_{17}\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_{17}\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{17}\right)\mathbf{a}_3 = -x_{17}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_{17}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{17}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H VIII}$$

$$\mathbf{B}_{117} = x_{18}\,\mathbf{a}_1 + y_{18}\,\mathbf{a}_2 + z_{18}\,\mathbf{a}_3 = x_{18}a\,\hat{\mathbf{x}} + y_{18}b\,\hat{\mathbf{y}} + z_{18}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H IX}$$

$$\mathbf{B}_{118} = \left(\tfrac{1}{2}-x_{18}\right)\mathbf{a}_1 - y_{18}\,\mathbf{a}_2 + z_{18}\,\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{18}\right)a\,\hat{\mathbf{x}} - y_{18}b\,\hat{\mathbf{y}} + z_{18}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H IX}$$

$$\mathbf{B}_{119} = \left(\tfrac{1}{2}-x_{18}\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_{18}\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_{18}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{18}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_{18}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_{18}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H IX}$$

$$\mathbf{B}_{120} = x_{18}\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_{18}\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_{18}\right)\mathbf{a}_3 = x_{18}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_{18}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_{18}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H IX}$$

$$\mathbf{B}_{121} = -x_{18}\,\mathbf{a}_1 - y_{18}\,\mathbf{a}_2 - z_{18}\,\mathbf{a}_3 = -x_{18}a\,\hat{\mathbf{x}} - y_{18}b\,\hat{\mathbf{y}} - z_{18}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H IX}$$

$$\mathbf{B}_{122} = \left(\tfrac{1}{2}+x_{18}\right)\mathbf{a}_1 + y_{18}\,\mathbf{a}_2 - z_{18}\,\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{18}\right)a\,\hat{\mathbf{x}} + y_{18}b\,\hat{\mathbf{y}} - z_{18}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H IX}$$

$$\mathbf{B}_{123} = \left(\tfrac{1}{2}+x_{18}\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_{18}\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{18}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{18}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_{18}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{18}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H IX}$$

$$\mathbf{B}_{124} = -x_{18}\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_{18}\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{18}\right)\mathbf{a}_3 = -x_{18}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_{18}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{18}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H IX}$$

$$\mathbf{B}_{125} = x_{19}\,\mathbf{a}_1 + y_{19}\,\mathbf{a}_2 + z_{19}\,\mathbf{a}_3 = x_{19}a\,\hat{\mathbf{x}} + y_{19}b\,\hat{\mathbf{y}} + z_{19}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H X}$$

$$\mathbf{B}_{126} = \left(\tfrac{1}{2}-x_{19}\right)\mathbf{a}_1 - y_{19}\,\mathbf{a}_2 + z_{19}\,\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{19}\right)a\,\hat{\mathbf{x}} - y_{19}b\,\hat{\mathbf{y}} + z_{19}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H X}$$

$$\mathbf{B}_{127} = \left(\tfrac{1}{2}-x_{19}\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_{19}\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_{19}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{19}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_{19}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_{19}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H X}$$

$$\mathbf{B}_{128} = x_{19}\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_{19}\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_{19}\right)\mathbf{a}_3 = x_{19}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_{19}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_{19}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H X}$$

$$\mathbf{B}_{129} = -x_{19}\,\mathbf{a}_1 - y_{19}\,\mathbf{a}_2 - z_{19}\,\mathbf{a}_3 = -x_{19}a\,\hat{\mathbf{x}} - y_{19}b\,\hat{\mathbf{y}} - z_{19}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H X}$$

$$\mathbf{B}_{130} = \left(\tfrac{1}{2}+x_{19}\right)\mathbf{a}_1 + y_{19}\,\mathbf{a}_2 - z_{19}\,\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{19}\right)a\,\hat{\mathbf{x}} + y_{19}b\,\hat{\mathbf{y}} - z_{19}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H X}$$

$$\mathbf{B}_{131} = \left(\tfrac{1}{2}+x_{19}\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_{19}\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{19}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{19}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_{19}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{19}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H X}$$

$$\mathbf{B}_{132} = -x_{19}\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_{19}\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{19}\right)\mathbf{a}_3 = -x_{19}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_{19}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{19}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H X}$$

$$\mathbf{B}_{133} = x_{20}\,\mathbf{a}_1 + y_{20}\,\mathbf{a}_2 + z_{20}\,\mathbf{a}_3 = x_{20}a\,\hat{\mathbf{x}} + y_{20}b\,\hat{\mathbf{y}} + z_{20}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XI}$$

$$\mathbf{B}_{134} = \left(\tfrac{1}{2}-x_{20}\right)\mathbf{a}_1 - y_{20}\,\mathbf{a}_2 + z_{20}\,\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{20}\right)a\,\hat{\mathbf{x}} - y_{20}b\,\hat{\mathbf{y}} + z_{20}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XI}$$

$$\mathbf{B}_{135} = \left(\tfrac{1}{2}-x_{20}\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_{20}\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_{20}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{20}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_{20}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_{20}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XI}$$

$$\mathbf{B}_{136} = x_{20}\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_{20}\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_{20}\right)\mathbf{a}_3 = x_{20}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_{20}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_{20}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XI}$$

$$\mathbf{B}_{137} = -x_{20}\,\mathbf{a}_1 - y_{20}\,\mathbf{a}_2 - z_{20}\,\mathbf{a}_3 = -x_{20}a\,\hat{\mathbf{x}} - y_{20}b\,\hat{\mathbf{y}} - z_{20}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XI}$$

$$\mathbf{B}_{138} = \left(\tfrac{1}{2}+x_{20}\right)\mathbf{a}_1 + y_{20}\,\mathbf{a}_2 - z_{20}\,\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{20}\right)a\,\hat{\mathbf{x}} + y_{20}b\,\hat{\mathbf{y}} - z_{20}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XI}$$

$$\mathbf{B}_{139} = \left(\tfrac{1}{2}+x_{20}\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_{20}\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{20}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{20}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_{20}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{20}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XI}$$

$$\mathbf{B}_{140} = -x_{20}\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_{20}\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{20}\right)\mathbf{a}_3 = -x_{20}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_{20}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{20}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XI}$$

$$\mathbf{B}_{141} = x_{21}\,\mathbf{a}_1 + y_{21}\,\mathbf{a}_2 + z_{21}\,\mathbf{a}_3 = x_{21}a\,\hat{\mathbf{x}} + y_{21}b\,\hat{\mathbf{y}} + z_{21}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XII}$$

$$\mathbf{B}_{142} = \left(\tfrac{1}{2} - x_{21}\right)\mathbf{a}_1 - y_{21}\,\mathbf{a}_2 + z_{21}\,\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{21}\right)a\,\hat{\mathbf{x}} - y_{21}b\,\hat{\mathbf{y}} + z_{21}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XII}$$

$$\mathbf{B}_{143} = \left(\tfrac{1}{2} - x_{21}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{21}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{21}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{21}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{21}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{21}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XII}$$

$$\mathbf{B}_{144} = x_{21}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{21}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{21}\right)\mathbf{a}_3 = x_{21}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{21}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{21}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XII}$$

$$\mathbf{B}_{145} = -x_{21}\,\mathbf{a}_1 - y_{21}\,\mathbf{a}_2 - z_{21}\,\mathbf{a}_3 = -x_{21}a\,\hat{\mathbf{x}} - y_{21}b\,\hat{\mathbf{y}} - z_{21}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XII}$$

$$\mathbf{B}_{146} = \left(\tfrac{1}{2} + x_{21}\right)\mathbf{a}_1 + y_{21}\,\mathbf{a}_2 - z_{21}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{21}\right)a\,\hat{\mathbf{x}} + y_{21}b\,\hat{\mathbf{y}} - z_{21}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XII}$$

$$\mathbf{B}_{147} = \left(\tfrac{1}{2} + x_{21}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{21}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{21}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{21}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{21}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{21}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XII}$$

$$\mathbf{B}_{148} = -x_{21}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{21}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{21}\right)\mathbf{a}_3 = -x_{21}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{21}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{21}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XII}$$

$$\mathbf{B}_{149} = x_{22}\,\mathbf{a}_1 + y_{22}\,\mathbf{a}_2 + z_{22}\,\mathbf{a}_3 = x_{22}a\,\hat{\mathbf{x}} + y_{22}b\,\hat{\mathbf{y}} + z_{22}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XIII}$$

$$\mathbf{B}_{150} = \left(\tfrac{1}{2} - x_{22}\right)\mathbf{a}_1 - y_{22}\,\mathbf{a}_2 + z_{22}\,\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{22}\right)a\,\hat{\mathbf{x}} - y_{22}b\,\hat{\mathbf{y}} + z_{22}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XIII}$$

$$\mathbf{B}_{151} = \left(\tfrac{1}{2} - x_{22}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{22}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{22}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{22}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{22}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{22}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XIII}$$

$$\mathbf{B}_{152} = x_{22}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{22}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{22}\right)\mathbf{a}_3 = x_{22}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{22}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{22}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XIII}$$

$$\mathbf{B}_{153} = -x_{22}\,\mathbf{a}_1 - y_{22}\,\mathbf{a}_2 - z_{22}\,\mathbf{a}_3 = -x_{22}a\,\hat{\mathbf{x}} - y_{22}b\,\hat{\mathbf{y}} - z_{22}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XIII}$$

$$\mathbf{B}_{154} = \left(\tfrac{1}{2} + x_{22}\right)\mathbf{a}_1 + y_{22}\,\mathbf{a}_2 - z_{22}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{22}\right)a\,\hat{\mathbf{x}} + y_{22}b\,\hat{\mathbf{y}} - z_{22}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XIII}$$

$$\mathbf{B}_{155} = \left(\tfrac{1}{2} + x_{22}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{22}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{22}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{22}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{22}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{22}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XIII}$$

$$\mathbf{B}_{156} = -x_{22}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{22}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{22}\right)\mathbf{a}_3 = -x_{22}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{22}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{22}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XIII}$$

$$\mathbf{B}_{157} = x_{23}\,\mathbf{a}_1 + y_{23}\,\mathbf{a}_2 + z_{23}\,\mathbf{a}_3 = x_{23}a\,\hat{\mathbf{x}} + y_{23}b\,\hat{\mathbf{y}} + z_{23}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XIV}$$

$$\mathbf{B}_{158} = \left(\tfrac{1}{2} - x_{23}\right)\mathbf{a}_1 - y_{23}\,\mathbf{a}_2 + z_{23}\,\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{23}\right)a\,\hat{\mathbf{x}} - y_{23}b\,\hat{\mathbf{y}} + z_{23}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XIV}$$

$$\mathbf{B}_{159} = \left(\tfrac{1}{2} - x_{23}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{23}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{23}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{23}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{23}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{23}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XIV}$$

$$\mathbf{B}_{160} = x_{23}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{23}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{23}\right)\mathbf{a}_3 = x_{23}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{23}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{23}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XIV}$$

$$\mathbf{B}_{161} = -x_{23}\,\mathbf{a}_1 - y_{23}\,\mathbf{a}_2 - z_{23}\,\mathbf{a}_3 = -x_{23}a\,\hat{\mathbf{x}} - y_{23}b\,\hat{\mathbf{y}} - z_{23}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XIV}$$

$$\mathbf{B}_{162} = \left(\tfrac{1}{2} + x_{23}\right)\mathbf{a}_1 + y_{23}\,\mathbf{a}_2 - z_{23}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{23}\right)a\,\hat{\mathbf{x}} + y_{23}b\,\hat{\mathbf{y}} - z_{23}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XIV}$$

$$\mathbf{B}_{163} = \left(\tfrac{1}{2} + x_{23}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{23}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{23}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{23}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{23}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{23}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XIV}$$

$$\mathbf{B}_{164} = -x_{23}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{23}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{23}\right)\mathbf{a}_3 = -x_{23}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{23}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{23}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XIV}$$

$$\mathbf{B}_{165} = x_{24}\,\mathbf{a}_1 + y_{24}\,\mathbf{a}_2 + z_{24}\,\mathbf{a}_3 = x_{24}a\,\hat{\mathbf{x}} + y_{24}b\,\hat{\mathbf{y}} + z_{24}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XV}$$

$$\mathbf{B}_{166} = \left(\tfrac{1}{2} - x_{24}\right)\mathbf{a}_1 - y_{24}\,\mathbf{a}_2 + z_{24}\,\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{24}\right)a\,\hat{\mathbf{x}} - y_{24}b\,\hat{\mathbf{y}} + z_{24}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XV}$$

$$\mathbf{B}_{167} = \left(\tfrac{1}{2} - x_{24}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{24}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{24}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{24}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{24}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{24}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XV}$$

$$\mathbf{B}_{168} = x_{24}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{24}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{24}\right)\mathbf{a}_3 = x_{24}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{24}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{24}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XV}$$

$$\mathbf{B}_{169} = -x_{24}\,\mathbf{a}_1 - y_{24}\,\mathbf{a}_2 - z_{24}\,\mathbf{a}_3 = -x_{24}a\,\hat{\mathbf{x}} - y_{24}b\,\hat{\mathbf{y}} - z_{24}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XV}$$

$$\mathbf{B}_{170} = \left(\tfrac{1}{2} + x_{24}\right)\mathbf{a}_1 + y_{24}\,\mathbf{a}_2 - z_{24}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{24}\right)a\,\hat{\mathbf{x}} + y_{24}b\,\hat{\mathbf{y}} - z_{24}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XV}$$

$$\mathbf{B}_{171} = \left(\tfrac{1}{2} + x_{24}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{24}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{24}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{24}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{24}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{24}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XV}$$

$$\mathbf{B}_{172} = -x_{24}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{24}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{24}\right)\mathbf{a}_3 = -x_{24}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{24}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{24}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XV}$$

$$\mathbf{B}_{173} = x_{25}\,\mathbf{a}_1 + y_{25}\,\mathbf{a}_2 + z_{25}\,\mathbf{a}_3 = x_{25}a\,\hat{\mathbf{x}} + y_{25}b\,\hat{\mathbf{y}} + z_{25}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XVI}$$

$$\mathbf{B}_{174} = \left(\tfrac{1}{2} - x_{25}\right)\mathbf{a}_1 - y_{25}\,\mathbf{a}_2 + z_{25}\,\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{25}\right)a\,\hat{\mathbf{x}} - y_{25}b\,\hat{\mathbf{y}} + z_{25}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XVI}$$

$$\mathbf{B}_{175} = \left(\tfrac{1}{2} - x_{25}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{25}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{25}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{25}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{25}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{25}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XVI}$$

$$\mathbf{B}_{176} = x_{25}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{25}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{25}\right)\mathbf{a}_3 = x_{25}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{25}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{25}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XVI}$$

$$\mathbf{B}_{177} = -x_{25}\,\mathbf{a}_1 - y_{25}\,\mathbf{a}_2 - z_{25}\,\mathbf{a}_3 = -x_{25}a\,\hat{\mathbf{x}} - y_{25}b\,\hat{\mathbf{y}} - z_{25}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XVI}$$

$$\mathbf{B}_{178} = \left(\tfrac{1}{2} + x_{25}\right)\mathbf{a}_1 + y_{25}\,\mathbf{a}_2 - z_{25}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{25}\right)a\,\hat{\mathbf{x}} + y_{25}b\,\hat{\mathbf{y}} - z_{25}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XVI}$$

$$\mathbf{B}_{179} = \left(\tfrac{1}{2} + x_{25}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{25}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{25}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{25}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{25}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{25}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XVI}$$

$$\mathbf{B}_{180} = -x_{25}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{25}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{25}\right)\mathbf{a}_3 = -x_{25}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{25}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{25}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XVI}$$

$$\mathbf{B}_{181} = x_{26}\,\mathbf{a}_1 + y_{26}\,\mathbf{a}_2 + z_{26}\,\mathbf{a}_3 = x_{26}a\,\hat{\mathbf{x}} + y_{26}b\,\hat{\mathbf{y}} + z_{26}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XVII}$$

$$\mathbf{B}_{182} = \left(\tfrac{1}{2} - x_{26}\right)\mathbf{a}_1 - y_{26}\,\mathbf{a}_2 + z_{26}\,\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{26}\right)a\,\hat{\mathbf{x}} - y_{26}b\,\hat{\mathbf{y}} + z_{26}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XVII}$$

$$\mathbf{B}_{183} = \left(\tfrac{1}{2} - x_{26}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{26}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{26}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{26}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{26}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{26}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XVII}$$

$$\mathbf{B}_{184} = x_{26}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{26}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{26}\right)\mathbf{a}_3 = x_{26}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{26}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{26}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XVII}$$

$$\mathbf{B}_{185} = -x_{26}\,\mathbf{a}_1 - y_{26}\,\mathbf{a}_2 - z_{26}\,\mathbf{a}_3 = -x_{26}a\,\hat{\mathbf{x}} - y_{26}b\,\hat{\mathbf{y}} - z_{26}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XVII}$$

$$\mathbf{B}_{186} = \left(\tfrac{1}{2} + x_{26}\right)\mathbf{a}_1 + y_{26}\,\mathbf{a}_2 - z_{26}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{26}\right)a\,\hat{\mathbf{x}} + y_{26}b\,\hat{\mathbf{y}} - z_{26}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XVII}$$

$$\mathbf{B}_{187} = \left(\tfrac{1}{2} + x_{26}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{26}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{26}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{26}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{26}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{26}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XVII}$$

$$\mathbf{B}_{188} = -x_{26}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{26}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{26}\right)\mathbf{a}_3 = -x_{26}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{26}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{26}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XVII}$$

$$\mathbf{B}_{189} = x_{27}\,\mathbf{a}_1 + y_{27}\,\mathbf{a}_2 + z_{27}\,\mathbf{a}_3 = x_{27}a\,\hat{\mathbf{x}} + y_{27}b\,\hat{\mathbf{y}} + z_{27}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XVIII}$$

$$\mathbf{B}_{190} = \left(\tfrac{1}{2} - x_{27}\right)\mathbf{a}_1 - y_{27}\,\mathbf{a}_2 + z_{27}\,\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{27}\right)a\,\hat{\mathbf{x}} - y_{27}b\,\hat{\mathbf{y}} + z_{27}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XVIII}$$

$$\mathbf{B}_{191} = \left(\tfrac{1}{2} - x_{27}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{27}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{27}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{27}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{27}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{27}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XVIII}$$

$$\mathbf{B}_{192} = x_{27}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{27}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{27}\right)\mathbf{a}_3 = x_{27}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{27}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{27}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XVIII}$$

$$\mathbf{B}_{193} = -x_{27}\,\mathbf{a}_1 - y_{27}\,\mathbf{a}_2 - z_{27}\,\mathbf{a}_3 = -x_{27}a\,\hat{\mathbf{x}} - y_{27}b\,\hat{\mathbf{y}} - z_{27}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XVIII}$$

$$\mathbf{B}_{194} = \left(\tfrac{1}{2} + x_{27}\right)\mathbf{a}_1 + y_{27}\,\mathbf{a}_2 - z_{27}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{27}\right)a\,\hat{\mathbf{x}} + y_{27}b\,\hat{\mathbf{y}} - z_{27}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XVIII}$$

$$\mathbf{B}_{195} = \left(\tfrac{1}{2} + x_{27}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{27}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{27}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{27}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{27}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{27}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XVIII}$$

$$\mathbf{B}_{196} = -x_{27}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{27}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{27}\right)\mathbf{a}_3 = -x_{27}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{27}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{27}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{H XVIII}$$

$$\mathbf{B}_{197} = x_{28}\,\mathbf{a}_1 + y_{28}\,\mathbf{a}_2 + z_{28}\,\mathbf{a}_3 = x_{28}a\,\hat{\mathbf{x}} + y_{28}b\,\hat{\mathbf{y}} + z_{28}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{K II}$$

$$\mathbf{B}_{198} = \left(\tfrac{1}{2} - x_{28}\right)\mathbf{a}_1 - y_{28}\,\mathbf{a}_2 + z_{28}\,\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{28}\right)a\,\hat{\mathbf{x}} - y_{28}b\,\hat{\mathbf{y}} + z_{28}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{K II}$$

$$\mathbf{B}_{199} = \left(\tfrac{1}{2} - x_{28}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{28}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{28}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{28}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{28}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{28}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{K II}$$

$$\mathbf{B}_{200} = x_{28}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{28}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{28}\right)\mathbf{a}_3 = x_{28}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{28}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{28}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{K II}$$

$$\mathbf{B}_{201} = -x_{28}\,\mathbf{a}_1 - y_{28}\,\mathbf{a}_2 - z_{28}\,\mathbf{a}_3 = -x_{28}a\,\hat{\mathbf{x}} - y_{28}b\,\hat{\mathbf{y}} - z_{28}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{K II}$$

$$\mathbf{B}_{202} = \left(\tfrac{1}{2} + x_{28}\right)\mathbf{a}_1 + y_{28}\,\mathbf{a}_2 - z_{28}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{28}\right)a\,\hat{\mathbf{x}} + y_{28}b\,\hat{\mathbf{y}} - z_{28}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{K II}$$

$$\mathbf{B}_{203} = \left(\tfrac{1}{2} + x_{28}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{28}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{28}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{28}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{28}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{28}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{K II}$$

$$\mathbf{B}_{204} = -x_{28}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{28}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{28}\right)\mathbf{a}_3 = -x_{28}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{28}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{28}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{K II}$$

$$\mathbf{B}_{205} = x_{29}\,\mathbf{a}_1 + y_{29}\,\mathbf{a}_2 + z_{29}\,\mathbf{a}_3 = x_{29}a\,\hat{\mathbf{x}} + y_{29}b\,\hat{\mathbf{y}} + z_{29}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Mg II}$$

$$\mathbf{B}_{206} = \left(\tfrac{1}{2} - x_{29}\right)\mathbf{a}_1 - y_{29}\,\mathbf{a}_2 + z_{29}\,\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{29}\right)a\,\hat{\mathbf{x}} - y_{29}b\,\hat{\mathbf{y}} + z_{29}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Mg II}$$

$$\mathbf{B}_{207} = \left(\tfrac{1}{2} - x_{29}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{29}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{29}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{29}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{29}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{29}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Mg II}$$

$$\mathbf{B}_{208} = x_{29}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{29}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{29}\right)\mathbf{a}_3 = x_{29}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{29}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{29}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Mg II}$$

$$\mathbf{B}_{209} = -x_{29}\,\mathbf{a}_1 - y_{29}\,\mathbf{a}_2 - z_{29}\,\mathbf{a}_3 = -x_{29}a\,\hat{\mathbf{x}} - y_{29}b\,\hat{\mathbf{y}} - z_{29}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Mg II}$$

$$\mathbf{B}_{210} = \left(\tfrac{1}{2} + x_{29}\right)\mathbf{a}_1 + y_{29}\,\mathbf{a}_2 - z_{29}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{29}\right)a\,\hat{\mathbf{x}} + y_{29}b\,\hat{\mathbf{y}} - z_{29}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Mg II}$$

$$\mathbf{B}_{211} = \left(\tfrac{1}{2} + x_{29}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{29}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{29}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{29}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{29}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{29}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Mg II}$$

$$\mathbf{B}_{212} = -x_{29}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{29}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{29}\right)\mathbf{a}_3 = -x_{29}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{29}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{29}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Mg II}$$

$$\mathbf{B}_{213} = x_{30}\,\mathbf{a}_1 + y_{30}\,\mathbf{a}_2 + z_{30}\,\mathbf{a}_3 = x_{30}a\,\hat{\mathbf{x}} + y_{30}b\,\hat{\mathbf{y}} + z_{30}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O III}$$

$$\mathbf{B}_{214} = \left(\tfrac{1}{2} - x_{30}\right)\mathbf{a}_1 - y_{30}\,\mathbf{a}_2 + z_{30}\,\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{30}\right)a\,\hat{\mathbf{x}} - y_{30}b\,\hat{\mathbf{y}} + z_{30}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O III}$$

$$\mathbf{B}_{215} = \left(\tfrac{1}{2} - x_{30}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{30}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{30}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{30}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{30}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{30}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O III}$$

$$\mathbf{B}_{216} = x_{30}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{30}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{30}\right)\mathbf{a}_3 = x_{30}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{30}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{30}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O III}$$

$$\mathbf{B}_{217} = -x_{30}\,\mathbf{a}_1 - y_{30}\,\mathbf{a}_2 - z_{30}\,\mathbf{a}_3 = -x_{30}a\,\hat{\mathbf{x}} - y_{30}b\,\hat{\mathbf{y}} - z_{30}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O III}$$

$$\mathbf{B}_{218} = \left(\tfrac{1}{2} + x_{30}\right)\mathbf{a}_1 + y_{30}\,\mathbf{a}_2 - z_{30}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{30}\right)a\,\hat{\mathbf{x}} + y_{30}b\,\hat{\mathbf{y}} - z_{30}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O III}$$

$$\mathbf{B}_{219} = \left(\tfrac{1}{2} + x_{30}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{30}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{30}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{30}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{30}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{30}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O III}$$

$$\mathbf{B}_{220} = -x_{30}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{30}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{30}\right)\mathbf{a}_3 = -x_{30}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{30}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{30}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O III}$$

$$\mathbf{B}_{221} = x_{31}\,\mathbf{a}_1 + y_{31}\,\mathbf{a}_2 + z_{31}\,\mathbf{a}_3 = x_{31}a\,\hat{\mathbf{x}} + y_{31}b\,\hat{\mathbf{y}} + z_{31}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O IV}$$

$$\mathbf{B}_{222} = \left(\tfrac{1}{2} - x_{31}\right)\mathbf{a}_1 - y_{31}\,\mathbf{a}_2 + z_{31}\,\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{31}\right)a\,\hat{\mathbf{x}} - y_{31}b\,\hat{\mathbf{y}} + z_{31}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O IV}$$

$$\mathbf{B}_{223} = \left(\tfrac{1}{2} - x_{31}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{31}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{31}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{31}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{31}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{31}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O IV}$$

$$\mathbf{B}_{224} = x_{31}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{31}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{31}\right)\mathbf{a}_3 = x_{31}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{31}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{31}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O IV}$$

$$\mathbf{B}_{225} = -x_{31}\,\mathbf{a}_1 - y_{31}\,\mathbf{a}_2 - z_{31}\,\mathbf{a}_3 = -x_{31}a\,\hat{\mathbf{x}} - y_{31}b\,\hat{\mathbf{y}} - z_{31}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O IV}$$

$$\mathbf{B}_{226} = \left(\tfrac{1}{2} + x_{31}\right)\mathbf{a}_1 + y_{31}\,\mathbf{a}_2 - z_{31}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{31}\right)a\,\hat{\mathbf{x}} + y_{31}b\,\hat{\mathbf{y}} - z_{31}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O IV}$$

$$\mathbf{B}_{227} = \left(\tfrac{1}{2} + x_{31}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{31}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{31}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{31}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{31}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{31}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O IV}$$

$$\mathbf{B}_{228} = -x_{31}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{31}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{31}\right)\mathbf{a}_3 = -x_{31}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{31}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{31}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O IV}$$

$$\mathbf{B}_{229} = x_{32}\,\mathbf{a}_1 + y_{32}\,\mathbf{a}_2 + z_{32}\,\mathbf{a}_3 = x_{32}a\,\hat{\mathbf{x}} + y_{32}b\,\hat{\mathbf{y}} + z_{32}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O V}$$

$$\mathbf{B}_{230} = \left(\tfrac{1}{2} - x_{32}\right)\mathbf{a}_1 - y_{32}\,\mathbf{a}_2 + z_{32}\,\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{32}\right)a\,\hat{\mathbf{x}} - y_{32}b\,\hat{\mathbf{y}} + z_{32}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O V}$$

$$\mathbf{B}_{231} = \left(\tfrac{1}{2} - x_{32}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{32}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{32}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{32}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{32}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{32}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O V}$$

$$\mathbf{B}_{232} = x_{32}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{32}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{32}\right)\mathbf{a}_3 = x_{32}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{32}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{32}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O V}$$

$$\mathbf{B}_{233} = -x_{32}\,\mathbf{a}_1 - y_{32}\,\mathbf{a}_2 - z_{32}\,\mathbf{a}_3 = -x_{32}a\,\hat{\mathbf{x}} - y_{32}b\,\hat{\mathbf{y}} - z_{32}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O V}$$

$$\mathbf{B}_{234} = \left(\tfrac{1}{2} + x_{32}\right)\mathbf{a}_1 + y_{32}\,\mathbf{a}_2 - z_{32}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{32}\right)a\,\hat{\mathbf{x}} + y_{32}b\,\hat{\mathbf{y}} - z_{32}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O V}$$

$$\mathbf{B}_{235} = \left(\tfrac{1}{2} + x_{32}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{32}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{32}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{32}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{32}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{32}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O V}$$

$$\mathbf{B}_{236} = -x_{32}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{32}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{32}\right)\mathbf{a}_3 = -x_{32}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{32}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{32}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O V}$$

$$\mathbf{B}_{237} = x_{33}\,\mathbf{a}_1 + y_{33}\,\mathbf{a}_2 + z_{33}\,\mathbf{a}_3 = x_{33}a\,\hat{\mathbf{x}} + y_{33}b\,\hat{\mathbf{y}} + z_{33}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O VI}$$

$$\mathbf{B}_{238} = \left(\tfrac{1}{2} - x_{33}\right)\mathbf{a}_1 - y_{33}\,\mathbf{a}_2 + z_{33}\,\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{33}\right)a\,\hat{\mathbf{x}} - y_{33}b\,\hat{\mathbf{y}} + z_{33}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O VI}$$

$$\mathbf{B}_{239} = \left(\tfrac{1}{2} - x_{33}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{33}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{33}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{33}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{33}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{33}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O VI}$$

$$\mathbf{B}_{240} = x_{33}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{33}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{33}\right)\mathbf{a}_3 = x_{33}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{33}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{33}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O VI}$$

$$\mathbf{B}_{241} = -x_{33}\,\mathbf{a}_1 - y_{33}\,\mathbf{a}_2 - z_{33}\,\mathbf{a}_3 = -x_{33}a\,\hat{\mathbf{x}} - y_{33}b\,\hat{\mathbf{y}} - z_{33}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O VI}$$

$$\mathbf{B}_{242} = \left(\tfrac{1}{2} + x_{33}\right)\mathbf{a}_1 + y_{33}\,\mathbf{a}_2 - z_{33}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{33}\right)a\,\hat{\mathbf{x}} + y_{33}b\,\hat{\mathbf{y}} - z_{33}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O VI}$$

$$\mathbf{B}_{243} = \left(\tfrac{1}{2} + x_{33}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{33}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{33}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{33}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{33}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{33}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O VI}$$

$$\mathbf{B}_{244} = -x_{33}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{33}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{33}\right)\mathbf{a}_3 = -x_{33}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{33}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{33}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O VI}$$

$$\mathbf{B}_{245} = x_{34}\,\mathbf{a}_1 + y_{34}\,\mathbf{a}_2 + z_{34}\,\mathbf{a}_3 = x_{34}a\,\hat{\mathbf{x}} + y_{34}b\,\hat{\mathbf{y}} + z_{34}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O VII}$$

$$\mathbf{B}_{246} = \left(\tfrac{1}{2} - x_{34}\right)\mathbf{a}_1 - y_{34}\,\mathbf{a}_2 + z_{34}\,\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{34}\right)a\,\hat{\mathbf{x}} - y_{34}b\,\hat{\mathbf{y}} + z_{34}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O VII}$$

$$\mathbf{B}_{247} = \left(\tfrac{1}{2} - x_{34}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{34}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{34}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{34}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{34}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{34}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O VII}$$

$$\mathbf{B}_{248} = x_{34}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{34}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{34}\right)\mathbf{a}_3 = x_{34}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{34}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{34}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O VII}$$

$$\mathbf{B}_{249} = -x_{34}\,\mathbf{a}_1 - y_{34}\,\mathbf{a}_2 - z_{34}\,\mathbf{a}_3 = -x_{34}a\,\hat{\mathbf{x}} - y_{34}b\,\hat{\mathbf{y}} - z_{34}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O VII}$$

$$\mathbf{B}_{250} = \left(\tfrac{1}{2} + x_{34}\right)\mathbf{a}_1 + y_{34}\,\mathbf{a}_2 - z_{34}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{34}\right)a\,\hat{\mathbf{x}} + y_{34}b\,\hat{\mathbf{y}} - z_{34}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O VII}$$

$$\mathbf{B}_{251} = \left(\tfrac{1}{2} + x_{34}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{34}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{34}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{34}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{34}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{34}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O VII}$$

$$\mathbf{B}_{252} = -x_{34}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{34}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{34}\right)\mathbf{a}_3 = -x_{34}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{34}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{34}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{O VII}$$

$\mathbf{B}_{253}$	=	$x_{35}\,\mathbf{a}_1 + y_{35}\,\mathbf{a}_2 + z_{35}\,\mathbf{a}_3$	=	$x_{35}a\,\hat{\mathbf{x}} + y_{35}b\,\hat{\mathbf{y}} + z_{35}c\,\hat{\mathbf{z}}$	(8*e*)	O VIII
$\mathbf{B}_{254}$	=	$\left(\frac{1}{2} - x_{35}\right)\mathbf{a}_1 - y_{35}\,\mathbf{a}_2 + z_{35}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{35}\right)a\,\hat{\mathbf{x}} - y_{35}b\,\hat{\mathbf{y}} + z_{35}c\,\hat{\mathbf{z}}$	(8*e*)	O VIII
$\mathbf{B}_{255}$	=	$\left(\frac{1}{2} - x_{35}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{35}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{35}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{35}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{35}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{35}\right)c\,\hat{\mathbf{z}}$	(8*e*)	O VIII
$\mathbf{B}_{256}$	=	$x_{35}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{35}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{35}\right)\mathbf{a}_3$	=	$x_{35}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{35}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{35}\right)c\,\hat{\mathbf{z}}$	(8*e*)	O VIII
$\mathbf{B}_{257}$	=	$-x_{35}\,\mathbf{a}_1 - y_{35}\,\mathbf{a}_2 - z_{35}\,\mathbf{a}_3$	=	$-x_{35}a\,\hat{\mathbf{x}} - y_{35}b\,\hat{\mathbf{y}} - z_{35}c\,\hat{\mathbf{z}}$	(8*e*)	O VIII
$\mathbf{B}_{258}$	=	$\left(\frac{1}{2} + x_{35}\right)\mathbf{a}_1 + y_{35}\,\mathbf{a}_2 - z_{35}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{35}\right)a\,\hat{\mathbf{x}} + y_{35}b\,\hat{\mathbf{y}} - z_{35}c\,\hat{\mathbf{z}}$	(8*e*)	O VIII
$\mathbf{B}_{259}$	=	$\left(\frac{1}{2} + x_{35}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{35}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{35}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{35}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{35}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{35}\right)c\,\hat{\mathbf{z}}$	(8*e*)	O VIII
$\mathbf{B}_{260}$	=	$-x_{35}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{35}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{35}\right)\mathbf{a}_3$	=	$-x_{35}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{35}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{35}\right)c\,\hat{\mathbf{z}}$	(8*e*)	O VIII
$\mathbf{B}_{261}$	=	$x_{36}\,\mathbf{a}_1 + y_{36}\,\mathbf{a}_2 + z_{36}\,\mathbf{a}_3$	=	$x_{36}a\,\hat{\mathbf{x}} + y_{36}b\,\hat{\mathbf{y}} + z_{36}c\,\hat{\mathbf{z}}$	(8*e*)	O IX
$\mathbf{B}_{262}$	=	$\left(\frac{1}{2} - x_{36}\right)\mathbf{a}_1 - y_{36}\,\mathbf{a}_2 + z_{36}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{36}\right)a\,\hat{\mathbf{x}} - y_{36}b\,\hat{\mathbf{y}} + z_{36}c\,\hat{\mathbf{z}}$	(8*e*)	O IX
$\mathbf{B}_{263}$	=	$\left(\frac{1}{2} - x_{36}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{36}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{36}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{36}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{36}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{36}\right)c\,\hat{\mathbf{z}}$	(8*e*)	O IX
$\mathbf{B}_{264}$	=	$x_{36}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{36}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{36}\right)\mathbf{a}_3$	=	$x_{36}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{36}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{36}\right)c\,\hat{\mathbf{z}}$	(8*e*)	O IX
$\mathbf{B}_{265}$	=	$-x_{36}\,\mathbf{a}_1 - y_{36}\,\mathbf{a}_2 - z_{36}\,\mathbf{a}_3$	=	$-x_{36}a\,\hat{\mathbf{x}} - y_{36}b\,\hat{\mathbf{y}} - z_{36}c\,\hat{\mathbf{z}}$	(8*e*)	O IX
$\mathbf{B}_{266}$	=	$\left(\frac{1}{2} + x_{36}\right)\mathbf{a}_1 + y_{36}\,\mathbf{a}_2 - z_{36}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{36}\right)a\,\hat{\mathbf{x}} + y_{36}b\,\hat{\mathbf{y}} - z_{36}c\,\hat{\mathbf{z}}$	(8*e*)	O IX
$\mathbf{B}_{267}$	=	$\left(\frac{1}{2} + x_{36}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{36}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{36}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{36}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{36}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{36}\right)c\,\hat{\mathbf{z}}$	(8*e*)	O IX
$\mathbf{B}_{268}$	=	$-x_{36}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{36}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{36}\right)\mathbf{a}_3$	=	$-x_{36}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{36}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{36}\right)c\,\hat{\mathbf{z}}$	(8*e*)	O IX
$\mathbf{B}_{269}$	=	$x_{37}\,\mathbf{a}_1 + y_{37}\,\mathbf{a}_2 + z_{37}\,\mathbf{a}_3$	=	$x_{37}a\,\hat{\mathbf{x}} + y_{37}b\,\hat{\mathbf{y}} + z_{37}c\,\hat{\mathbf{z}}$	(8*e*)	O X
$\mathbf{B}_{270}$	=	$\left(\frac{1}{2} - x_{37}\right)\mathbf{a}_1 - y_{37}\,\mathbf{a}_2 + z_{37}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{37}\right)a\,\hat{\mathbf{x}} - y_{37}b\,\hat{\mathbf{y}} + z_{37}c\,\hat{\mathbf{z}}$	(8*e*)	O X
$\mathbf{B}_{271}$	=	$\left(\frac{1}{2} - x_{37}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{37}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{37}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{37}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{37}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{37}\right)c\,\hat{\mathbf{z}}$	(8*e*)	O X
$\mathbf{B}_{272}$	=	$x_{37}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{37}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{37}\right)\mathbf{a}_3$	=	$x_{37}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{37}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{37}\right)c\,\hat{\mathbf{z}}$	(8*e*)	O X
$\mathbf{B}_{273}$	=	$-x_{37}\,\mathbf{a}_1 - y_{37}\,\mathbf{a}_2 - z_{37}\,\mathbf{a}_3$	=	$-x_{37}a\,\hat{\mathbf{x}} - y_{37}b\,\hat{\mathbf{y}} - z_{37}c\,\hat{\mathbf{z}}$	(8*e*)	O X
$\mathbf{B}_{274}$	=	$\left(\frac{1}{2} + x_{37}\right)\mathbf{a}_1 + y_{37}\,\mathbf{a}_2 - z_{37}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{37}\right)a\,\hat{\mathbf{x}} + y_{37}b\,\hat{\mathbf{y}} - z_{37}c\,\hat{\mathbf{z}}$	(8*e*)	O X
$\mathbf{B}_{275}$	=	$\left(\frac{1}{2} + x_{37}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{37}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{37}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{37}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{37}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{37}\right)c\,\hat{\mathbf{z}}$	(8*e*)	O X
$\mathbf{B}_{276}$	=	$-x_{37}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{37}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{37}\right)\mathbf{a}_3$	=	$-x_{37}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{37}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{37}\right)c\,\hat{\mathbf{z}}$	(8*e*)	O X

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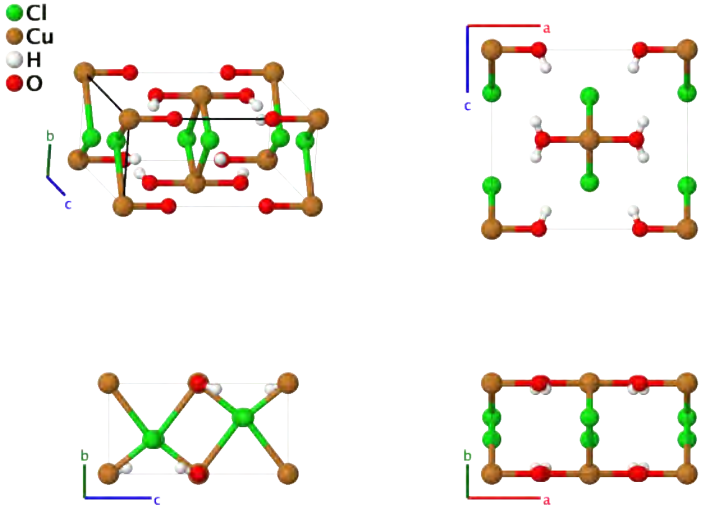
- K. Herrmann, ed., *Strukturbericht Band VII 1939* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1943).

Geometry files:
- CIF: pp. 1607
- POSCAR: pp. 1608

Eriochalcite ($CuCl_2{\cdot}2H_2O$, *C*45) Structure: A2BC4D2_oP18_53_h_a_i_e

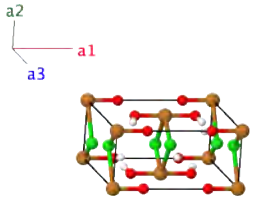
http://aflow.org/prototype-encyclopedia/A2BC4D2_oP18_53_h_a_i_e



Prototype	:	$Cl_2CuH_4O_2$
AFLOW prototype label	:	A2BC4D2_oP18_53_h_a_i_e
***Strukturbericht* designation**	:	*C*45
Pearson symbol	:	oP18
Space group number	:	53
Space group symbol	:	*Pmna*
AFLOW prototype command	:	`aflow --proto=A2BC4D2_oP18_53_h_a_i_e --params=`$a, b/a, c/a, x_2, y_3, z_3, x_4, y_4, z_4$

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(2*a*)	Cu
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(2*a*)	Cu
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1$	=	$x_2 a\,\hat{\mathbf{x}}$	(4*e*)	O

$\mathbf{B}_4$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \frac{1}{2}\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*e*)	O
$\mathbf{B}_5$	=	$-x_2\,\mathbf{a}_1$	=	$-x_2 a\,\hat{\mathbf{x}}$	(4*e*)	O
$\mathbf{B}_6$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \frac{1}{2}\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*e*)	O
$\mathbf{B}_7$	=	$y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*h*)	Cl
$\mathbf{B}_8$	=	$\frac{1}{2}\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(4*h*)	Cl
$\mathbf{B}_9$	=	$\frac{1}{2}\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(4*h*)	Cl
$\mathbf{B}_{10}$	=	$-y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-y_3 b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(4*h*)	Cl
$\mathbf{B}_{11}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*i*)	H
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(8*i*)	H
$\mathbf{B}_{13}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(8*i*)	H
$\mathbf{B}_{14}$	=	$x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*i*)	H
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*i*)	H
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(8*i*)	H
$\mathbf{B}_{17}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(8*i*)	H
$\mathbf{B}_{18}$	=	$-x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*i*)	H

References:

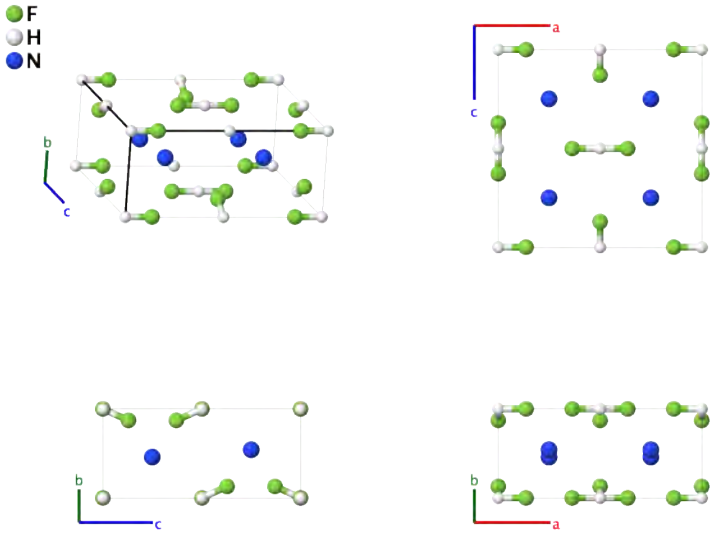
- S. Brownstein, N. F. Han, E. Gabe, and Y. LePage, *A redetermination of the crystal structure of cupric chloride dihydrate*, Zeitschrift für Kristallographie - Crystalline Materials **189**, 13–15 (1989), doi:10.1524/zkri.1989.189.1-2.13.

Geometry files:

- CIF: pp. 1609
- POSCAR: pp. 1610

NH_4HF_2 ($F5_8$) Structure: A2BC_oP16_53_eh_ab_g

http://aflow.org/prototype-encyclopedia/A2BC_oP16_53_eh_ab_g

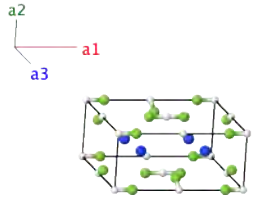


Prototype	:	F_2H_5N
AFLOW prototype label	:	A2BC_oP16_53_eh_ab_g
***Strukturbericht* designation**	:	$F5_8$
Pearson symbol	:	oP16
Space group number	:	53
Space group symbol	:	*Pmna*
AFLOW prototype command	:	`aflow --proto=A2BC_oP16_53_eh_ab_g --params=`$a, b/a, c/a, x_3, y_4, y_5, z_5$

- This structure was first investigated by (Pauling, 1933) and assigned *Strukturbericht* designation $F5_8$ by (Gottfried, 1937). It was reinvestigated by (Rogers, 1940). Neither paper notes the positions of the hydrogen atoms, but under the assumption that the structure is similar to KHF_2 ($F5_2$), (Downs, 2003) puts some of them between pairs of fluorine atoms. The remaining hydrogen atoms are part of the NH_4 radical.
- The crystal structure was given in the *Pman* setting of space group #53. We used FINDSYM to change it to the standard *Pmna* structure.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(2*a*)	H I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(2*a*)	H I
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	(2*b*)	H II
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	(2*b*)	H II
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1$	=	$x_3 a\,\hat{\mathbf{x}}$	(4*e*)	F I
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*e*)	F I
$\mathbf{B}_7$	=	$-x_3\,\mathbf{a}_1$	=	$-x_3 a\,\hat{\mathbf{x}}$	(4*e*)	F I
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*e*)	F I
$\mathbf{B}_9$	=	$\frac{1}{4}\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*g*)	NH_4
$\mathbf{B}_{10}$	=	$\frac{1}{4}\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(4*g*)	NH_4
$\mathbf{B}_{11}$	=	$\frac{3}{4}\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(4*g*)	NH_4
$\mathbf{B}_{12}$	=	$\frac{3}{4}\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*g*)	NH_4
$\mathbf{B}_{13}$	=	$y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(4*h*)	F II
$\mathbf{B}_{14}$	=	$\frac{1}{2}\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	(4*h*)	F II
$\mathbf{B}_{15}$	=	$\frac{1}{2}\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	(4*h*)	F II
$\mathbf{B}_{16}$	=	$-y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-y_5 b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(4*h*)	F II

References:

- M. T. Rogers and L. Helmholz, *A Redetermination of the Parameters in Ammonium Bifluoride*, J. Am. Chem. Soc. **62**, 1533–1536 (1940), doi:10.1021/ja01863a057.
- L. Pauling, *The Crystal Structure of Ammonium Hydrogen Fluoride, NH_4HF_2*, Zeitschrift für Kristallographie - Crystalline Materials **85**, 380–391 (1933), doi:10.1524/zkri.1933.85.1.380.
- C. Gottfried and F. Schossberger, eds., *Strukturbericht Band III 1933-1935* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Found in:

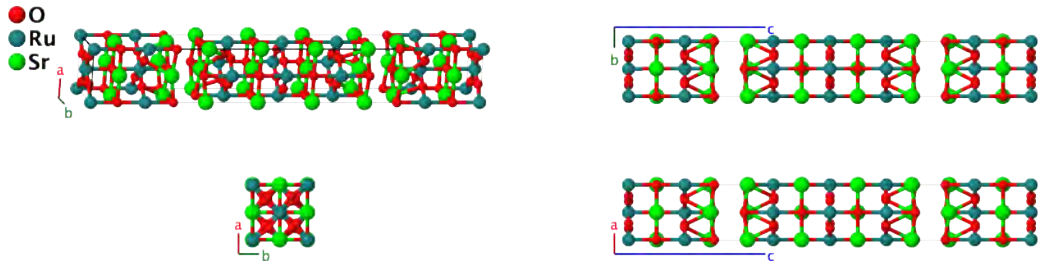
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1610
- POSCAR: pp. 1610

Orthorhombic $Sr_4Ru_3O_{10}$ Structure: A10B3C4_oP68_55_2e2fgh2i_adef_2e2f

http://aflow.org/prototype-encyclopedia/A10B3C4_oP68_55_2e2fgh2i_adef_2e2f

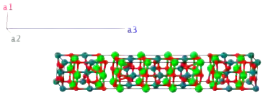


Prototype	:	$O_{10}Ru_3Sr_4$
AFLOW prototype label	:	A10B3C4_oP68_55_2e2fgh2i_adef_2e2f
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP68
Space group number	:	55
Space group symbol	:	*Pbam*
AFLOW prototype command	:	`aflow --proto=A10B3C4_oP68_55_2e2fgh2i_adef_2e2f --params=`$a, b/a, c/a, z_3, z_4, z_5, z_6, z_7, z_8, z_9, z_{10}, z_{11}, z_{12}, x_{13}, y_{13}, x_{14}, y_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}$

- This structure consists of triple-layer ruthenate structures separated by 2.37 Å from each other. In the *Pbam* #55 space group shown here there are two inequivalent stacks in the orthorhombic cell.
- This cell is very problematic. (Crawford, 2002) note that the x-ray scattering intensities are pseudo body-centered, but found that refining this structure in a body-centered cell with space group *Bbcm* (*Cmca* #64 in our standard orientation) led to non-positive definite thermal parameters. However, if we allow a lattice and atom position uncertainty of only 0.01 Å in the atomic coordinates, AFLOW-SYM and FINDSYM place the structure in base-centered space group *Cmca* #64, oC68. In that case there is only one triple-layer stack in the primitive cell, and the two stacks in the conventional orthorhombic cell are equivalent.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	Ru I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$	$(2a)$	Ru I
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(2d)$	Ru II
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(2d)$	Ru II
$\mathbf{B}_5$	=	$z_3\,\mathbf{a}_3$	=	$z_3 c\,\hat{\mathbf{z}}$	$(4e)$	O I

$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_7$	=	$-z_3\,\mathbf{a}_3$	=	$-z_3 c\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_8$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(4e)$	O I
$\mathbf{B}_9$	=	$z_4\,\mathbf{a}_3$	=	$z_4 c\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_{10}$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_{11}$	=	$-z_4\,\mathbf{a}_3$	=	$-z_4 c\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_{12}$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_{13}$	=	$z_5\,\mathbf{a}_3$	=	$z_5 c\,\hat{\mathbf{z}}$	$(4e)$	Ru III
$\mathbf{B}_{14}$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	$(4e)$	Ru III
$\mathbf{B}_{15}$	=	$-z_5\,\mathbf{a}_3$	=	$-z_5 c\,\hat{\mathbf{z}}$	$(4e)$	Ru III
$\mathbf{B}_{16}$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(4e)$	Ru III
$\mathbf{B}_{17}$	=	$z_6\,\mathbf{a}_3$	=	$z_6 c\,\hat{\mathbf{z}}$	$(4e)$	Sr I
$\mathbf{B}_{18}$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	$(4e)$	Sr I
$\mathbf{B}_{19}$	=	$-z_6\,\mathbf{a}_3$	=	$-z_6 c\,\hat{\mathbf{z}}$	$(4e)$	Sr I
$\mathbf{B}_{20}$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(4e)$	Sr I
$\mathbf{B}_{21}$	=	$z_7\,\mathbf{a}_3$	=	$z_7 c\,\hat{\mathbf{z}}$	$(4e)$	Sr II
$\mathbf{B}_{22}$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	$(4e)$	Sr II
$\mathbf{B}_{23}$	=	$-z_7\,\mathbf{a}_3$	=	$-z_7 c\,\hat{\mathbf{z}}$	$(4e)$	Sr II
$\mathbf{B}_{24}$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	$(4e)$	Sr II
$\mathbf{B}_{25}$	=	$\frac{1}{2}\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$\frac{1}{2}b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	$(4f)$	O III
$\mathbf{B}_{26}$	=	$\frac{1}{2}\,\mathbf{a}_1 - z_8\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - z_8 c\,\hat{\mathbf{z}}$	$(4f)$	O III
$\mathbf{B}_{27}$	=	$\frac{1}{2}\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$\frac{1}{2}b\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	$(4f)$	O III
$\mathbf{B}_{28}$	=	$\frac{1}{2}\,\mathbf{a}_1 + z_8\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + z_8 c\,\hat{\mathbf{z}}$	$(4f)$	O III
$\mathbf{B}_{29}$	=	$\frac{1}{2}\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$\frac{1}{2}b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	$(4f)$	O IV
$\mathbf{B}_{30}$	=	$\frac{1}{2}\,\mathbf{a}_1 - z_9\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - z_9 c\,\hat{\mathbf{z}}$	$(4f)$	O IV
$\mathbf{B}_{31}$	=	$\frac{1}{2}\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$\frac{1}{2}b\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}}$	$(4f)$	O IV
$\mathbf{B}_{32}$	=	$\frac{1}{2}\,\mathbf{a}_1 + z_9\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + z_9 c\,\hat{\mathbf{z}}$	$(4f)$	O IV
$\mathbf{B}_{33}$	=	$\frac{1}{2}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$\frac{1}{2}b\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}}$	$(4f)$	Ru IV
$\mathbf{B}_{34}$	=	$\frac{1}{2}\,\mathbf{a}_1 - z_{10}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - z_{10} c\,\hat{\mathbf{z}}$	$(4f)$	Ru IV
$\mathbf{B}_{35}$	=	$\frac{1}{2}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3$	=	$\frac{1}{2}b\,\hat{\mathbf{y}} - z_{10} c\,\hat{\mathbf{z}}$	$(4f)$	Ru IV
$\mathbf{B}_{36}$	=	$\frac{1}{2}\,\mathbf{a}_1 + z_{10}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + z_{10} c\,\hat{\mathbf{z}}$	$(4f)$	Ru IV
$\mathbf{B}_{37}$	=	$\frac{1}{2}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$\frac{1}{2}b\,\hat{\mathbf{y}} + z_{11} c\,\hat{\mathbf{z}}$	$(4f)$	Sr III
$\mathbf{B}_{38}$	=	$\frac{1}{2}\,\mathbf{a}_1 - z_{11}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - z_{11} c\,\hat{\mathbf{z}}$	$(4f)$	Sr III
$\mathbf{B}_{39}$	=	$\frac{1}{2}\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3$	=	$\frac{1}{2}b\,\hat{\mathbf{y}} - z_{11} c\,\hat{\mathbf{z}}$	$(4f)$	Sr III
$\mathbf{B}_{40}$	=	$\frac{1}{2}\,\mathbf{a}_1 + z_{11}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + z_{11} c\,\hat{\mathbf{z}}$	$(4f)$	Sr III
$\mathbf{B}_{41}$	=	$\frac{1}{2}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3$	=	$\frac{1}{2}b\,\hat{\mathbf{y}} + z_{12} c\,\hat{\mathbf{z}}$	$(4f)$	Sr IV

$\mathbf{B}_{42}$	=	$\frac{1}{2}\,\mathbf{a}_1 - z_{12}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - z_{12}c\,\hat{\mathbf{z}}$	$(4f)$	Sr IV
$\mathbf{B}_{43}$	=	$\frac{1}{2}\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3$	=	$\frac{1}{2}b\,\hat{\mathbf{y}} - z_{12}c\,\hat{\mathbf{z}}$	$(4f)$	Sr IV
$\mathbf{B}_{44}$	=	$\frac{1}{2}\,\mathbf{a}_1 + z_{12}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + z_{12}c\,\hat{\mathbf{z}}$	$(4f)$	Sr IV
$\mathbf{B}_{45}$	=	$x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2$	=	$x_{13}a\,\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}}$	$(4g)$	O V
$\mathbf{B}_{46}$	=	$-x_{13}\,\mathbf{a}_1 - y_{13}\,\mathbf{a}_2$	=	$-x_{13}a\,\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}}$	$(4g)$	O V
$\mathbf{B}_{47}$	=	$\left(\frac{1}{2} - x_{13}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{13}\right)\mathbf{a}_2$	=	$\left(\frac{1}{2} - x_{13}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{13}\right)b\,\hat{\mathbf{y}}$	$(4g)$	O V
$\mathbf{B}_{48}$	=	$\left(\frac{1}{2} + x_{13}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{13}\right)\mathbf{a}_2$	=	$\left(\frac{1}{2} + x_{13}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{13}\right)b\,\hat{\mathbf{y}}$	$(4g)$	O V
$\mathbf{B}_{49}$	=	$x_{14}\,\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$x_{14}a\,\hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(4h)$	O VI
$\mathbf{B}_{50}$	=	$-x_{14}\,\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_{14}a\,\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(4h)$	O VI
$\mathbf{B}_{51}$	=	$\left(\frac{1}{2} - x_{14}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{14}\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{14}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{14}\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(4h)$	O VI
$\mathbf{B}_{52}$	=	$\left(\frac{1}{2} + x_{14}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{14}\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{14}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{14}\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(4h)$	O VI
$\mathbf{B}_{53}$	=	$x_{15}\,\mathbf{a}_1 + y_{15}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3$	=	$x_{15}a\,\hat{\mathbf{x}} + y_{15}b\,\hat{\mathbf{y}} + z_{15}c\,\hat{\mathbf{z}}$	$(8i)$	O VII
$\mathbf{B}_{54}$	=	$-x_{15}\,\mathbf{a}_1 - y_{15}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3$	=	$-x_{15}a\,\hat{\mathbf{x}} - y_{15}b\,\hat{\mathbf{y}} + z_{15}c\,\hat{\mathbf{z}}$	$(8i)$	O VII
$\mathbf{B}_{55}$	=	$\left(\frac{1}{2} - x_{15}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{15}\right)\mathbf{a}_2 - z_{15}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{15}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{15}\right)b\,\hat{\mathbf{y}} - z_{15}c\,\hat{\mathbf{z}}$	$(8i)$	O VII
$\mathbf{B}_{56}$	=	$\left(\frac{1}{2} + x_{15}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{15}\right)\mathbf{a}_2 - z_{15}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{15}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{15}\right)b\,\hat{\mathbf{y}} - z_{15}c\,\hat{\mathbf{z}}$	$(8i)$	O VII
$\mathbf{B}_{57}$	=	$-x_{15}\,\mathbf{a}_1 - y_{15}\,\mathbf{a}_2 - z_{15}\,\mathbf{a}_3$	=	$-x_{15}a\,\hat{\mathbf{x}} - y_{15}b\,\hat{\mathbf{y}} - z_{15}c\,\hat{\mathbf{z}}$	$(8i)$	O VII
$\mathbf{B}_{58}$	=	$x_{15}\,\mathbf{a}_1 + y_{15}\,\mathbf{a}_2 - z_{15}\,\mathbf{a}_3$	=	$x_{15}a\,\hat{\mathbf{x}} + y_{15}b\,\hat{\mathbf{y}} - z_{15}c\,\hat{\mathbf{z}}$	$(8i)$	O VII
$\mathbf{B}_{59}$	=	$\left(\frac{1}{2} + x_{15}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{15}\right)\mathbf{a}_2 + z_{15}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{15}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{15}\right)b\,\hat{\mathbf{y}} + z_{15}c\,\hat{\mathbf{z}}$	$(8i)$	O VII
$\mathbf{B}_{60}$	=	$\left(\frac{1}{2} - x_{15}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{15}\right)\mathbf{a}_2 + z_{15}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{15}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{15}\right)b\,\hat{\mathbf{y}} + z_{15}c\,\hat{\mathbf{z}}$	$(8i)$	O VII
$\mathbf{B}_{61}$	=	$x_{16}\,\mathbf{a}_1 + y_{16}\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3$	=	$x_{16}a\,\hat{\mathbf{x}} + y_{16}b\,\hat{\mathbf{y}} + z_{16}c\,\hat{\mathbf{z}}$	$(8i)$	O VIII
$\mathbf{B}_{62}$	=	$-x_{16}\,\mathbf{a}_1 - y_{16}\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3$	=	$-x_{16}a\,\hat{\mathbf{x}} - y_{16}b\,\hat{\mathbf{y}} + z_{16}c\,\hat{\mathbf{z}}$	$(8i)$	O VIII
$\mathbf{B}_{63}$	=	$\left(\frac{1}{2} - x_{16}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{16}\right)\mathbf{a}_2 - z_{16}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{16}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{16}\right)b\,\hat{\mathbf{y}} - z_{16}c\,\hat{\mathbf{z}}$	$(8i)$	O VIII
$\mathbf{B}_{64}$	=	$\left(\frac{1}{2} + x_{16}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{16}\right)\mathbf{a}_2 - z_{16}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{16}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{16}\right)b\,\hat{\mathbf{y}} - z_{16}c\,\hat{\mathbf{z}}$	$(8i)$	O VIII
$\mathbf{B}_{65}$	=	$-x_{16}\,\mathbf{a}_1 - y_{16}\,\mathbf{a}_2 - z_{16}\,\mathbf{a}_3$	=	$-x_{16}a\,\hat{\mathbf{x}} - y_{16}b\,\hat{\mathbf{y}} - z_{16}c\,\hat{\mathbf{z}}$	$(8i)$	O VIII
$\mathbf{B}_{66}$	=	$x_{16}\,\mathbf{a}_1 + y_{16}\,\mathbf{a}_2 - z_{16}\,\mathbf{a}_3$	=	$x_{16}a\,\hat{\mathbf{x}} + y_{16}b\,\hat{\mathbf{y}} - z_{16}c\,\hat{\mathbf{z}}$	$(8i)$	O VIII
$\mathbf{B}_{67}$	=	$\left(\frac{1}{2} + x_{16}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{16}\right)\mathbf{a}_2 + z_{16}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{16}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{16}\right)b\,\hat{\mathbf{y}} + z_{16}c\,\hat{\mathbf{z}}$	$(8i)$	O VIII
$\mathbf{B}_{68}$	=	$\left(\frac{1}{2} - x_{16}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{16}\right)\mathbf{a}_2 + z_{16}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{16}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{16}\right)b\,\hat{\mathbf{y}} + z_{16}c\,\hat{\mathbf{z}}$	$(8i)$	O VIII

References:

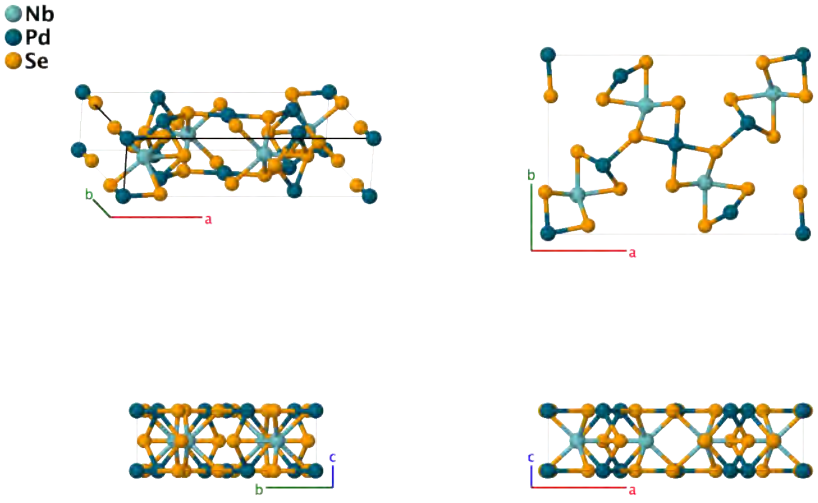
- M. K. Crawford, R. L. Harlow, W. Marshall, Z. Li, G. Cao, R. L. Lindstrom, Q. Huang, and J. W. Lynn, *Structure and magnetism of single crystal $Sr_4Ru_3O_{10}$: A ferromagnetic triple-layer ruthenate*, Phys. Rev. B **65**, 214412 (2002), doi:10.1103/PhysRevB.65.214412.

Geometry files:

- CIF: pp. 1610
- POSCAR: pp. 1611

$Nb_2Pd_3Se_8$ Structure: A2B3C8_oP26_55_h_ag_2g2h

http://aflow.org/prototype-encyclopedia/A2B3C8_oP26_55_h_ag_2g2h



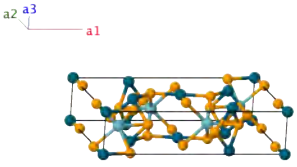
Prototype	:	$Nb_2Pd_3Se_8$
AFLOW prototype label	:	A2B3C8_oP26_55_h_ag_2g2h
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP26
Space group number	:	55
Space group symbol	:	*Pbam*
AFLOW prototype command	:	`aflow --proto=A2B3C8_oP26_55_h_ag_2g2h --params=`$a, b/a, c/a, x_2, y_2, x_3, y_3, x_4, y_4, x_5, y_5, x_6, y_6, x_7, y_7$

Other compounds with this structure

- $Ta_2Pd_3Se_8$

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	Pd I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$	$(2a)$	Pd I
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2$	=	$x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}}$	$(4g)$	Pd II

$\mathbf{B}_4$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2$	=	$-x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}}$	$(4g)$	Pd II
$\mathbf{B}_5$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2$	=	$\left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right) b\,\hat{\mathbf{y}}$	$(4g)$	Pd II
$\mathbf{B}_6$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2$	=	$\left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right) b\,\hat{\mathbf{y}}$	$(4g)$	Pd II
$\mathbf{B}_7$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}}$	$(4g)$	Se I
$\mathbf{B}_8$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}}$	$(4g)$	Se I
$\mathbf{B}_9$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2$	=	$\left(\frac{1}{2} - x_3\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right) b\,\hat{\mathbf{y}}$	$(4g)$	Se I
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2$	=	$\left(\frac{1}{2} + x_3\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right) b\,\hat{\mathbf{y}}$	$(4g)$	Se I
$\mathbf{B}_{11}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}}$	$(4g)$	Se II
$\mathbf{B}_{12}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}}$	$(4g)$	Se II
$\mathbf{B}_{13}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2$	=	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right) b\,\hat{\mathbf{y}}$	$(4g)$	Se II
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2$	=	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right) b\,\hat{\mathbf{y}}$	$(4g)$	Se II
$\mathbf{B}_{15}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \frac{1}{2} c\,\hat{\mathbf{z}}$	$(4h)$	Nb
$\mathbf{B}_{16}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \frac{1}{2} c\,\hat{\mathbf{z}}$	$(4h)$	Nb
$\mathbf{B}_{17}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right) b\,\hat{\mathbf{y}} + \frac{1}{2} c\,\hat{\mathbf{z}}$	$(4h)$	Nb
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right) b\,\hat{\mathbf{y}} + \frac{1}{2} c\,\hat{\mathbf{z}}$	$(4h)$	Nb
$\mathbf{B}_{19}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \frac{1}{2} c\,\hat{\mathbf{z}}$	$(4h)$	Se III
$\mathbf{B}_{20}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \frac{1}{2} c\,\hat{\mathbf{z}}$	$(4h)$	Se III
$\mathbf{B}_{21}$	=	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right) b\,\hat{\mathbf{y}} + \frac{1}{2} c\,\hat{\mathbf{z}}$	$(4h)$	Se III
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right) b\,\hat{\mathbf{y}} + \frac{1}{2} c\,\hat{\mathbf{z}}$	$(4h)$	Se III
$\mathbf{B}_{23}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + \frac{1}{2} c\,\hat{\mathbf{z}}$	$(4h)$	Se IV
$\mathbf{B}_{24}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \frac{1}{2} c\,\hat{\mathbf{z}}$	$(4h)$	Se IV
$\mathbf{B}_{25}$	=	$\left(\frac{1}{2} - x_7\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_7\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right) b\,\hat{\mathbf{y}} + \frac{1}{2} c\,\hat{\mathbf{z}}$	$(4h)$	Se IV
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} + x_7\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_7\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_7\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_7\right) b\,\hat{\mathbf{y}} + \frac{1}{2} c\,\hat{\mathbf{z}}$	$(4h)$	Se IV

References:

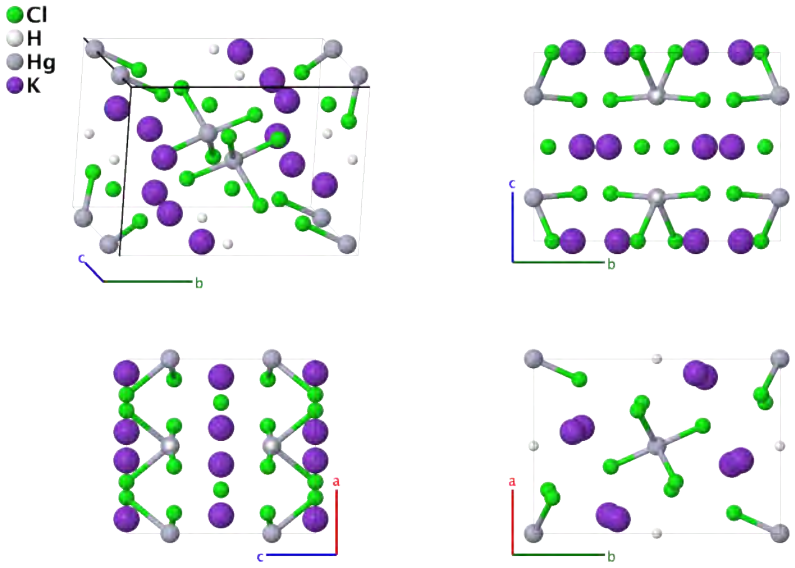
- D. A. Keszler and J. A. Ibers, *A new structural type in ternary chalcogenide chemistry: Structure and properties of* $Nb_2Pd_3Se_8$, J. Solid State Chem. **52**, 73–79 (1984), doi:10.1016/0022-4596(84)90200-7.

Geometry files:

- CIF: pp. 1611
- POSCAR: pp. 1611

$K_2HgCl_4{\cdot}H_2O$ ($E3_4$) Structure: A4BCD2_oP32_55_ghi_f_e_gh

http://aflow.org/prototype-encyclopedia/A4BCD2_oP32_55_ghi_f_e_gh

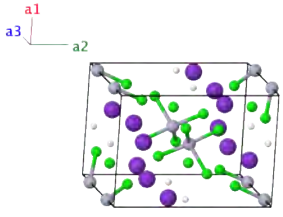


Prototype	:	$Cl_4(H_2O)HgK_2$
AFLOW prototype label	:	A4BCD2_oP32_55_ghi_f_e_gh
***Strukturbericht* designation**	:	$E3_4$
Pearson symbol	:	oP32
Space group number	:	55
Space group symbol	:	*Pbam*
AFLOW prototype command	:	`aflow --proto=A4BCD2_oP32_55_ghi_f_e_gh --params=`$a, b/a, c/a, z_1, z_2, x_3, y_3, x_4, y_4, x_5, y_5, x_6, y_6, x_7, y_7, z_7$

- The positions of the hydrogen atoms in the water molecules have not been determined.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$z_1\,\mathbf{a}_3$	=	$z_1 c\,\hat{\mathbf{z}}$	($4e$)	Hg
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	($4e$)	Hg

$\mathbf{B}_3$	=	$-z_1\,\mathbf{a}_3$	=	$-z_1 c\,\hat{\mathbf{z}}$	$(4e)$	Hg
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	$(4e)$	Hg
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(4f)$	H_2O
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_1 - z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - z_2 c\,\hat{\mathbf{z}}$	$(4f)$	H_2O
$\mathbf{B}_7$	=	$\frac{1}{2}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\frac{1}{2}b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	$(4f)$	H_2O
$\mathbf{B}_8$	=	$\frac{1}{2}\,\mathbf{a}_1 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + z_2 c\,\hat{\mathbf{z}}$	$(4f)$	H_2O
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}}$	$(4g)$	Cl I
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}}$	$(4g)$	Cl I
$\mathbf{B}_{11}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)b\,\hat{\mathbf{y}}$	$(4g)$	Cl I
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)b\,\hat{\mathbf{y}}$	$(4g)$	Cl I
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}}$	$(4g)$	K I
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}}$	$(4g)$	K I
$\mathbf{B}_{15}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)b\,\hat{\mathbf{y}}$	$(4g)$	K I
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)b\,\hat{\mathbf{y}}$	$(4g)$	K I
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(4h)$	Cl II
$\mathbf{B}_{18}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(4h)$	Cl II
$\mathbf{B}_{19}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(4h)$	Cl II
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(4h)$	Cl II
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(4h)$	K II
$\mathbf{B}_{22}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(4h)$	K II
$\mathbf{B}_{23}$	=	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(4h)$	K II
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(4h)$	K II
$\mathbf{B}_{25}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	$(8i)$	Cl III
$\mathbf{B}_{26}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	$(8i)$	Cl III
$\mathbf{B}_{27}$	=	$\left(\frac{1}{2} - x_7\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right)b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	$(8i)$	Cl III
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} + x_7\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_7\right)\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_7\right)b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	$(8i)$	Cl III
$\mathbf{B}_{29}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	$(8i)$	Cl III
$\mathbf{B}_{30}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	$(8i)$	Cl III
$\mathbf{B}_{31}$	=	$\left(\frac{1}{2} + x_7\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_7\right)\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_7\right)b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	$(8i)$	Cl III
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2} - x_7\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right)b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	$(8i)$	Cl III

References:

- K. Aurivillius and C. Stålhandske, *An X-Ray Single Crystal Study of $K_2HgCl_4{\cdot}H_2O$*, Acta Chem. Scand. **27**, 1086–1088 (1973), doi:10.3891/acta.chem.scand.27-1086.

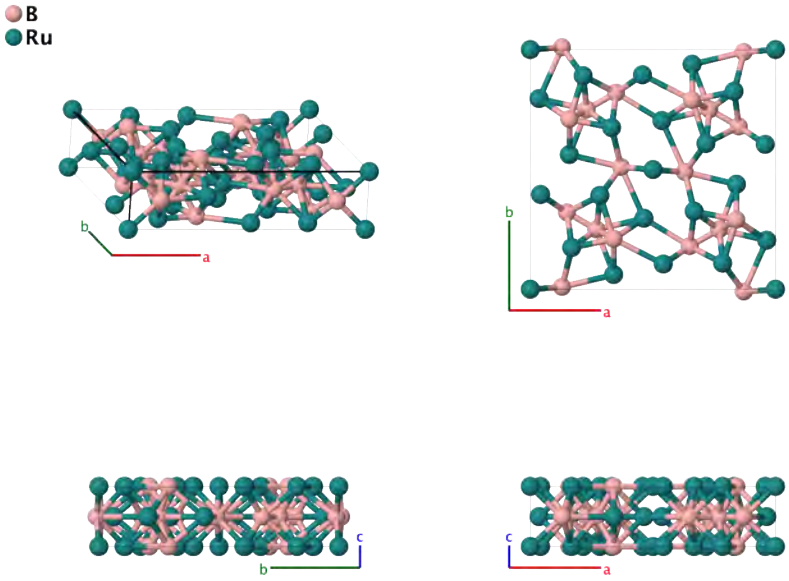
Geometry files:

- CIF: pp. 1612

- POSCAR: pp. 1612

$Ru_{11}B_8$ Structure: A8B11_oP38_55_g3h_a3g2h

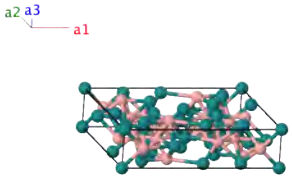
http://aflow.org/prototype-encyclopedia/A8B11_oP38_55_g3h_a3g2h



Prototype	:	B_8Ru_{11}
AFLOW prototype label	:	A8B11_oP38_55_g3h_a3g2h
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP38
Space group number	:	55
Space group symbol	:	*Pbam*
AFLOW prototype command	:	`aflow --proto=A8B11_oP38_55_g3h_a3g2h --params=`$a, b/a, c/a, x_2, y_2, x_3, y_3, x_4, y_4, x_5, y_5, x_6, y_6, x_7, y_7, x_8, y_8, x_9, y_9, x_{10}, y_{10}$

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$ =	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($2a$)	Ru I

$$
\begin{array}{lcccccc}
\mathbf{B}_{2} & = & \frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 & = & \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} & (2a) & \text{Ru I} \\
\mathbf{B}_{3} & = & x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 & = & x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} & (4g) & \text{B I} \\
\mathbf{B}_{4} & = & -x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 & = & -x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} & (4g) & \text{B I} \\
\mathbf{B}_{5} & = & \left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 & = & \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right)b\,\hat{\mathbf{y}} & (4g) & \text{B I} \\
\mathbf{B}_{6} & = & \left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 & = & \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right)b\,\hat{\mathbf{y}} & (4g) & \text{B I} \\
\mathbf{B}_{7} & = & x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 & = & x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} & (4g) & \text{Ru II} \\
\mathbf{B}_{8} & = & -x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 & = & -x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} & (4g) & \text{Ru II} \\
\mathbf{B}_{9} & = & \left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 & = & \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)b\,\hat{\mathbf{y}} & (4g) & \text{Ru II} \\
\mathbf{B}_{10} & = & \left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 & = & \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)b\,\hat{\mathbf{y}} & (4g) & \text{Ru II} \\
\mathbf{B}_{11} & = & x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 & = & x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} & (4g) & \text{Ru III} \\
\mathbf{B}_{12} & = & -x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 & = & -x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} & (4g) & \text{Ru III} \\
\mathbf{B}_{13} & = & \left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 & = & \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)b\,\hat{\mathbf{y}} & (4g) & \text{Ru III} \\
\mathbf{B}_{14} & = & \left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 & = & \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)b\,\hat{\mathbf{y}} & (4g) & \text{Ru III} \\
\mathbf{B}_{15} & = & x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 & = & x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} & (4g) & \text{Ru IV} \\
\mathbf{B}_{16} & = & -x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 & = & -x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} & (4g) & \text{Ru IV} \\
\mathbf{B}_{17} & = & \left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 & = & \left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right)b\,\hat{\mathbf{y}} & (4g) & \text{Ru IV} \\
\mathbf{B}_{18} & = & \left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 & = & \left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right)b\,\hat{\mathbf{y}} & (4g) & \text{Ru IV} \\
\mathbf{B}_{19} & = & x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3 & = & x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}} & (4h) & \text{B II} \\
\mathbf{B}_{20} & = & -x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3 & = & -x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}} & (4h) & \text{B II} \\
\mathbf{B}_{21} & = & \left(\frac{1}{2} - x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3 & = & \left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}} & (4h) & \text{B II} \\
\mathbf{B}_{22} & = & \left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3 & = & \left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}} & (4h) & \text{B II} \\
\mathbf{B}_{23} & = & x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3 & = & x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}} & (4h) & \text{B III} \\
\mathbf{B}_{24} & = & -x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3 & = & -x_7 a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}} & (4h) & \text{B III} \\
\mathbf{B}_{25} & = & \left(\frac{1}{2} - x_7\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3 & = & \left(\frac{1}{2} - x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}} & (4h) & \text{B III} \\
\mathbf{B}_{26} & = & \left(\frac{1}{2} + x_7\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_7\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3 & = & \left(\frac{1}{2} + x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_7\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}} & (4h) & \text{B III} \\
\mathbf{B}_{27} & = & x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3 & = & x_8 a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}} & (4h) & \text{B IV} \\
\mathbf{B}_{28} & = & -x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3 & = & -x_8 a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}} & (4h) & \text{B IV} \\
\mathbf{B}_{29} & = & \left(\frac{1}{2} - x_8\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_8\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3 & = & \left(\frac{1}{2} - x_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_8\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}} & (4h) & \text{B IV} \\
\mathbf{B}_{30} & = & \left(\frac{1}{2} + x_8\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_8\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3 & = & \left(\frac{1}{2} + x_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_8\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}} & (4h) & \text{B IV} \\
\mathbf{B}_{31} & = & x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3 & = & x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}} & (4h) & \text{Ru V} \\
\mathbf{B}_{32} & = & -x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3 & = & -x_9 a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}} & (4h) & \text{Ru V} \\
\mathbf{B}_{33} & = & \left(\frac{1}{2} - x_9\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_9\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3 & = & \left(\frac{1}{2} - x_9\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_9\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}} & (4h) & \text{Ru V} \\
\mathbf{B}_{34} & = & \left(\frac{1}{2} + x_9\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_9\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3 & = & \left(\frac{1}{2} + x_9\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_9\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}} & (4h) & \text{Ru V} \\
\mathbf{B}_{35} & = & x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3 & = & x_{10} a\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}} & (4h) & \text{Ru VI} \\
\mathbf{B}_{36} & = & -x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3 & = & -x_{10} a\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}} & (4h) & \text{Ru VI} \\
\mathbf{B}_{37} & = & \left(\frac{1}{2} - x_{10}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{10}\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3 & = & \left(\frac{1}{2} - x_{10}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{10}\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}} & (4h) & \text{Ru VI}
\end{array}
$$

$$\mathbf{B}_{38} = \left(\tfrac{1}{2} + x_{10}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{10}\right)\mathbf{a}_2 + \tfrac{1}{2}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{10}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{10}\right)b\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}} \quad (4h) \quad \text{Ru VI}$$

References:

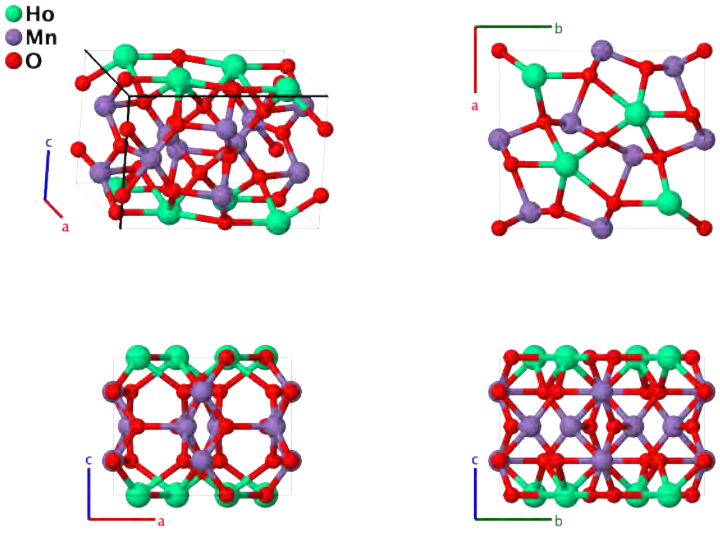
- J. Åselius, *The Crystal Structure of* $Ru_{11}B_8$, Acta Chem. Scand. **14**, 2169–2176 (1960), doi:10.3891/acta.chem.scand.14-2169.

Geometry files:

- CIF: pp. 1612
- POSCAR: pp. 1613

$HoMn_2O_5$ Structure: AB2C5_oP32_55_g_fh_eghi

http://aflow.org/prototype-encyclopedia/AB2C5_oP32_55_g_fh_eghi



Prototype	:	$HoMn_2O_5$
AFLOW prototype label	:	AB2C5_oP32_55_g_fh_eghi
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP32
Space group number	:	55
Space group symbol	:	*Pbam*
AFLOW prototype command	:	`aflow --proto=AB2C5_oP32_55_g_fh_eghi --params=`$a, b/a, c/a, z_1, z_2, x_3, y_3, x_4, y_4, x_5, y_5, x_6, y_6, x_7, y_7, z_7$

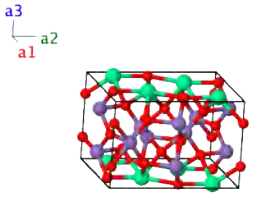
Other compounds with this structure

- $DyMn_2O_5$, $ErMn_2O_5$, $EuMn_2O_5$, $LaMn_2O_5$, $NdMn_2O_5$, $PrMn_2O_5$, $SmMn_2O_5$, and $TbMn_2O_5$

- We found no definitive definition for the prototype of the structure XMn_2O_5, where X is a rare earth metal. (Quezel-Ambrunaz, 1964) has the earliest description of the structure we could find, so we use $HoMn_2O_5$ as the prototype.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$\mathbf{B}_{1}$	=	$z_1\,\mathbf{a}_3$	=	$z_1 c\,\hat{\mathbf{z}}$	($4e$)	O I
$\mathbf{B}_{2}$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	($4e$)	O I
$\mathbf{B}_{3}$	=	$-z_1\,\mathbf{a}_3$	=	$-z_1 c\,\hat{\mathbf{z}}$	($4e$)	O I
$\mathbf{B}_{4}$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	($4e$)	O I
$\mathbf{B}_{5}$	=	$\frac{1}{2}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	($4f$)	Mn I
$\mathbf{B}_{6}$	=	$\frac{1}{2}\,\mathbf{a}_1 - z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - z_2 c\,\hat{\mathbf{z}}$	($4f$)	Mn I
$\mathbf{B}_{7}$	=	$\frac{1}{2}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\frac{1}{2}b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	($4f$)	Mn I
$\mathbf{B}_{8}$	=	$\frac{1}{2}\,\mathbf{a}_1 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + z_2 c\,\hat{\mathbf{z}}$	($4f$)	Mn I
$\mathbf{B}_{9}$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}}$	($4g$)	Ho
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}}$	($4g$)	Ho
$\mathbf{B}_{11}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)b\,\hat{\mathbf{y}}$	($4g$)	Ho
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)b\,\hat{\mathbf{y}}$	($4g$)	Ho
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}}$	($4g$)	O II
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}}$	($4g$)	O II
$\mathbf{B}_{15}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)b\,\hat{\mathbf{y}}$	($4g$)	O II
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)b\,\hat{\mathbf{y}}$	($4g$)	O II
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4h$)	Mn II
$\mathbf{B}_{18}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4h$)	Mn II
$\mathbf{B}_{19}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4h$)	Mn II
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4h$)	Mn II
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4h$)	O III
$\mathbf{B}_{22}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4h$)	O III
$\mathbf{B}_{23}$	=	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4h$)	O III
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4h$)	O III
$\mathbf{B}_{25}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	($8i$)	O IV
$\mathbf{B}_{26}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	($8i$)	O IV
$\mathbf{B}_{27}$	=	$\left(\frac{1}{2} - x_7\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right)b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	($8i$)	O IV
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} + x_7\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_7\right)\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_7\right)b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	($8i$)	O IV
$\mathbf{B}_{29}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	($8i$)	O IV
$\mathbf{B}_{30}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	($8i$)	O IV
$\mathbf{B}_{31}$	=	$\left(\frac{1}{2} + x_7\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_7\right)\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_7\right)b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	($8i$)	O IV
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2} - x_7\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right)b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	($8i$)	O IV

References:

- S. Quezel-Ambrunaz, F. Bertaut, and G. Buisson, *Structure des composés d'oxydes de terres rares et de manganèse de formule TMn_2O_5*, C. R. Acad. Sci. **258**, 3025–3027 (1964).
http://gallica.bnf.fr/ark:/12148/bpt6k4011c?rk=85837;2.

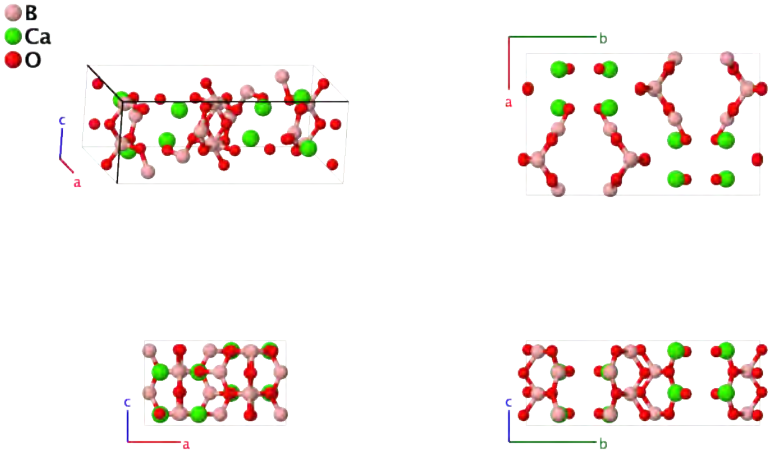
Found in:

- P. Euzen, P. Leone, C. Gueho, and P. Palvadeau, *Structure of* $NdMn_2O_5$, Acta Crystallogr. C **49**, 1875–1877 (1993), doi:10.1107/S0108270193003221.

Geometry files:
- CIF: pp. 1613
- POSCAR: pp. 1613

Calciborite (CaB_2O_4 II) Structure: A2BC4_oP56_56_2e_e_4e

http://aflow.org/prototype-encyclopedia/A2BC4_oP56_56_2e_e_4e

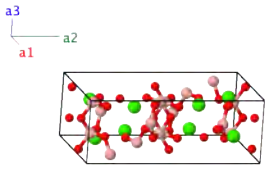


Prototype	:	B_2CaO_4
AFLOW prototype label	:	A2BC4_oP56_56_2e_e_4e
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP56
Space group number	:	56
Space group symbol	:	*Pccn*
AFLOW prototype command	:	`aflow --proto=A2BC4_oP56_56_2e_e_4e --params=`$a, b/a, c/a, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7$

- CaB_2O_4 exists in at least four phases (Marezio, 1969):
- I - The ground state, stable below 1.2 GPa, *Strukturbericht* $E3_2$.
- II – Orthorhombic high pressure structure, stable between 1.2 and 1.5 GPa, presumably calciborite (this structure).
- III – Orthorhombic high pressure structure, stable between 1.5 and 2.5 GPa.
- IV – Cubic high pressure structure, stable between 2.5 and 4.0 GPa.
- This structure has the same orthorhombic lattice constants as the CaB_2O_4 II structure identified in (Marezio, 1969ab), however they state that the structure is stable between 1.2 and 1.5 GPa, while this mineral was obtained "From drillcore into a contact metasomatized limestone near a quartz diorite intrusion associated with a copper deposit in skarn." (Downs, 2003).

Simple Orthorhombic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= a\,\hat{\mathbf{x}}\\ \mathbf{a}_2 &= b\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	$x_1 a\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(8*e*)	B I
$\mathbf{B}_2$	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_1\right)\mathbf{a}_2 + z_1\,\mathbf{a}_3$	$\left(\frac{1}{2} - x_1\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_1\right)b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(8*e*)	B I
$\mathbf{B}_3$	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2} + y_1\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	$-x_1 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right)c\,\hat{\mathbf{z}}$	(8*e*)	B I
$\mathbf{B}_4$	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_1\right)a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right)c\,\hat{\mathbf{z}}$	(8*e*)	B I
$\mathbf{B}_5$	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	$-x_1 a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	(8*e*)	B I
$\mathbf{B}_6$	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_1\right)\mathbf{a}_2 - z_1\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_1\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_1\right)b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	(8*e*)	B I
$\mathbf{B}_7$	$x_1\,\mathbf{a}_1 + \left(\frac{1}{2} - y_1\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	$x_1 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\,\hat{\mathbf{z}}$	(8*e*)	B I
$\mathbf{B}_8$	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + y_1\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_1\right)a\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\,\hat{\mathbf{z}}$	(8*e*)	B I
$\mathbf{B}_9$	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	$x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(8*e*)	B II
$\mathbf{B}_{10}$	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 + z_2\,\mathbf{a}_3$	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right)b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(8*e*)	B II
$\mathbf{B}_{11}$	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	$-x_2 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	(8*e*)	B II
$\mathbf{B}_{12}$	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	(8*e*)	B II
$\mathbf{B}_{13}$	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	$-x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(8*e*)	B II
$\mathbf{B}_{14}$	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 - z_2\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right)b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(8*e*)	B II
$\mathbf{B}_{15}$	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	$x_2 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	(8*e*)	B II
$\mathbf{B}_{16}$	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	(8*e*)	B II
$\mathbf{B}_{17}$	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(8*e*)	Ca
$\mathbf{B}_{18}$	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(8*e*)	Ca
$\mathbf{B}_{19}$	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	$-x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(8*e*)	Ca
$\mathbf{B}_{20}$	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(8*e*)	Ca
$\mathbf{B}_{21}$	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	$-x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(8*e*)	Ca
$\mathbf{B}_{22}$	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 - z_3\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(8*e*)	Ca
$\mathbf{B}_{23}$	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	$x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(8*e*)	Ca
$\mathbf{B}_{24}$	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(8*e*)	Ca
$\mathbf{B}_{25}$	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{26}$	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{27}$	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	$-x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{28}$	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(8*e*)	O I

$\mathbf{B}_{29}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2}+x_4\right)\mathbf{a}_1 + \left(\frac{1}{2}+y_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_4\right)b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{31}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2}-y_4\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_4\right)c\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2}-x_4\right)\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_4\right)c\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{33}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(8*e*)	O II
$\mathbf{B}_{34}$	=	$\left(\frac{1}{2}-x_5\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_5\right)\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_5\right)b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(8*e*)	O II
$\mathbf{B}_{35}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2}+y_5\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_5\right)\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_5\right)c\,\hat{\mathbf{z}}$	(8*e*)	O II
$\mathbf{B}_{36}$	=	$\left(\frac{1}{2}+x_5\right)\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\frac{1}{2}-z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_5\right)c\,\hat{\mathbf{z}}$	(8*e*)	O II
$\mathbf{B}_{37}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(8*e*)	O II
$\mathbf{B}_{38}$	=	$\left(\frac{1}{2}+x_5\right)\mathbf{a}_1 + \left(\frac{1}{2}+y_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_5\right)b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(8*e*)	O II
$\mathbf{B}_{39}$	=	$x_5\,\mathbf{a}_1 + \left(\frac{1}{2}-y_5\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	(8*e*)	O II
$\mathbf{B}_{40}$	=	$\left(\frac{1}{2}-x_5\right)\mathbf{a}_1 + y_5\,\mathbf{a}_2 + \left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	(8*e*)	O II
$\mathbf{B}_{41}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(8*e*)	O III
$\mathbf{B}_{42}$	=	$\left(\frac{1}{2}-x_6\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_6\right)\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_6\right)b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(8*e*)	O III
$\mathbf{B}_{43}$	=	$-x_6\,\mathbf{a}_1 + \left(\frac{1}{2}+y_6\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_6\right)\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_6\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_6\right)c\,\hat{\mathbf{z}}$	(8*e*)	O III
$\mathbf{B}_{44}$	=	$\left(\frac{1}{2}+x_6\right)\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\frac{1}{2}-z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_6\right)c\,\hat{\mathbf{z}}$	(8*e*)	O III
$\mathbf{B}_{45}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(8*e*)	O III
$\mathbf{B}_{46}$	=	$\left(\frac{1}{2}+x_6\right)\mathbf{a}_1 + \left(\frac{1}{2}+y_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_6\right)b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(8*e*)	O III
$\mathbf{B}_{47}$	=	$x_6\,\mathbf{a}_1 + \left(\frac{1}{2}-y_6\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_6\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	(8*e*)	O III
$\mathbf{B}_{48}$	=	$\left(\frac{1}{2}-x_6\right)\mathbf{a}_1 + y_6\,\mathbf{a}_2 + \left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	(8*e*)	O III
$\mathbf{B}_{49}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(8*e*)	O IV
$\mathbf{B}_{50}$	=	$\left(\frac{1}{2}-x_7\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_7\right)\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_7\right)b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(8*e*)	O IV
$\mathbf{B}_{51}$	=	$-x_7\,\mathbf{a}_1 + \left(\frac{1}{2}+y_7\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_7\right)\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_7\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_7\right)c\,\hat{\mathbf{z}}$	(8*e*)	O IV
$\mathbf{B}_{52}$	=	$\left(\frac{1}{2}+x_7\right)\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \left(\frac{1}{2}-z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_7\right)c\,\hat{\mathbf{z}}$	(8*e*)	O IV
$\mathbf{B}_{53}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	(8*e*)	O IV
$\mathbf{B}_{54}$	=	$\left(\frac{1}{2}+x_7\right)\mathbf{a}_1 + \left(\frac{1}{2}+y_7\right)\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_7\right)b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	(8*e*)	O IV
$\mathbf{B}_{55}$	=	$x_7\,\mathbf{a}_1 + \left(\frac{1}{2}-y_7\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_7\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	(8*e*)	O IV
$\mathbf{B}_{56}$	=	$\left(\frac{1}{2}-x_7\right)\mathbf{a}_1 + y_7\,\mathbf{a}_2 + \left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_7\right)a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	(8*e*)	O IV

References:

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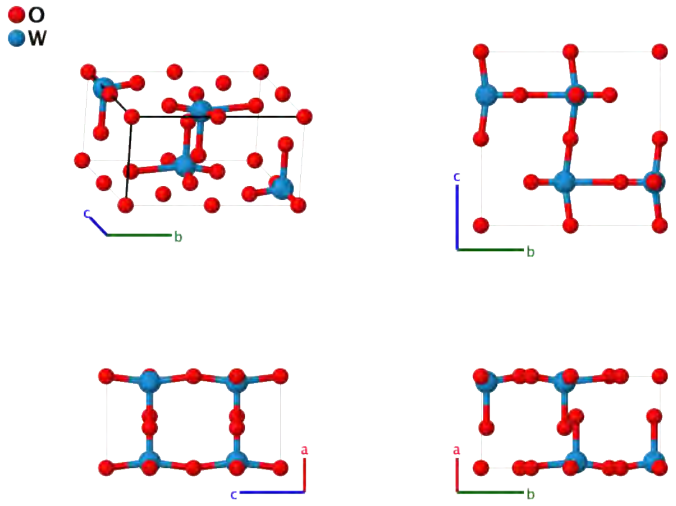
Found in:

- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:
- CIF: pp. 1613
- POSCAR: pp. 1614

$D0_{10}$ (WO_3) (*obsolete*) Structure: A3B_oP16_57_a2d_d

http://aflow.org/prototype-encyclopedia/A3B_oP16_57_a2d_d



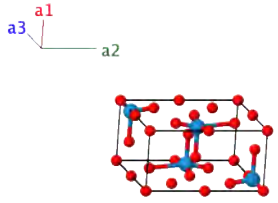
Prototype	:	O_3W
AFLOW prototype label	:	A3B_oP16_57_a2d_d
***Strukturbericht* designation**	:	$D0_{10}$
Pearson symbol	:	oP16
Space group number	:	57
Space group symbol	:	*Pbcm*
AFLOW prototype command	:	`aflow --proto=A3B_oP16_57_a2d_d --params=`$a, b/a, c/a, x_2, y_2, x_3, y_3, x_4, y_4$

- All stable phases of WO_3 are distortions of the cubic α-ReO_3 ($D0_9$) phase. Based on (Woodward, 1997 and Vogt, 1999), the known stable phases and their approximate temperature ranges are:
  - α-WO_3 (1010-1170 K) (Vogt, 1999)
  - β-WO_3 (600-1170 K) (Vogt, 1999)
  - γ-WO_3 (290-600 K) (Vogt, 1999)
  - δ-WO_3 (230-290 K) (Diehl, 1978)
  - ϵ-WO_3 (below 23 K) (Woodward, 1997)

  Woodward's caption says "The transition temperatures display large hysteresis effects and universal agreement is not found in the literature."
- In addition, several other structures have been proposed and/or found:
  - The original $D0_{10}$ structure (Bräkken, 1931), (Hermann, 1937), this structure, superseded by δ-WO_3
  - Original β-WO_3 (Salje, 1977)
  - Hexagonal WO_3 (Gerand, 1979) (metastable)

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}b\,\hat{\mathbf{y}}$	(4*a*)	O I
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$x_2a\,\hat{\mathbf{x}} + y_2b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*d*)	O II
$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}} - y_2b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(4*d*)	O II
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right)b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*d*)	O II
$\mathbf{B}_8$	=	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$x_2a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right)b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(4*d*)	O II
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + y_3b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*d*)	O III
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - y_3b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(4*d*)	O III
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*d*)	O III
$\mathbf{B}_{12}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(4*d*)	O III
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + y_4b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*d*)	W
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} - y_4b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(4*d*)	W
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*d*)	W
$\mathbf{B}_{16}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(4*d*)	W

References:

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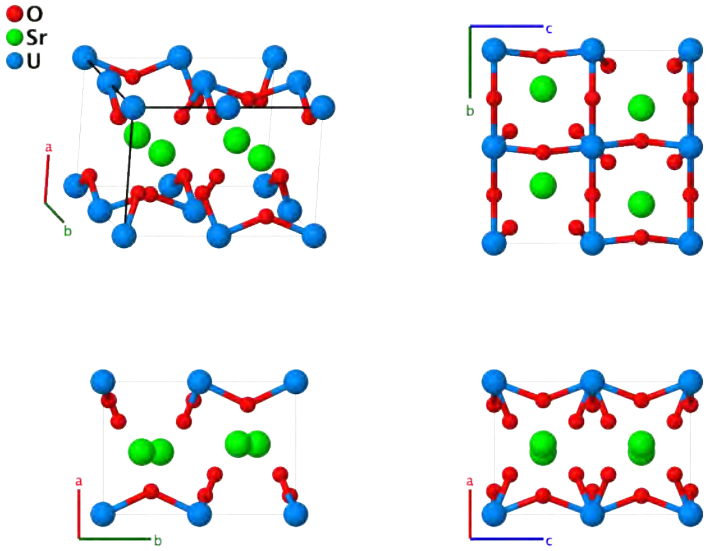
Geometry files:
- CIF: pp. 1614
- POSCAR: pp. 1615

$SrUO_4$ Structure: A4BC_oP24_57_cde_d_a

http://aflow.org/prototype-encyclopedia/A4BC_oP24_57_cde_d_a



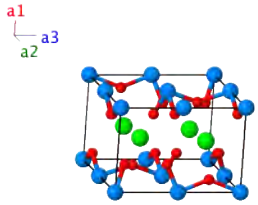
Prototype	:	O_4SrU
AFLOW prototype label	:	A4BC_oP24_57_cde_d_a
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP24
Space group number	:	57
Space group symbol	:	*Pbcm*
AFLOW prototype command	:	`aflow --proto=A4BC_oP24_57_cde_d_a --params=`$a, b/a, c/a, x_2, x_3, y_3, x_4, y_4, x_5, y_5, z_5$

Other compounds with this structure

- $BaUO_4$

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(4*a*)	U
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	(4*a*)	U

$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*a*)	U
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}b\,\hat{\mathbf{y}}$	(4*a*)	U
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2$	=	$x_2a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}}$	(4*c*)	O I
$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2$	=	$-x_2a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}}$	(4*c*)	O I
$\mathbf{B}_8$	=	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$x_2a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + y_3b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*d*)	O II
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - y_3b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(4*d*)	O II
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*d*)	O II
$\mathbf{B}_{12}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(4*d*)	O II
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + y_4b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*d*)	Sr
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} - y_4b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(4*d*)	Sr
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*d*)	Sr
$\mathbf{B}_{16}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(4*d*)	Sr
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}} + y_5b\,\hat{\mathbf{y}} + z_5c\,\hat{\mathbf{z}}$	(8*e*)	O III
$\mathbf{B}_{18}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}} - y_5b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	(8*e*)	O III
$\mathbf{B}_{19}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	(8*e*)	O III
$\mathbf{B}_{20}$	=	$x_5\,\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right)b\,\hat{\mathbf{y}} - z_5c\,\hat{\mathbf{z}}$	(8*e*)	O III
$\mathbf{B}_{21}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}} - y_5b\,\hat{\mathbf{y}} - z_5c\,\hat{\mathbf{z}}$	(8*e*)	O III
$\mathbf{B}_{22}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}} + y_5b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	(8*e*)	O III
$\mathbf{B}_{23}$	=	$x_5\,\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	(8*e*)	O III
$\mathbf{B}_{24}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right)b\,\hat{\mathbf{y}} + z_5c\,\hat{\mathbf{z}}$	(8*e*)	O III

References:

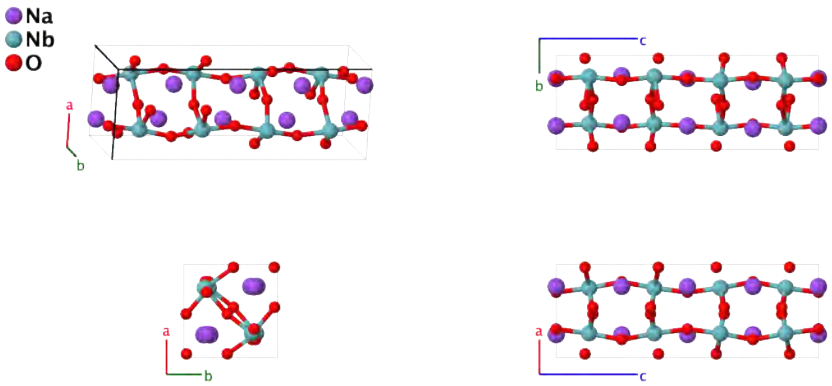
- B. O. Loopstra and H. M. Rietveld, *The structure of some alkaline-earth metal uranates*, Acta Crystallogr. Sect. B Struct. Sci. **25**, 787–791 (1969), doi:10.1107/S0567740869002974.

Geometry files:

- CIF: pp. 1615
- POSCAR: pp. 1615

Lueshite ($NaNbO_3$) Structure: ABC3_oP40_57_cd_e_cd2e

http://aflow.org/prototype-encyclopedia/ABC3_oP40_57_cd_e_cd2e



Prototype	:	$NaNbO_3$
AFLOW prototype label	:	ABC3_oP40_57_cd_e_cd2e
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP40
Space group number	:	57
Space group symbol	:	*Pbcm*
AFLOW prototype command	:	`aflow --proto=ABC3_oP40_57_cd_e_cd2e --params=`$a, b/a, c/a, x_1, x_2, x_3, y_3, x_4, y_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7$

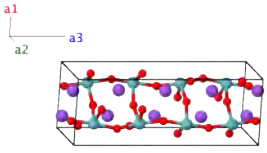
Other compounds with this structure

- $NaNb_{1-x}Ti_xO_3$

- If the AFLOW parameters are set to `--params=`$a, 1, \sqrt{8}, 1/4, 3/4, 1/4, 1/4, 1/4, 3/4, 1/4, 3/4, 5/8, 1/2, 1/2, 5/8, 0, 0, 5/8$ then the structure is equivalent to Cubic Perovskite $E2_1$.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2$	=	$x_1 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}}$	($4c$)	Na I
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \frac{1}{2} c\,\hat{\mathbf{z}}$	($4c$)	Na I
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2$	=	$-x_1 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}}$	($4c$)	Na I
$\mathbf{B}_4$	=	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \frac{1}{2} c\,\hat{\mathbf{z}}$	($4c$)	Na I
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2$	=	$x_2 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}}$	($4c$)	O I

$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \frac{1}{2} c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2$	=	$-x_2 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}}$	(4*c*)	O I
$\mathbf{B}_8$	=	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \frac{1}{2} c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \frac{1}{4} c\,\hat{\mathbf{z}}$	(4*d*)	Na II
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \frac{3}{4} c\,\hat{\mathbf{z}}$	(4*d*)	Na II
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right) b\,\hat{\mathbf{y}} + \frac{1}{4} c\,\hat{\mathbf{z}}$	(4*d*)	Na II
$\mathbf{B}_{12}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right) b\,\hat{\mathbf{y}} + \frac{3}{4} c\,\hat{\mathbf{z}}$	(4*d*)	Na II
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \frac{1}{4} c\,\hat{\mathbf{z}}$	(4*d*)	O II
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \frac{3}{4} c\,\hat{\mathbf{z}}$	(4*d*)	O II
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right) b\,\hat{\mathbf{y}} + \frac{1}{4} c\,\hat{\mathbf{z}}$	(4*d*)	O II
$\mathbf{B}_{16}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right) b\,\hat{\mathbf{y}} + \frac{3}{4} c\,\hat{\mathbf{z}}$	(4*d*)	O II
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(8*e*)	Nb
$\mathbf{B}_{18}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	(8*e*)	Nb
$\mathbf{B}_{19}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right) c\,\hat{\mathbf{z}}$	(8*e*)	Nb
$\mathbf{B}_{20}$	=	$x_5\,\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right) b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(8*e*)	Nb
$\mathbf{B}_{21}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(8*e*)	Nb
$\mathbf{B}_{22}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right) c\,\hat{\mathbf{z}}$	(8*e*)	Nb
$\mathbf{B}_{23}$	=	$x_5\,\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	(8*e*)	Nb
$\mathbf{B}_{24}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right) b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(8*e*)	Nb
$\mathbf{B}_{25}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(8*e*)	O III
$\mathbf{B}_{26}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	(8*e*)	O III
$\mathbf{B}_{27}$	=	$-x_6\,\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right) c\,\hat{\mathbf{z}}$	(8*e*)	O III
$\mathbf{B}_{28}$	=	$x_6\,\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right) b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(8*e*)	O III
$\mathbf{B}_{29}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(8*e*)	O III
$\mathbf{B}_{30}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right) c\,\hat{\mathbf{z}}$	(8*e*)	O III
$\mathbf{B}_{31}$	=	$x_6\,\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	(8*e*)	O III
$\mathbf{B}_{32}$	=	$-x_6\,\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right) b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(8*e*)	O III
$\mathbf{B}_{33}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(8*e*)	O IV
$\mathbf{B}_{34}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right) c\,\hat{\mathbf{z}}$	(8*e*)	O IV
$\mathbf{B}_{35}$	=	$-x_7\,\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right) c\,\hat{\mathbf{z}}$	(8*e*)	O IV
$\mathbf{B}_{36}$	=	$x_7\,\mathbf{a}_1 + \left(\frac{1}{2} - y_7\right)\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_7\right) b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	(8*e*)	O IV
$\mathbf{B}_{37}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	(8*e*)	O IV
$\mathbf{B}_{38}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right) c\,\hat{\mathbf{z}}$	(8*e*)	O IV
$\mathbf{B}_{39}$	=	$x_7\,\mathbf{a}_1 + \left(\frac{1}{2} - y_7\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_7\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right) c\,\hat{\mathbf{z}}$	(8*e*)	O IV
$\mathbf{B}_{40}$	=	$-x_7\,\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right) b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(8*e*)	O IV

References:

- A. C. Sakowski-Cowley, K. Lukaszewicz, and H. D. Megaw, *The structure of sodium niobate at room temperature, and the problem of reliability in pseudosymmetric structures*, Acta Crystallogr. Sect. B Struct. Sci. **25**, 851–865 (1969), doi:10.1107/S0567740869003141.

Found in:

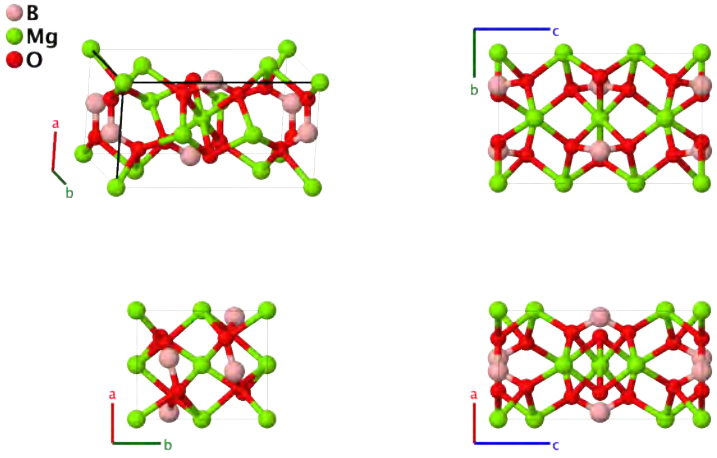
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1615
- POSCAR: pp. 1616

Kotoite ($Mg_3(BO_3)_2$) Structure: A2B3C6_oP22_58_g_af_gh

http://aflow.org/prototype-encyclopedia/A2B3C6_oP22_58_g_af_gh



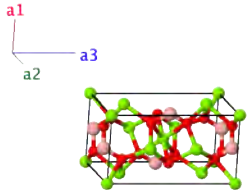
Prototype	:	$B_2Mg_3O_6$
AFLOW prototype label	:	A2B3C6_oP22_58_g_af_gh
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP22
Space group number	:	58
Space group symbol	:	*Pnnm*
AFLOW prototype command	:	`aflow --proto=A2B3C6_oP22_58_g_af_gh --params=`$a, b/a, c/a, z_2, x_3, y_3, x_4, y_4, x_5, y_5, z_5$

Other compounds with this structure

- $Co_3(BO_3)_2$, $Ni_3(BO_3)_2$, $Mn_3(BO_3)_2$ (jimboite), $CoNi_2(BO_3)_2$, $KNa_2(BO_3)_2$, and $Eu_3(BO_3)_2$

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	Mg I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(2a)$	Mg I
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(4f)$	Mg II

$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \left(\frac{1}{2} - z_2\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	$(4f)$	Mg II
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\frac{1}{2}b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	$(4f)$	Mg II
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_1 + \left(\frac{1}{2} + z_2\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	$(4f)$	Mg II
$\mathbf{B}_7$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}}$	$(4g)$	B
$\mathbf{B}_8$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}}$	$(4g)$	B
$\mathbf{B}_9$	=	$\left(\frac{1}{2} - x_3\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(4g)$	B
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} + x_3\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(4g)$	B
$\mathbf{B}_{11}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}}$	$(4g)$	O I
$\mathbf{B}_{12}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}}$	$(4g)$	O I
$\mathbf{B}_{13}$	=	$\left(\frac{1}{2} - x_4\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(4g)$	O I
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} + x_4\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(4g)$	O I
$\mathbf{B}_{15}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(8h)$	O II
$\mathbf{B}_{16}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(8h)$	O II
$\mathbf{B}_{17}$	=	$\left(\frac{1}{2} - x_5\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	$(8h)$	O II
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} + x_5\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	$(8h)$	O II
$\mathbf{B}_{19}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	$(8h)$	O II
$\mathbf{B}_{20}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	$(8h)$	O II
$\mathbf{B}_{21}$	=	$\left(\frac{1}{2} + x_5\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	$(8h)$	O II
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} - x_5\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	$(8h)$	O II

References:

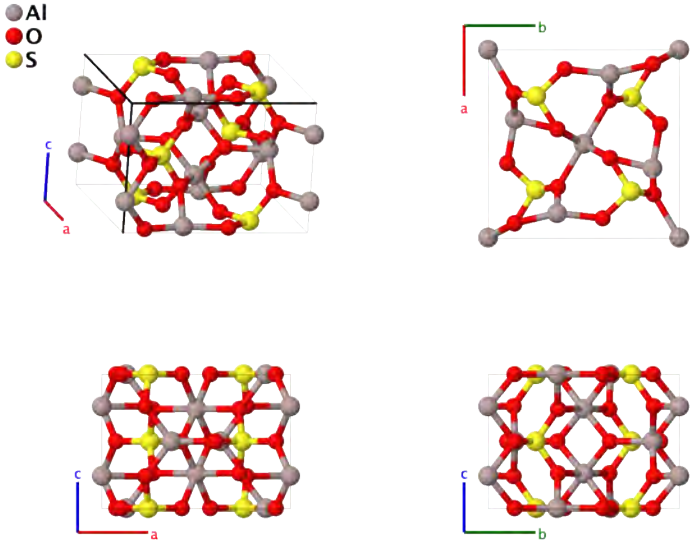
- S. V. Berger, *The Crystal Structure of the Isomorphous Orthoborates of Cobalt and Magnesium*, Acta Chem. Scand. **3**, 660–675 (1949), doi:10.3891/acta.chem.scand.03-0660.

Geometry files:

- CIF: pp. 1616
- POSCAR: pp. 1616

Andalusite (Al_2SiO_5, $S0_2$) Structure: A2B5C_oP32_58_eg_3gh_g

http://aflow.org/prototype-encyclopedia/A2B5C_oP32_58_eg_3gh_g

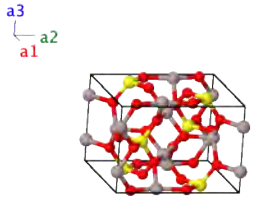


Prototype	:	Al_2O_5Si
AFLOW prototype label	:	A2B5C_oP32_58_eg_3gh_g
***Strukturbericht* designation**	:	$S0_2$
Pearson symbol	:	oP32
Space group number	:	58
Space group symbol	:	*Pnnm*
AFLOW prototype command	:	`aflow --proto=A2B5C_oP32_58_eg_3gh_g --params=`$a, b/a, c/a, z_1, x_2, y_2, x_3, y_3, x_4, y_4, x_5, y_5, x_6, y_6, x_7, y_7, z_7$

- Three crystal polymorphs of Al_2SiO_5 have been characterized: kyanite ($S0_1$), space group $P\bar{1}$ #2, andalusite ($S0_2$), space group *Pnnm* #58, and sillimanite ($S0_3$), space group *Pnma* #62. All are characterized chains of edge-sharing SiO_6 tetrahedra and Al octahedra.
- We use the ambient pressure data of (Yang, 1997).
- (Hermann, 1937) listed this as $S0_2$, but also gave it the $H5_3$ designation in the index.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$z_1\,\mathbf{a}_3$	=	$z_1 c\,\hat{\mathbf{z}}$	($4e$)	Al I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right)c\,\hat{\mathbf{z}}$	($4e$)	Al I
$\mathbf{B}_3$	=	$-z_1\,\mathbf{a}_3$	=	$-z_1 c\,\hat{\mathbf{z}}$	($4e$)	Al I
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\,\hat{\mathbf{z}}$	($4e$)	Al I
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2$	=	$x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}}$	($4g$)	Al II
$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2$	=	$-x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}}$	($4g$)	Al II
$\mathbf{B}_7$	=	$\left(\frac{1}{2} - x_2\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4g$)	Al II
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4g$)	Al II
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}}$	($4g$)	O I
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}}$	($4g$)	O I
$\mathbf{B}_{11}$	=	$\left(\frac{1}{2} - x_3\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4g$)	O I
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4g$)	O I
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}}$	($4g$)	O II
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}}$	($4g$)	O II
$\mathbf{B}_{15}$	=	$\left(\frac{1}{2} - x_4\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4g$)	O II
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_4\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4g$)	O II
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}}$	($4g$)	O III
$\mathbf{B}_{18}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}}$	($4g$)	O III
$\mathbf{B}_{19}$	=	$\left(\frac{1}{2} - x_5\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4g$)	O III
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} + x_5\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4g$)	O III
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2$	=	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}}$	($4g$)	S
$\mathbf{B}_{22}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2$	=	$-x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}}$	($4g$)	S
$\mathbf{B}_{23}$	=	$\left(\frac{1}{2} - x_6\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4g$)	S
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} + x_6\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4g$)	S
$\mathbf{B}_{25}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	($8h$)	O IV
$\mathbf{B}_{26}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	($8h$)	O IV
$\mathbf{B}_{27}$	=	$\left(\frac{1}{2} - x_7\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right)c\,\hat{\mathbf{z}}$	($8h$)	O IV
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} + x_7\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - y_7\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_7\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right)c\,\hat{\mathbf{z}}$	($8h$)	O IV
$\mathbf{B}_{29}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	($8h$)	O IV
$\mathbf{B}_{30}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	($8h$)	O IV
$\mathbf{B}_{31}$	=	$\left(\frac{1}{2} + x_7\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - y_7\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_7\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right)c\,\hat{\mathbf{z}}$	($8h$)	O IV

$\mathbf{B}_{32} = \left(\frac{1}{2} - x_7\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3 = \left(\frac{1}{2} - x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right)c\,\hat{\mathbf{z}}$ (8*h*) O IV

References:

- J. K. Winter and S. Ghose, *Thermal expansion and high-temperaturce crystal chemistry of the Al_2SiO_4 polymorphs*, Am. Mineral. **64**, 573–586 (1979).
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Found in:

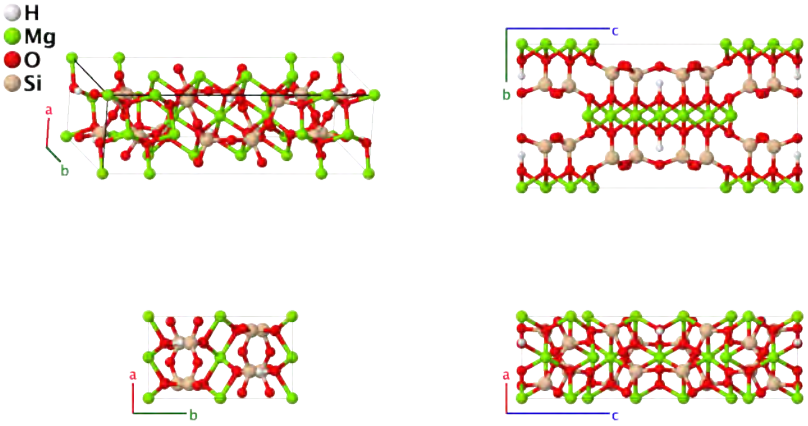
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1616
- POSCAR: pp. 1617

Protoanthophyllite ($H_2Mg_7Si_8O_{24}$) Structure: A2B7C24D8_oP82_58_g_ae2f_2g5h_2h

http://aflow.org/prototype-encyclopedia/A2B7C24D8_oP82_58_g_ae2f_2g5h_2h

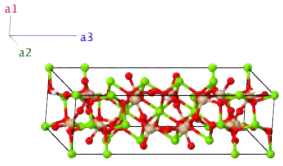


Prototype	:	$H_2Mg_7O_{24}Si_8$
AFLOW prototype label	:	A2B7C24D8_oP82_58_g_ae2f_2g5h_2h
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP82
Space group number	:	58
Space group symbol	:	*Pnnm*
AFLOW prototype command	:	`aflow --proto=A2B7C24D8_oP82_58_g_ae2f_2g5h_2h --params=`$a, b/a, c/a, z_2, z_3, z_4, x_5, y_5, x_6, y_6, x_7, y_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}$

- This structure is approximately one-half of the unit cell of anthophyllite ($S4_4$).
- Like anthophyllite, iron is sometimes mixed with magnesium. For this sample, (Konishi, 2003) found that the Mg-I ($2a$) site contains 2.5% iron, Mg-III ($4f$) contains 2.1%, and Mg-IV ($4f$) contains 27.7% iron.
- (Konishi, 2003) give the crystal structure in the *Pnmn* setting of space group #58. We used FINDSYM to transform this to the standard *Pnnm* orientation.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$ =	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($2a$)	Mg I

$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(2*a*)	Mg I
$\mathbf{B}_3$	=	$z_2\,\mathbf{a}_3$	=	$z_2 c\,\hat{\mathbf{z}}$	(4*e*)	Mg II
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	(4*e*)	Mg II
$\mathbf{B}_5$	=	$-z_2\,\mathbf{a}_3$	=	$-z_2 c\,\hat{\mathbf{z}}$	(4*e*)	Mg II
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	(4*e*)	Mg II
$\mathbf{B}_7$	=	$\frac{1}{2}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{2}b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*f*)	Mg III
$\mathbf{B}_8$	=	$\frac{1}{2}\,\mathbf{a}_1 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(4*f*)	Mg III
$\mathbf{B}_9$	=	$\frac{1}{2}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\frac{1}{2}b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(4*f*)	Mg III
$\mathbf{B}_{10}$	=	$\frac{1}{2}\,\mathbf{a}_1 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(4*f*)	Mg III
$\mathbf{B}_{11}$	=	$\frac{1}{2}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\frac{1}{2}b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(4*f*)	Mg IV
$\mathbf{B}_{12}$	=	$\frac{1}{2}\,\mathbf{a}_1 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(4*f*)	Mg IV
$\mathbf{B}_{13}$	=	$\frac{1}{2}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\frac{1}{2}b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(4*f*)	Mg IV
$\mathbf{B}_{14}$	=	$\frac{1}{2}\,\mathbf{a}_1 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(4*f*)	Mg IV
$\mathbf{B}_{15}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}}$	(4*g*)	H
$\mathbf{B}_{16}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}}$	(4*g*)	H
$\mathbf{B}_{17}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*g*)	H
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*g*)	H
$\mathbf{B}_{19}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2$	=	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}}$	(4*g*)	O I
$\mathbf{B}_{20}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2$	=	$-x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}}$	(4*g*)	O I
$\mathbf{B}_{21}$	=	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*g*)	O I
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*g*)	O I
$\mathbf{B}_{23}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2$	=	$x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}}$	(4*g*)	O II
$\mathbf{B}_{24}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2$	=	$-x_7 a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}}$	(4*g*)	O II
$\mathbf{B}_{25}$	=	$\left(\frac{1}{2} - x_7\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*g*)	O II
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} + x_7\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_7\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_7\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*g*)	O II
$\mathbf{B}_{27}$	=	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(8*h*)	O III
$\mathbf{B}_{28}$	=	$-x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(8*h*)	O III
$\mathbf{B}_{29}$	=	$\left(\frac{1}{2} - x_8\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_8\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_8\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right)c\,\hat{\mathbf{z}}$	(8*h*)	O III
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2} + x_8\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_8\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_8\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right)c\,\hat{\mathbf{z}}$	(8*h*)	O III
$\mathbf{B}_{31}$	=	$-x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(8*h*)	O III
$\mathbf{B}_{32}$	=	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(8*h*)	O III
$\mathbf{B}_{33}$	=	$\left(\frac{1}{2} + x_8\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_8\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_8\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right)c\,\hat{\mathbf{z}}$	(8*h*)	O III
$\mathbf{B}_{34}$	=	$\left(\frac{1}{2} - x_8\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_8\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_8\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right)c\,\hat{\mathbf{z}}$	(8*h*)	O III

$\mathbf{B}_{35}$	=	$x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(8*h*)	O IV
$\mathbf{B}_{36}$	=	$-x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(8*h*)	O IV
$\mathbf{B}_{37}$	=	$\left(\frac{1}{2} - x_9\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_9\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_9\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_9\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_9\right)c\,\hat{\mathbf{z}}$	(8*h*)	O IV
$\mathbf{B}_{38}$	=	$\left(\frac{1}{2} + x_9\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_9\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_9\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_9\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_9\right)c\,\hat{\mathbf{z}}$	(8*h*)	O IV
$\mathbf{B}_{39}$	=	$-x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}}$	(8*h*)	O IV
$\mathbf{B}_{40}$	=	$x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}}$	(8*h*)	O IV
$\mathbf{B}_{41}$	=	$\left(\frac{1}{2} + x_9\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_9\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_9\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_9\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right)c\,\hat{\mathbf{z}}$	(8*h*)	O IV
$\mathbf{B}_{42}$	=	$\left(\frac{1}{2} - x_9\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_9\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_9\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_9\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right)c\,\hat{\mathbf{z}}$	(8*h*)	O IV
$\mathbf{B}_{43}$	=	$x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$x_{10} a\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}}$	(8*h*)	O V
$\mathbf{B}_{44}$	=	$-x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$-x_{10} a\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}}$	(8*h*)	O V
$\mathbf{B}_{45}$	=	$\left(\frac{1}{2} - x_{10}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{10}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{10}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{10}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{10}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{10}\right)c\,\hat{\mathbf{z}}$	(8*h*)	O V
$\mathbf{B}_{46}$	=	$\left(\frac{1}{2} + x_{10}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{10}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{10}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{10}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{10}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{10}\right)c\,\hat{\mathbf{z}}$	(8*h*)	O V
$\mathbf{B}_{47}$	=	$-x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3$	=	$-x_{10} a\,\hat{\mathbf{x}} - y_{10} b\,\hat{\mathbf{y}} - z_{10} c\,\hat{\mathbf{z}}$	(8*h*)	O V
$\mathbf{B}_{48}$	=	$x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3$	=	$x_{10} a\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} - z_{10} c\,\hat{\mathbf{z}}$	(8*h*)	O V
$\mathbf{B}_{49}$	=	$\left(\frac{1}{2} + x_{10}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{10}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{10}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{10}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{10}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{10}\right)c\,\hat{\mathbf{z}}$	(8*h*)	O V
$\mathbf{B}_{50}$	=	$\left(\frac{1}{2} - x_{10}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{10}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{10}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{10}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{10}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{10}\right)c\,\hat{\mathbf{z}}$	(8*h*)	O V
$\mathbf{B}_{51}$	=	$x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$x_{11} a\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} + z_{11} c\,\hat{\mathbf{z}}$	(8*h*)	O VI
$\mathbf{B}_{52}$	=	$-x_{11}\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$-x_{11} a\,\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} + z_{11} c\,\hat{\mathbf{z}}$	(8*h*)	O VI
$\mathbf{B}_{53}$	=	$\left(\frac{1}{2} - x_{11}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{11}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{11}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{11}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{11}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{11}\right)c\,\hat{\mathbf{z}}$	(8*h*)	O VI
$\mathbf{B}_{54}$	=	$\left(\frac{1}{2} + x_{11}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{11}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{11}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{11}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{11}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{11}\right)c\,\hat{\mathbf{z}}$	(8*h*)	O VI
$\mathbf{B}_{55}$	=	$-x_{11}\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3$	=	$-x_{11} a\,\hat{\mathbf{x}} - y_{11} b\,\hat{\mathbf{y}} - z_{11} c\,\hat{\mathbf{z}}$	(8*h*)	O VI
$\mathbf{B}_{56}$	=	$x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3$	=	$x_{11} a\,\hat{\mathbf{x}} + y_{11} b\,\hat{\mathbf{y}} - z_{11} c\,\hat{\mathbf{z}}$	(8*h*)	O VI
$\mathbf{B}_{57}$	=	$\left(\frac{1}{2} + x_{11}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{11}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{11}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{11}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{11}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{11}\right)c\,\hat{\mathbf{z}}$	(8*h*)	O VI
$\mathbf{B}_{58}$	=	$\left(\frac{1}{2} - x_{11}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{11}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{11}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{11}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{11}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{11}\right)c\,\hat{\mathbf{z}}$	(8*h*)	O VI
$\mathbf{B}_{59}$	=	$x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3$	=	$x_{12} a\,\hat{\mathbf{x}} + y_{12} b\,\hat{\mathbf{y}} + z_{12} c\,\hat{\mathbf{z}}$	(8*h*)	O VII
$\mathbf{B}_{60}$	=	$-x_{12}\,\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3$	=	$-x_{12} a\,\hat{\mathbf{x}} - y_{12} b\,\hat{\mathbf{y}} + z_{12} c\,\hat{\mathbf{z}}$	(8*h*)	O VII

$\mathbf{B}_{61}$	=	$\left(\frac{1}{2}-x_{12}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{12}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{12}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{12}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{12}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{12}\right)c\,\hat{\mathbf{z}}$	(8*h*)	O VII
$\mathbf{B}_{62}$	=	$\left(\frac{1}{2}+x_{12}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{12}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{12}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{12}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{12}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{12}\right)c\,\hat{\mathbf{z}}$	(8*h*)	O VII
$\mathbf{B}_{63}$	=	$-x_{12}\,\mathbf{a}_1-y_{12}\,\mathbf{a}_2-z_{12}\,\mathbf{a}_3$	=	$-x_{12}a\,\hat{\mathbf{x}}-y_{12}b\,\hat{\mathbf{y}}-z_{12}c\,\hat{\mathbf{z}}$	(8*h*)	O VII
$\mathbf{B}_{64}$	=	$x_{12}\,\mathbf{a}_1+y_{12}\,\mathbf{a}_2-z_{12}\,\mathbf{a}_3$	=	$x_{12}a\,\hat{\mathbf{x}}+y_{12}b\,\hat{\mathbf{y}}-z_{12}c\,\hat{\mathbf{z}}$	(8*h*)	O VII
$\mathbf{B}_{65}$	=	$\left(\frac{1}{2}+x_{12}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{12}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{12}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{12}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{12}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{12}\right)c\,\hat{\mathbf{z}}$	(8*h*)	O VII
$\mathbf{B}_{66}$	=	$\left(\frac{1}{2}-x_{12}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{12}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{12}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{12}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{12}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{12}\right)c\,\hat{\mathbf{z}}$	(8*h*)	O VII
$\mathbf{B}_{67}$	=	$x_{13}\,\mathbf{a}_1+y_{13}\,\mathbf{a}_2+z_{13}\,\mathbf{a}_3$	=	$x_{13}a\,\hat{\mathbf{x}}+y_{13}b\,\hat{\mathbf{y}}+z_{13}c\,\hat{\mathbf{z}}$	(8*h*)	Si I
$\mathbf{B}_{68}$	=	$-x_{13}\,\mathbf{a}_1-y_{13}\,\mathbf{a}_2+z_{13}\,\mathbf{a}_3$	=	$-x_{13}a\,\hat{\mathbf{x}}-y_{13}b\,\hat{\mathbf{y}}+z_{13}c\,\hat{\mathbf{z}}$	(8*h*)	Si I
$\mathbf{B}_{69}$	=	$\left(\frac{1}{2}-x_{13}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{13}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{13}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{13}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{13}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{13}\right)c\,\hat{\mathbf{z}}$	(8*h*)	Si I
$\mathbf{B}_{70}$	=	$\left(\frac{1}{2}+x_{13}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{13}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{13}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{13}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{13}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{13}\right)c\,\hat{\mathbf{z}}$	(8*h*)	Si I
$\mathbf{B}_{71}$	=	$-x_{13}\,\mathbf{a}_1-y_{13}\,\mathbf{a}_2-z_{13}\,\mathbf{a}_3$	=	$-x_{13}a\,\hat{\mathbf{x}}-y_{13}b\,\hat{\mathbf{y}}-z_{13}c\,\hat{\mathbf{z}}$	(8*h*)	Si I
$\mathbf{B}_{72}$	=	$x_{13}\,\mathbf{a}_1+y_{13}\,\mathbf{a}_2-z_{13}\,\mathbf{a}_3$	=	$x_{13}a\,\hat{\mathbf{x}}+y_{13}b\,\hat{\mathbf{y}}-z_{13}c\,\hat{\mathbf{z}}$	(8*h*)	Si I
$\mathbf{B}_{73}$	=	$\left(\frac{1}{2}+x_{13}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{13}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{13}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{13}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{13}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{13}\right)c\,\hat{\mathbf{z}}$	(8*h*)	Si I
$\mathbf{B}_{74}$	=	$\left(\frac{1}{2}-x_{13}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{13}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{13}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{13}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{13}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{13}\right)c\,\hat{\mathbf{z}}$	(8*h*)	Si I
$\mathbf{B}_{75}$	=	$x_{14}\,\mathbf{a}_1+y_{14}\,\mathbf{a}_2+z_{14}\,\mathbf{a}_3$	=	$x_{14}a\,\hat{\mathbf{x}}+y_{14}b\,\hat{\mathbf{y}}+z_{14}c\,\hat{\mathbf{z}}$	(8*h*)	Si II
$\mathbf{B}_{76}$	=	$-x_{14}\,\mathbf{a}_1-y_{14}\,\mathbf{a}_2+z_{14}\,\mathbf{a}_3$	=	$-x_{14}a\,\hat{\mathbf{x}}-y_{14}b\,\hat{\mathbf{y}}+z_{14}c\,\hat{\mathbf{z}}$	(8*h*)	Si II
$\mathbf{B}_{77}$	=	$\left(\frac{1}{2}-x_{14}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{14}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{14}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{14}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{14}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{14}\right)c\,\hat{\mathbf{z}}$	(8*h*)	Si II
$\mathbf{B}_{78}$	=	$\left(\frac{1}{2}+x_{14}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{14}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{14}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{14}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{14}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{14}\right)c\,\hat{\mathbf{z}}$	(8*h*)	Si II
$\mathbf{B}_{79}$	=	$-x_{14}\,\mathbf{a}_1-y_{14}\,\mathbf{a}_2-z_{14}\,\mathbf{a}_3$	=	$-x_{14}a\,\hat{\mathbf{x}}-y_{14}b\,\hat{\mathbf{y}}-z_{14}c\,\hat{\mathbf{z}}$	(8*h*)	Si II
$\mathbf{B}_{80}$	=	$x_{14}\,\mathbf{a}_1+y_{14}\,\mathbf{a}_2-z_{14}\,\mathbf{a}_3$	=	$x_{14}a\,\hat{\mathbf{x}}+y_{14}b\,\hat{\mathbf{y}}-z_{14}c\,\hat{\mathbf{z}}$	(8*h*)	Si II
$\mathbf{B}_{81}$	=	$\left(\frac{1}{2}+x_{14}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{14}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{14}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{14}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{14}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{14}\right)c\,\hat{\mathbf{z}}$	(8*h*)	Si II
$\mathbf{B}_{82}$	=	$\left(\frac{1}{2}-x_{14}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{14}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{14}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{14}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{14}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{14}\right)c\,\hat{\mathbf{z}}$	(8*h*)	Si II

References:

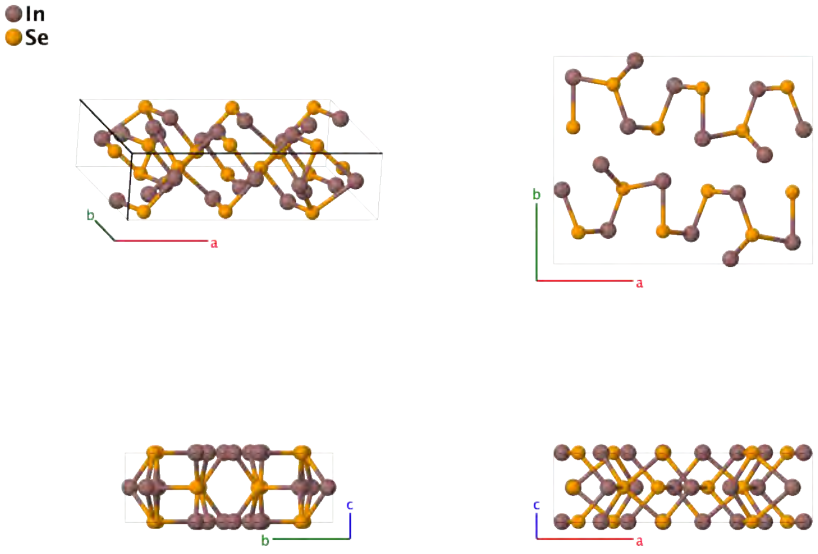
- H. Konishi, T. L. Groy, I. Dódony, R. Miyawaki, S. Matsubara, and P. R. Buseck, *Crystal structure of protoanthophyllite: A new mineral from the Takase ultramafic complex, Japan*, Am. Mineral. **88**, 1718–1723 (2003), doi:10.2138/am-2003-11-1212.

Geometry files:

- CIF: pp. 1617
- POSCAR: pp. 1617

In_4Se_3 Structure: A4B3_oP28_58_4g_3g

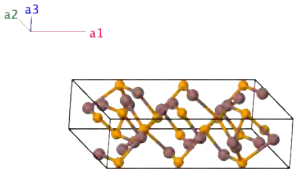
http://aflow.org/prototype-encyclopedia/A4B3_oP28_58_4g_3g



Prototype	:	In_4Se_3
AFLOW prototype label	:	A4B3_oP28_58_4g_3g
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP28
Space group number	:	58
Space group symbol	:	*Pnnm*
AFLOW prototype command	:	`aflow --proto=A4B3_oP28_58_4g_3g --params=`$a, b/a, c/a, x_1, y_1, x_2, y_2, x_3, y_3, x_4, y_4, x_5, y_5, x_6, y_6, x_7, y_7$

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2$	=	$x_1 a\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}}$	$(4g)$	In I
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2$	=	$-x_1 a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}}$	$(4g)$	In I
$\mathbf{B}_3$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_1\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_1\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(4g)$	In I

$\mathbf{B}_4$	=	$\left(\frac{1}{2}+x_1\right)\mathbf{a}_1+\left(\frac{1}{2}-y_1\right)\mathbf{a}_2+\frac{1}{2}\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_1\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_1\right)b\,\hat{\mathbf{y}}+\frac{1}{2}c\,\hat{\mathbf{z}}$	(4*g*)	In I
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1+y_2\,\mathbf{a}_2$	=	$x_2a\,\hat{\mathbf{x}}+y_2b\,\hat{\mathbf{y}}$	(4*g*)	In II
$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1-y_2\,\mathbf{a}_2$	=	$-x_2a\,\hat{\mathbf{x}}-y_2b\,\hat{\mathbf{y}}$	(4*g*)	In II
$\mathbf{B}_7$	=	$\left(\frac{1}{2}-x_2\right)\mathbf{a}_1+\left(\frac{1}{2}+y_2\right)\mathbf{a}_2+\frac{1}{2}\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_2\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_2\right)b\,\hat{\mathbf{y}}+\frac{1}{2}c\,\hat{\mathbf{z}}$	(4*g*)	In II
$\mathbf{B}_8$	=	$\left(\frac{1}{2}+x_2\right)\mathbf{a}_1+\left(\frac{1}{2}-y_2\right)\mathbf{a}_2+\frac{1}{2}\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_2\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_2\right)b\,\hat{\mathbf{y}}+\frac{1}{2}c\,\hat{\mathbf{z}}$	(4*g*)	In II
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1+y_3\,\mathbf{a}_2$	=	$x_3a\,\hat{\mathbf{x}}+y_3b\,\hat{\mathbf{y}}$	(4*g*)	In III
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1-y_3\,\mathbf{a}_2$	=	$-x_3a\,\hat{\mathbf{x}}-y_3b\,\hat{\mathbf{y}}$	(4*g*)	In III
$\mathbf{B}_{11}$	=	$\left(\frac{1}{2}-x_3\right)\mathbf{a}_1+\left(\frac{1}{2}+y_3\right)\mathbf{a}_2+\frac{1}{2}\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_3\right)b\,\hat{\mathbf{y}}+\frac{1}{2}c\,\hat{\mathbf{z}}$	(4*g*)	In III
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2}+x_3\right)\mathbf{a}_1+\left(\frac{1}{2}-y_3\right)\mathbf{a}_2+\frac{1}{2}\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_3\right)b\,\hat{\mathbf{y}}+\frac{1}{2}c\,\hat{\mathbf{z}}$	(4*g*)	In III
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1+y_4\,\mathbf{a}_2$	=	$x_4a\,\hat{\mathbf{x}}+y_4b\,\hat{\mathbf{y}}$	(4*g*)	In IV
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1-y_4\,\mathbf{a}_2$	=	$-x_4a\,\hat{\mathbf{x}}-y_4b\,\hat{\mathbf{y}}$	(4*g*)	In IV
$\mathbf{B}_{15}$	=	$\left(\frac{1}{2}-x_4\right)\mathbf{a}_1+\left(\frac{1}{2}+y_4\right)\mathbf{a}_2+\frac{1}{2}\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_4\right)b\,\hat{\mathbf{y}}+\frac{1}{2}c\,\hat{\mathbf{z}}$	(4*g*)	In IV
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2}+x_4\right)\mathbf{a}_1+\left(\frac{1}{2}-y_4\right)\mathbf{a}_2+\frac{1}{2}\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_4\right)b\,\hat{\mathbf{y}}+\frac{1}{2}c\,\hat{\mathbf{z}}$	(4*g*)	In IV
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1+y_5\,\mathbf{a}_2$	=	$x_5a\,\hat{\mathbf{x}}+y_5b\,\hat{\mathbf{y}}$	(4*g*)	Se I
$\mathbf{B}_{18}$	=	$-x_5\,\mathbf{a}_1-y_5\,\mathbf{a}_2$	=	$-x_5a\,\hat{\mathbf{x}}-y_5b\,\hat{\mathbf{y}}$	(4*g*)	Se I
$\mathbf{B}_{19}$	=	$\left(\frac{1}{2}-x_5\right)\mathbf{a}_1+\left(\frac{1}{2}+y_5\right)\mathbf{a}_2+\frac{1}{2}\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_5\right)b\,\hat{\mathbf{y}}+\frac{1}{2}c\,\hat{\mathbf{z}}$	(4*g*)	Se I
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2}+x_5\right)\mathbf{a}_1+\left(\frac{1}{2}-y_5\right)\mathbf{a}_2+\frac{1}{2}\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_5\right)b\,\hat{\mathbf{y}}+\frac{1}{2}c\,\hat{\mathbf{z}}$	(4*g*)	Se I
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1+y_6\,\mathbf{a}_2$	=	$x_6a\,\hat{\mathbf{x}}+y_6b\,\hat{\mathbf{y}}$	(4*g*)	Se II
$\mathbf{B}_{22}$	=	$-x_6\,\mathbf{a}_1-y_6\,\mathbf{a}_2$	=	$-x_6a\,\hat{\mathbf{x}}-y_6b\,\hat{\mathbf{y}}$	(4*g*)	Se II
$\mathbf{B}_{23}$	=	$\left(\frac{1}{2}-x_6\right)\mathbf{a}_1+\left(\frac{1}{2}+y_6\right)\mathbf{a}_2+\frac{1}{2}\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_6\right)b\,\hat{\mathbf{y}}+\frac{1}{2}c\,\hat{\mathbf{z}}$	(4*g*)	Se II
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2}+x_6\right)\mathbf{a}_1+\left(\frac{1}{2}-y_6\right)\mathbf{a}_2+\frac{1}{2}\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_6\right)b\,\hat{\mathbf{y}}+\frac{1}{2}c\,\hat{\mathbf{z}}$	(4*g*)	Se II
$\mathbf{B}_{25}$	=	$x_7\,\mathbf{a}_1+y_7\,\mathbf{a}_2$	=	$x_7a\,\hat{\mathbf{x}}+y_7b\,\hat{\mathbf{y}}$	(4*g*)	Se III
$\mathbf{B}_{26}$	=	$-x_7\,\mathbf{a}_1-y_7\,\mathbf{a}_2$	=	$-x_7a\,\hat{\mathbf{x}}-y_7b\,\hat{\mathbf{y}}$	(4*g*)	Se III
$\mathbf{B}_{27}$	=	$\left(\frac{1}{2}-x_7\right)\mathbf{a}_1+\left(\frac{1}{2}+y_7\right)\mathbf{a}_2+\frac{1}{2}\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_7\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_7\right)b\,\hat{\mathbf{y}}+\frac{1}{2}c\,\hat{\mathbf{z}}$	(4*g*)	Se III
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2}+x_7\right)\mathbf{a}_1+\left(\frac{1}{2}-y_7\right)\mathbf{a}_2+\frac{1}{2}\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_7\right)b\,\hat{\mathbf{y}}+\frac{1}{2}c\,\hat{\mathbf{z}}$	(4*g*)	Se III

References:

- J. H. C. Hogg, H. H. Sutherland, and D. J. Williams, *The crystal structure of tetraindium triselenide*, Acta Crystallogr. Sect. B Struct. Sci. **29**, 1590–1593 (1973), doi:10.1107/S0567740873005108.

Found in:

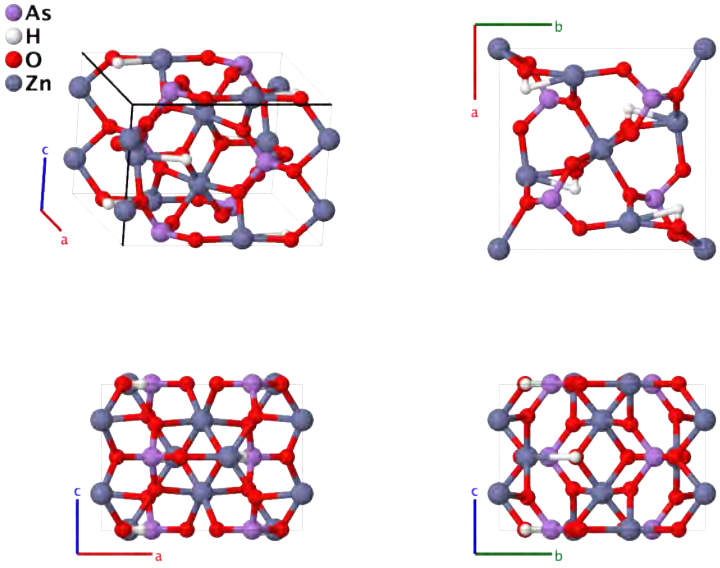
- N. Benramdane and R. H. Misho, *Structural and optical properties of In_4Se_3 thin films obtained by flash evaporation*, Sol. Energy Mater. Sol. Cells **37**, 367–377 (1995), doi:10.1016/0927-0248(95)00031-3.

Geometry files:

- CIF: pp. 1618
- POSCAR: pp. 1618

Adamite [$Zn_2(AsO_4)(OH)$, $H2_7$] Structure: ABC5D2_oP36_58_g_g_3gh_eg

http://aflow.org/prototype-encyclopedia/ABC5D2_oP36_58_g_g_3gh_eg



Prototype	:	$AsHO_5Zn_2$
AFLOW prototype label	:	ABC5D2_oP36_58_g_g_3gh_eg
***Strukturbericht* designation**	:	$H2_7$
Pearson symbol	:	oP36
Space group number	:	58
Space group symbol	:	*Pnnm*
AFLOW prototype command	:	`aflow --proto=ABC5D2_oP36_58_g_g_3gh_eg --params=`$a, b/a, c/a, z_1, x_2, y_2, x_3, y_3, x_4, y_4, x_5, y_5, x_6, y_6, x_7, y_7, x_8, y_8, z_8$

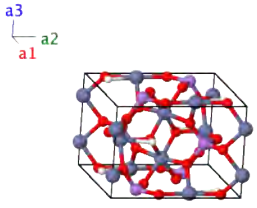
Other compounds with this structure

- $Cu_2(AsO_4)(OH)$ and $Cu_2(PO_4)(OH)$

- This structure was originally determined by (Kokkoros, 1937) and designated $H2_7$ by (Gottfried, 1940). (Hill, 1976) refined the structure, including the positions of the hydrogen atoms in the OH radical.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$z_1\,\mathbf{a}_3$	=	$z_1 c\,\hat{\mathbf{z}}$	($4e$)	Zn I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right)c\,\hat{\mathbf{z}}$	($4e$)	Zn I
$\mathbf{B}_3$	=	$-z_1\,\mathbf{a}_3$	=	$-z_1 c\,\hat{\mathbf{z}}$	($4e$)	Zn I
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\,\hat{\mathbf{z}}$	($4e$)	Zn I
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2$	=	$x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}}$	($4g$)	As
$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2$	=	$-x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}}$	($4g$)	As
$\mathbf{B}_7$	=	$\left(\frac{1}{2} - x_2\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4g$)	As
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4g$)	As
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}}$	($4g$)	H
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}}$	($4g$)	H
$\mathbf{B}_{11}$	=	$\left(\frac{1}{2} - x_3\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4g$)	H
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4g$)	H
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}}$	($4g$)	O I
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}}$	($4g$)	O I
$\mathbf{B}_{15}$	=	$\left(\frac{1}{2} - x_4\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4g$)	O I
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_4\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4g$)	O I
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}}$	($4g$)	O II
$\mathbf{B}_{18}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}}$	($4g$)	O II
$\mathbf{B}_{19}$	=	$\left(\frac{1}{2} - x_5\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4g$)	O II
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} + x_5\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4g$)	O II
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2$	=	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}}$	($4g$)	O III
$\mathbf{B}_{22}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2$	=	$-x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}}$	($4g$)	O III
$\mathbf{B}_{23}$	=	$\left(\frac{1}{2} - x_6\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4g$)	O III
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} + x_6\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4g$)	O III
$\mathbf{B}_{25}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2$	=	$x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}}$	($4g$)	Zn II
$\mathbf{B}_{26}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2$	=	$-x_7 a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}}$	($4g$)	Zn II
$\mathbf{B}_{27}$	=	$\left(\frac{1}{2} - x_7\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4g$)	Zn II
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} + x_7\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - y_7\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_7\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4g$)	Zn II
$\mathbf{B}_{29}$	=	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	($8h$)	O IV
$\mathbf{B}_{30}$	=	$-x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	($8h$)	O IV
$\mathbf{B}_{31}$	=	$\left(\frac{1}{2} - x_8\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + y_8\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_8\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right)c\,\hat{\mathbf{z}}$	($8h$)	O IV
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2} + x_8\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - y_8\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_8\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right)c\,\hat{\mathbf{z}}$	($8h$)	O IV

$\mathbf{B}_{33}$	=	$-x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(8*h*)	O IV
$\mathbf{B}_{34}$	=	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(8*h*)	O IV
$\mathbf{B}_{35}$	=	$\left(\frac{1}{2}+x_8\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_8\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_8\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_8\right)c\,\hat{\mathbf{z}}$	(8*h*)	O IV
$\mathbf{B}_{36}$	=	$\left(\frac{1}{2}-x_8\right)\mathbf{a}_1 + \left(\frac{1}{2}+y_8\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_8\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_8\right)c\,\hat{\mathbf{z}}$	(8*h*)	O IV

References:

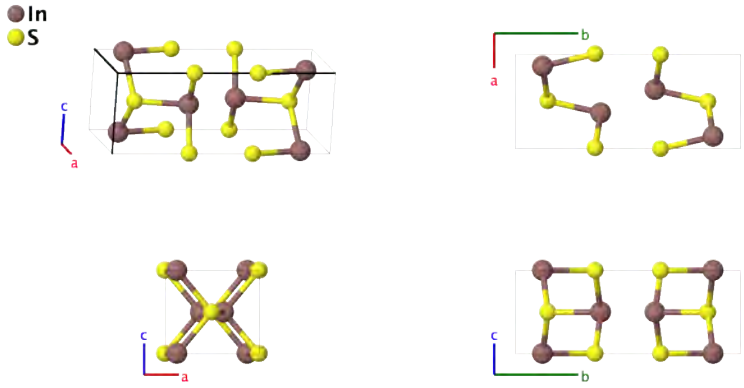
- R. J. Hill, *The crystal structure and infrared properties of adamite*, Am. Mineral. **61**, 979–986 (1976).
- P. Kokkoros, *Über die Struktur von Adamin*, Zeitschrift für Kristallographie - Crystalline Materials **96**, 417–434 (1937), doi:10.1524/zkri.1937.96.1.417.
- C. Gottfried, ed., *Strukturbericht Band V 1937* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1940).

Geometry files:

- CIF: pp. 1618
- POSCAR: pp. 1619

InS Structure: AB_oP8_58_g_g

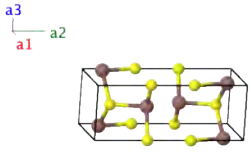
http://aflow.org/prototype-encyclopedia/AB_oP8_58_g_g



Prototype	:	InS
AFLOW prototype label	:	AB_oP8_58_g_g
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP8
Space group number	:	58
Space group symbol	:	*Pnnm*
AFLOW prototype command	:	`aflow --proto=AB_oP8_58_g_g --params=`$a, b/a, c/a, x_1, y_1, x_2, y_2$

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2$	=	$x_1 a\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}}$	$(4g)$	In
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2$	=	$-x_1 a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}}$	$(4g)$	In
$\mathbf{B}_3$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_1\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_1\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(4g)$	In
$\mathbf{B}_4$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_1\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_1\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(4g)$	In
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2$	=	$x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}}$	$(4g)$	S
$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2$	=	$-x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}}$	$(4g)$	S
$\mathbf{B}_7$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(4g)$	S
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right)b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(4g)$	S

References:

- K. Schubert, E. Dörre, and E. Günzel, *Kristallchemische Ergebnisse an Phasen aus B-Elementen*, Naturwissenschaften **41**, 448 (1954), doi:10.1007/BF00628872.

Found in:

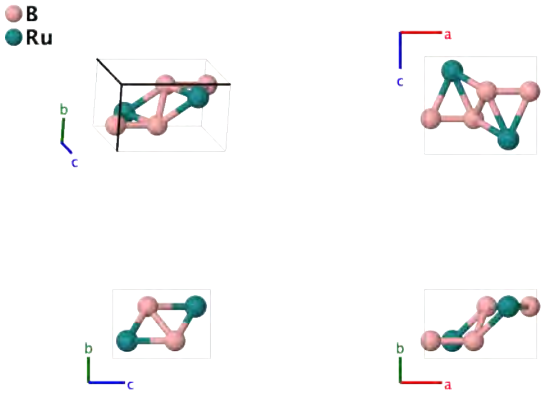
- P. Villars and L. Calvert, *Pearson's Handbook of Crystallographic Data for Intermetallic Phases* (ASM International, Materials Park, OH, 1991), 2nd edn.

Geometry files:

- CIF: pp. 1619
- POSCAR: pp. 1619

RuB_2 Structure: A2B_oP6_59_f_a

http://aflow.org/prototype-encyclopedia/A2B_oP6_59_f_a



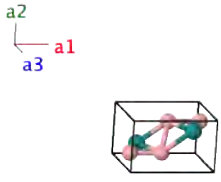
Prototype	:	B_2Ru
AFLOW prototype label	:	A2B_oP6_59_f_a
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP6
Space group number	:	59
Space group symbol	:	*Pmmn*
AFLOW prototype command	:	`aflow --proto=A2B_oP6_59_f_a --params=`$a, b/a, c/a, z_1, x_2, z_2$

Other compounds with this structure

- OsB_2

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	$(2a)$	Ru
$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	$(2a)$	Ru
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(4f)$	B
$\mathbf{B}_4$	=	$\left(\frac{1}{2} - x_2\right)\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(4f)$	B
$\mathbf{B}_5$	=	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	$(4f)$	B
$\mathbf{B}_6$	=	$\left(\frac{1}{2} + x_2\right)\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	$(4f)$	B

References:

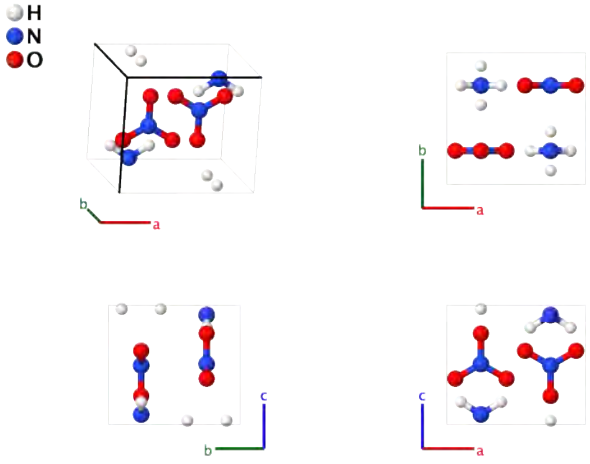
- B. Aronsson, *The Crystal Structure of RuB_2, OsB_2, and $IrB_{1.35}$ and Some General Comments on the Crystal Chemistry of Borides in the Composition Range MeB - MeB_3*, Acta Chem. Scand. **17**, 2036–2050 (1963), doi:10.3891/acta.chem.scand.17-2036.

Geometry files:

- CIF: pp. 1620
- POSCAR: pp. 1620

NH_4NO_3 IV ($G0_{11}$) Structure: A4B2C3_oP18_59_ef_ab_af

http://aflow.org/prototype-encyclopedia/A4B2C3_oP18_59_ef_ab_af



Prototype	:	$H_4N_2O_3$
AFLOW prototype label	:	A4B2C3_oP18_59_ef_ab_af
***Strukturbericht* designation**	:	$G0_{11}$
Pearson symbol	:	oP18
Space group number	:	59
Space group symbol	:	*Pmmn*
AFLOW prototype command	:	`aflow --proto=A4B2C3_oP18_59_ef_ab_af --params=`$a, b/a, c/a, z_1, z_2, z_3, y_4, z_4, x_5, z_5, x_6, z_6$

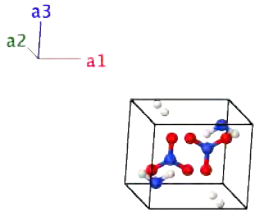
- Ammonium Nitrate exists in a variety of forms, (Hermann, 1937) depending on the temperature:

Phase	Temperature °C	*Strukturbericht*	Page	
I	125 – 170	$G0_8$	AB_cP2_221_a_b.NH4.NO3	
II	84 – 125	$G0_9$	ABC3_tP10_100_b_a_bc	
III	32 – 84	$G0_{10}$	ABC3_oP20_62_c_c_cd.N.NH4.O	
IV	-17 – 32	$G0_{11}$	A4B2C3_oP18_59_ef_ab_af	(this structure)
V	< -17	Gwihabaite	A4B2C3_tP72_77_8d_ab2c2d_6d2	

- In the original reference (West, 1932) did not determine the positions of the hydrogen atoms. Since the hydrogen atoms are in the same space group, we continue to designate this the $G0_{11}$ structure.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	$(2a)$	N I
$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	$(2a)$	N I
$\mathbf{B}_3$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(2a)$	O I
$\mathbf{B}_4$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	$(2a)$	O I
$\mathbf{B}_5$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(2b)$	N II
$\mathbf{B}_6$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(2b)$	N II
$\mathbf{B}_7$	=	$\frac{1}{4}\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(4e)$	H I
$\mathbf{B}_8$	=	$\frac{1}{4}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(4e)$	H I
$\mathbf{B}_9$	=	$\frac{3}{4}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	$(4e)$	H I
$\mathbf{B}_{10}$	=	$\frac{3}{4}\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	$(4e)$	H I
$\mathbf{B}_{11}$	=	$x_5\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(4f)$	H II
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(4f)$	H II
$\mathbf{B}_{13}$	=	$-x_5\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	$(4f)$	H II
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	$(4f)$	H II
$\mathbf{B}_{15}$	=	$x_6\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(4f)$	O II
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(4f)$	O II
$\mathbf{B}_{17}$	=	$-x_6\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	$(4f)$	O II
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	$(4f)$	O II

References:

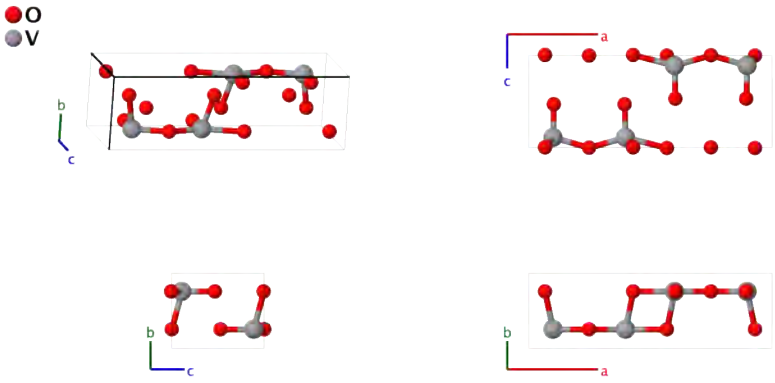
- C. S. Choi, J. E. Mapes, and E. Prince, *The structure of ammonium nitrate (IV)*, Acta Crystallogr. Sect. B Struct. Sci. **28**, 1357–1361 (1972), doi:10.1107/S0567740872004303.
- C. D. West, *The Crystal Structure of Rhombic Ammonium Nitrate*, J. Am. Chem. Soc. **54**, 2256–2260 (1932), doi:10.1021/ja01345a013.
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1620
- POSCAR: pp. 1620

Shcherbinaite (V_2O_5) (*revised*) Structure: A5B2_oP14_59_a2f_f

http://aflow.org/prototype-encyclopedia/A5B2_oP14_59_a2f_f

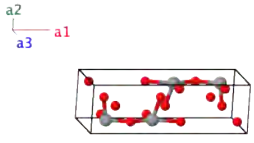


Prototype	:	O_5V_2
AFLOW prototype label	:	A5B2_oP14_59_a2f_f
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP14
Space group number	:	59
Space group symbol	:	*Pmmn*
AFLOW prototype command	:	`aflow --proto=A5B2_oP14_59_a2f_f --params=`$a, b/a, c/a, z_1, x_2, z_2, x_3, z_3, x_4, z_4$

- An earlier version of this structure found by (Ketelaar, 1936) was given the $D8_7$ *Strukturbericht* designation in (Gottfried, 1938). It was later realized that this structure had a rather unexpected arrangement of vanadium atoms, and the structure was revised by (Enjalbert, 1986) and others.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	$(2a)$	O I
$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	$(2a)$	O I
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(4f)$	O II
$\mathbf{B}_4$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(4f)$	O II
$\mathbf{B}_5$	=	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	$(4f)$	O II

$\mathbf{B}_6$	=	$\left(\frac{1}{2}+x_2\right)\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2-z_2\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_2\right)a\hat{\mathbf{x}}+\frac{3}{4}b\hat{\mathbf{y}}-z_2c\hat{\mathbf{z}}$	$(4f)$	O II
$\mathbf{B}_7$	=	$x_3\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+z_3\mathbf{a}_3$	=	$x_3a\hat{\mathbf{x}}+\frac{1}{4}b\hat{\mathbf{y}}+z_3c\hat{\mathbf{z}}$	$(4f)$	O III
$\mathbf{B}_8$	=	$\left(\frac{1}{2}-x_3\right)\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+z_3\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_3\right)a\hat{\mathbf{x}}+\frac{1}{4}b\hat{\mathbf{y}}+z_3c\hat{\mathbf{z}}$	$(4f)$	O III
$\mathbf{B}_9$	=	$-x_3\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2-z_3\mathbf{a}_3$	=	$-x_3a\hat{\mathbf{x}}+\frac{3}{4}b\hat{\mathbf{y}}-z_3c\hat{\mathbf{z}}$	$(4f)$	O III
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2}+x_3\right)\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2-z_3\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_3\right)a\hat{\mathbf{x}}+\frac{3}{4}b\hat{\mathbf{y}}-z_3c\hat{\mathbf{z}}$	$(4f)$	O III
$\mathbf{B}_{11}$	=	$x_4\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+z_4\mathbf{a}_3$	=	$x_4a\hat{\mathbf{x}}+\frac{1}{4}b\hat{\mathbf{y}}+z_4c\hat{\mathbf{z}}$	$(4f)$	V
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2}-x_4\right)\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+z_4\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_4\right)a\hat{\mathbf{x}}+\frac{1}{4}b\hat{\mathbf{y}}+z_4c\hat{\mathbf{z}}$	$(4f)$	V
$\mathbf{B}_{13}$	=	$-x_4\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2-z_4\mathbf{a}_3$	=	$-x_4a\hat{\mathbf{x}}+\frac{3}{4}b\hat{\mathbf{y}}-z_4c\hat{\mathbf{z}}$	$(4f)$	V
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2}+x_4\right)\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2-z_4\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_4\right)a\hat{\mathbf{x}}+\frac{3}{4}b\hat{\mathbf{y}}-z_4c\hat{\mathbf{z}}$	$(4f)$	V

References:

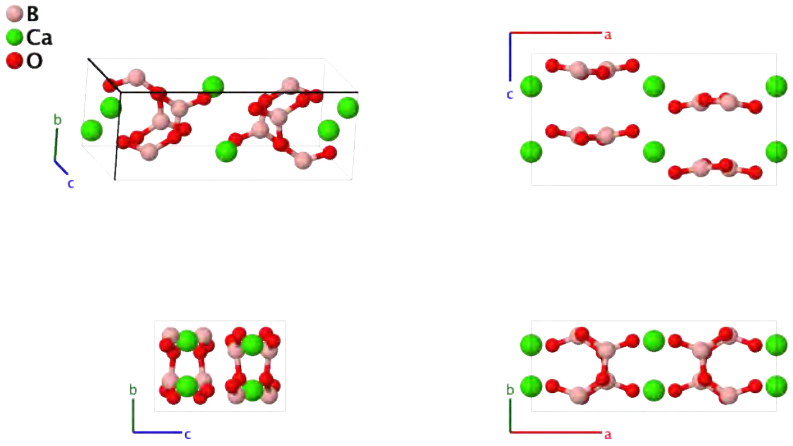
- R. Enjalbert and J. Galy, *A Refinement of the Structure of V_2O_5*, Acta Crystallogr. C **42**, 1467–1469 (1986), doi:10.1107/S0108270186091825.
- J. A. A. Ketelaar, *Crystal Structure and Shape of Colloidal Particles of Vanadium Pentoxide*, Nature **137**, 316 (1936), doi:10.1038/137316a0.
- C. Gottfried, ed., *Strukturbericht Band IV 1936* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1938).

Geometry files:

- CIF: pp. 1620
- POSCAR: pp. 1621

CaB_2O_4 I ($E3_2$) Structure: A2BC4_oP28_60_d_c_2d

http://aflow.org/prototype-encyclopedia/A2BC4_oP28_60_d_c_2d

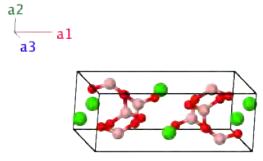


Prototype	:	B_2CaO_4
AFLOW prototype label	:	A2BC4_oP28_60_d_c_2d
***Strukturbericht* designation**	:	$E3_2$
Pearson symbol	:	oP28
Space group number	:	60
Space group symbol	:	*Pbcn*
AFLOW prototype command	:	`aflow --proto=A2BC4_oP28_60_d_c_2d --params=`$a, b/a, c/a, y_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4$

- CaB_2O_4 exists in at least four phases (Marezio, 1969):
- I - The ground state, stable below 1.2 GPa, *Strukturbericht* $E3_2$ (this structure).
- II – Orthorhombic high pressure structure, stable between 1.2 and 1.5 GPa, presumably calciborite.
- III – Orthorhombic high pressure structure, stable between 1.5 and 2.5 GPa.
- IV – Cubic high pressure structure, stable between 2.5 and 4.0 GPa.
- (Marezio, 1963) gives the crystal structure in the *Pnca* setting of space group #60. We have used FINDSYM to transform this to the standard *Pbcn* setting.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$y_1\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_1 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*c*)	Ca
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_1\right)\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_1\right)b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(4*c*)	Ca
$\mathbf{B}_3$	=	$-y_1\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_1 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(4*c*)	Ca
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_1\right)\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_1\right)b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*c*)	Ca
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(8*d*)	B
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	(8*d*)	B
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	(8*d*)	B
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right)b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(8*d*)	B
$\mathbf{B}_9$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(8*d*)	B
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	(8*d*)	B
$\mathbf{B}_{11}$	=	$x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	(8*d*)	B
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right)b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(8*d*)	B
$\mathbf{B}_{13}$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(8*d*)	O I
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(8*d*)	O I
$\mathbf{B}_{15}$	=	$-x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(8*d*)	O I
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(8*d*)	O I
$\mathbf{B}_{17}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(8*d*)	O I
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(8*d*)	O I
$\mathbf{B}_{19}$	=	$x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(8*d*)	O I
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(8*d*)	O I
$\mathbf{B}_{21}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{23}$	=	$-x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{25}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{27}$	=	$x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*d*)	O II

References:

- M. Marezio, H. A. Plettinger, and W. H. Zachariasen, *Refinement of the calcium metaborate structure*, Acta Cryst. **16**,

390–392 (1963), doi:10.1107/S0365110X63001031.

Found in:

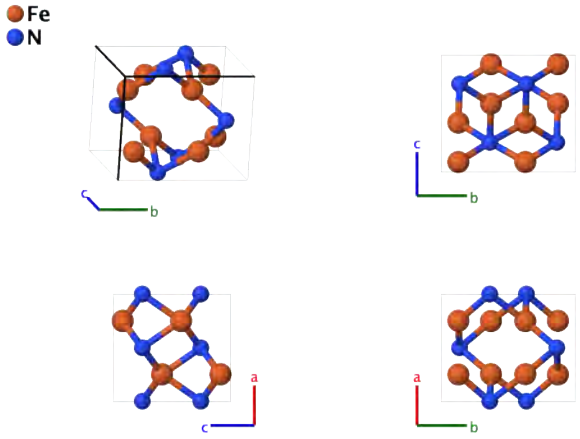
- M. Marezio, J. P. Remeika, and P. D. Dernier, *The crystal structure of the high-pressure phase CaB_2O_4(IV), and polymorphism in CaB_2O_4*, Acta Crystallogr. Sect. B Struct. Sci. **25**, 965–970 (1969), doi:10.1107/S0567740869003256.

Geometry files:

- CIF: pp. 1621
- POSCAR: pp. 1621

ζ-Fe_2N Structure: A2B_oP12_60_d_c

http://aflow.org/prototype-encyclopedia/A2B_oP12_60_d_c.Fe2N

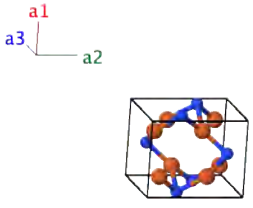


Prototype	:	Fe_2N
AFLOW prototype label	:	A2B_oP12_60_d_c
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP12
Space group number	:	60
Space group symbol	:	*Pbcn*
AFLOW prototype command	:	`aflow --proto=A2B_oP12_60_d_c --params=`$a, b/a, c/a, y_1, x_2, y_2, z_2$

- This structure has the same AFLOW label as α-PbO_2, but in that case the lead site is only 49% occupied, so the composition is actually closer to PbO_4.
- The structures are generated by the same symmetry operations with different sets of parameters (`--params`) specified in their corresponding CIF files.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$y_1\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_1 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($4c$)	N
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_1\right)\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_1\right)b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	($4c$)	N
$\mathbf{B}_3$	=	$-y_1\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_1 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	($4c$)	N

$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_1\right)\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_1\right)b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*c*)	N
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(8*d*)	Fe
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	(8*d*)	Fe
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	(8*d*)	Fe
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right)b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(8*d*)	Fe
$\mathbf{B}_9$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(8*d*)	Fe
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	(8*d*)	Fe
$\mathbf{B}_{11}$	=	$x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	(8*d*)	Fe
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right)b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(8*d*)	Fe

References:

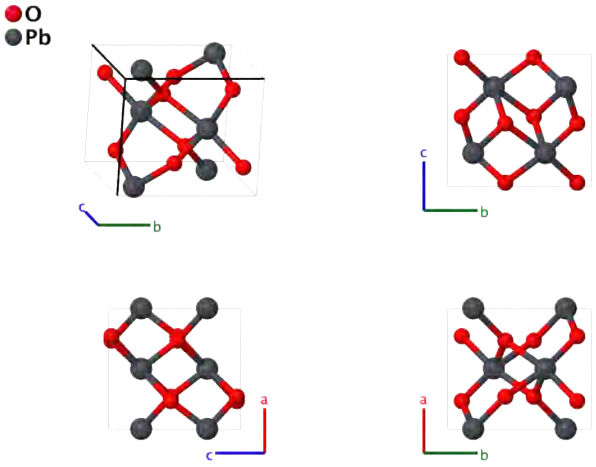
- D. Rechenbach and H. Jacobs, *Structure determination of ζ-Fe_2N by neutron and synchrotron powder diffraction*, J. Alloys Compd. **235**, 15–22 (1996), doi:10.1016/0925-8388(95)02097-7.

Geometry files:

- CIF: pp. 1621
- POSCAR: pp. 1622

α-PbO_2 Structure: A2B_oP12_60_d_c

http://aflow.org/prototype-encyclopedia/A2B_oP12_60_d_c



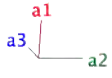
Prototype	:	O_2Pb
AFLOW prototype label	:	A2B_oP12_60_d_c
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP12
Space group number	:	60
Space group symbol	:	*Pbcn*
AFLOW prototype command	:	`aflow --proto=A2B_oP12_60_d_c --params=`$a, b/a, c/a, y_1, x_2, y_2, z_2$

Other compounds with this structure

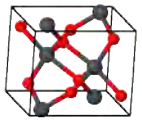
- (Ti,Zr)O_2 (srilankite)

- The experimental evidence shows that the Pb site is only 49% occupied, so stoichiometrically this compound is closer to PbO_4.
- This structure has the same AFLOW label as ζ-Fe_2N, but in that case the nitrogen site is fully occupied. The structures are generated by the same symmetry operations with different sets of parameters (`--params`) specified in their corresponding CIF files.

Simple Orthorhombic primitive vectors:



$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$y_1\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_1 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($4c$)	Pb
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_1\right)\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_1\right)b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	($4c$)	Pb
$\mathbf{B}_3$	=	$-y_1\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_1 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	($4c$)	Pb
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_1\right)\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_1\right)b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($4c$)	Pb
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	($8d$)	O
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	($8d$)	O
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	($8d$)	O
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right)b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	($8d$)	O
$\mathbf{B}_9$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	($8d$)	O
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	($8d$)	O
$\mathbf{B}_{11}$	=	$x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	($8d$)	O
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right)b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	($8d$)	O

References:

- R. J. Hill, *The Crystal Structures of Lead Dioxides from the Positive Plate of the Lead/Acid Battery*, Mater. Res. Bull. **17**, 769–784 (1982), doi:10.1016/0025-5408(82)90028-9.

Found in:

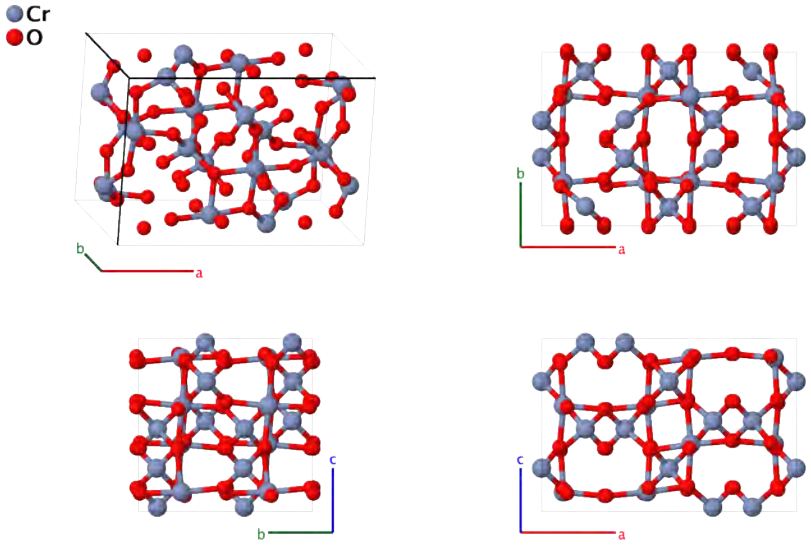
- P. Villars and L. Calvert, *Pearson's Handbook of Crystallographic Data for Intermetallic Phases* (ASM International, Materials Park, OK, 1991), vol. IV, chap. , p. 4745.

Geometry files:

- CIF: pp. 1622
- POSCAR: pp. 1622

Cr_5O_{12} Structure: A5B12_oP68_60_c2d_6d

http://aflow.org/prototype-encyclopedia/A5B12_oP68_60_c2d_6d

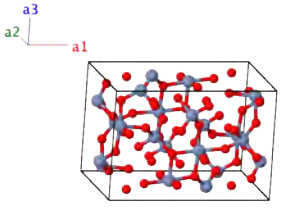


Prototype	:	Cr_5O_{12}
AFLOW prototype label	:	A5B12_oP68_60_c2d_6d
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP68
Space group number	:	60
Space group symbol	:	*Pbcn*
AFLOW prototype command	:	`aflow --proto=A5B12_oP68_60_c2d_6d --params=`$a, b/a, c/a, y_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9$

- (Wilhelmi, 1965) describes this as a distorted face-centered cubic lattice of oxygen atoms with the chromium atoms in interstitial sites. One can very clearly see the CrO_4 tetrahedra and CrO_6 octahedra, indicating the presence of both Cr^{2+} and Cr^{3+} ions.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$y_1\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_1 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($4c$)	Cr I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_1\right)\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_1\right)b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	($4c$)	Cr I

$\mathbf{B}_3$	=	$-y_1\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_1 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(4*c*)	Cr I
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_1\right)\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_1\right)b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*c*)	Cr I
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(8*d*)	Cr II
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	(8*d*)	Cr II
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	(8*d*)	Cr II
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right)b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(8*d*)	Cr II
$\mathbf{B}_9$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(8*d*)	Cr II
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	(8*d*)	Cr II
$\mathbf{B}_{11}$	=	$x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	(8*d*)	Cr II
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right)b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(8*d*)	Cr II
$\mathbf{B}_{13}$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(8*d*)	Cr III
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(8*d*)	Cr III
$\mathbf{B}_{15}$	=	$-x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(8*d*)	Cr III
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(8*d*)	Cr III
$\mathbf{B}_{17}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(8*d*)	Cr III
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(8*d*)	Cr III
$\mathbf{B}_{19}$	=	$x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(8*d*)	Cr III
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(8*d*)	Cr III
$\mathbf{B}_{21}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*d*)	O I
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(8*d*)	O I
$\mathbf{B}_{23}$	=	$-x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(8*d*)	O I
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*d*)	O I
$\mathbf{B}_{25}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*d*)	O I
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(8*d*)	O I
$\mathbf{B}_{27}$	=	$x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(8*d*)	O I
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*d*)	O I
$\mathbf{B}_{29}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{31}$	=	$-x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right)b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{33}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(8*d*)	O II

$\mathbf{B}_{34}$	=	$\left(\frac{1}{2}+x_5\right)\mathbf{a}_1+\left(\frac{1}{2}+y_5\right)\mathbf{a}_2+\left(\frac{1}{2}-z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_5\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_5\right)c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{35}$	=	$x_5\,\mathbf{a}_1-y_5\,\mathbf{a}_2+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}}-y_5b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{36}$	=	$\left(\frac{1}{2}-x_5\right)\mathbf{a}_1+\left(\frac{1}{2}+y_5\right)\mathbf{a}_2+z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_5\right)b\,\hat{\mathbf{y}}+z_5c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{37}$	=	$x_6\,\mathbf{a}_1+y_6\,\mathbf{a}_2+z_6\,\mathbf{a}_3$	=	$x_6a\,\hat{\mathbf{x}}+y_6b\,\hat{\mathbf{y}}+z_6c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{38}$	=	$\left(\frac{1}{2}-x_6\right)\mathbf{a}_1+\left(\frac{1}{2}-y_6\right)\mathbf{a}_2+\left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_6\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{39}$	=	$-x_6\,\mathbf{a}_1+y_6\,\mathbf{a}_2+\left(\frac{1}{2}-z_6\right)\mathbf{a}_3$	=	$-x_6a\,\hat{\mathbf{x}}+y_6b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_6\right)c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{40}$	=	$\left(\frac{1}{2}+x_6\right)\mathbf{a}_1+\left(\frac{1}{2}-y_6\right)\mathbf{a}_2-z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_6\right)b\,\hat{\mathbf{y}}-z_6c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{41}$	=	$-x_6\,\mathbf{a}_1-y_6\,\mathbf{a}_2-z_6\,\mathbf{a}_3$	=	$-x_6a\,\hat{\mathbf{x}}-y_6b\,\hat{\mathbf{y}}-z_6c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{42}$	=	$\left(\frac{1}{2}+x_6\right)\mathbf{a}_1+\left(\frac{1}{2}+y_6\right)\mathbf{a}_2+\left(\frac{1}{2}-z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_6\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_6\right)c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{43}$	=	$x_6\,\mathbf{a}_1-y_6\,\mathbf{a}_2+\left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	=	$x_6a\,\hat{\mathbf{x}}-y_6b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{44}$	=	$\left(\frac{1}{2}-x_6\right)\mathbf{a}_1+\left(\frac{1}{2}+y_6\right)\mathbf{a}_2+z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_6\right)b\,\hat{\mathbf{y}}+z_6c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{45}$	=	$x_7\,\mathbf{a}_1+y_7\,\mathbf{a}_2+z_7\,\mathbf{a}_3$	=	$x_7a\,\hat{\mathbf{x}}+y_7b\,\hat{\mathbf{y}}+z_7c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{46}$	=	$\left(\frac{1}{2}-x_7\right)\mathbf{a}_1+\left(\frac{1}{2}-y_7\right)\mathbf{a}_2+\left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_7\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_7\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{47}$	=	$-x_7\,\mathbf{a}_1+y_7\,\mathbf{a}_2+\left(\frac{1}{2}-z_7\right)\mathbf{a}_3$	=	$-x_7a\,\hat{\mathbf{x}}+y_7b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_7\right)c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{48}$	=	$\left(\frac{1}{2}+x_7\right)\mathbf{a}_1+\left(\frac{1}{2}-y_7\right)\mathbf{a}_2-z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_7\right)b\,\hat{\mathbf{y}}-z_7c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{49}$	=	$-x_7\,\mathbf{a}_1-y_7\,\mathbf{a}_2-z_7\,\mathbf{a}_3$	=	$-x_7a\,\hat{\mathbf{x}}-y_7b\,\hat{\mathbf{y}}-z_7c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{50}$	=	$\left(\frac{1}{2}+x_7\right)\mathbf{a}_1+\left(\frac{1}{2}+y_7\right)\mathbf{a}_2+\left(\frac{1}{2}-z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_7\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_7\right)c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{51}$	=	$x_7\,\mathbf{a}_1-y_7\,\mathbf{a}_2+\left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	=	$x_7a\,\hat{\mathbf{x}}-y_7b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{52}$	=	$\left(\frac{1}{2}-x_7\right)\mathbf{a}_1+\left(\frac{1}{2}+y_7\right)\mathbf{a}_2+z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_7\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_7\right)b\,\hat{\mathbf{y}}+z_7c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{53}$	=	$x_8\,\mathbf{a}_1+y_8\,\mathbf{a}_2+z_8\,\mathbf{a}_3$	=	$x_8a\,\hat{\mathbf{x}}+y_8b\,\hat{\mathbf{y}}+z_8c\,\hat{\mathbf{z}}$	(8*d*)	O V
$\mathbf{B}_{54}$	=	$\left(\frac{1}{2}-x_8\right)\mathbf{a}_1+\left(\frac{1}{2}-y_8\right)\mathbf{a}_2+\left(\frac{1}{2}+z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_8\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_8\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_8\right)c\,\hat{\mathbf{z}}$	(8*d*)	O V
$\mathbf{B}_{55}$	=	$-x_8\,\mathbf{a}_1+y_8\,\mathbf{a}_2+\left(\frac{1}{2}-z_8\right)\mathbf{a}_3$	=	$-x_8a\,\hat{\mathbf{x}}+y_8b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_8\right)c\,\hat{\mathbf{z}}$	(8*d*)	O V
$\mathbf{B}_{56}$	=	$\left(\frac{1}{2}+x_8\right)\mathbf{a}_1+\left(\frac{1}{2}-y_8\right)\mathbf{a}_2-z_8\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_8\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_8\right)b\,\hat{\mathbf{y}}-z_8c\,\hat{\mathbf{z}}$	(8*d*)	O V
$\mathbf{B}_{57}$	=	$-x_8\,\mathbf{a}_1-y_8\,\mathbf{a}_2-z_8\,\mathbf{a}_3$	=	$-x_8a\,\hat{\mathbf{x}}-y_8b\,\hat{\mathbf{y}}-z_8c\,\hat{\mathbf{z}}$	(8*d*)	O V
$\mathbf{B}_{58}$	=	$\left(\frac{1}{2}+x_8\right)\mathbf{a}_1+\left(\frac{1}{2}+y_8\right)\mathbf{a}_2+\left(\frac{1}{2}-z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_8\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_8\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_8\right)c\,\hat{\mathbf{z}}$	(8*d*)	O V
$\mathbf{B}_{59}$	=	$x_8\,\mathbf{a}_1-y_8\,\mathbf{a}_2+\left(\frac{1}{2}+z_8\right)\mathbf{a}_3$	=	$x_8a\,\hat{\mathbf{x}}-y_8b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_8\right)c\,\hat{\mathbf{z}}$	(8*d*)	O V
$\mathbf{B}_{60}$	=	$\left(\frac{1}{2}-x_8\right)\mathbf{a}_1+\left(\frac{1}{2}+y_8\right)\mathbf{a}_2+z_8\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_8\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_8\right)b\,\hat{\mathbf{y}}+z_8c\,\hat{\mathbf{z}}$	(8*d*)	O V
$\mathbf{B}_{61}$	=	$x_9\,\mathbf{a}_1+y_9\,\mathbf{a}_2+z_9\,\mathbf{a}_3$	=	$x_9a\,\hat{\mathbf{x}}+y_9b\,\hat{\mathbf{y}}+z_9c\,\hat{\mathbf{z}}$	(8*d*)	O VI
$\mathbf{B}_{62}$	=	$\left(\frac{1}{2}-x_9\right)\mathbf{a}_1+\left(\frac{1}{2}-y_9\right)\mathbf{a}_2+\left(\frac{1}{2}+z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_9\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_9\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_9\right)c\,\hat{\mathbf{z}}$	(8*d*)	O VI
$\mathbf{B}_{63}$	=	$-x_9\,\mathbf{a}_1+y_9\,\mathbf{a}_2+\left(\frac{1}{2}-z_9\right)\mathbf{a}_3$	=	$-x_9a\,\hat{\mathbf{x}}+y_9b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_9\right)c\,\hat{\mathbf{z}}$	(8*d*)	O VI

$\mathbf{B}_{64}$	=	$\left(\frac{1}{2}+x_9\right)\mathbf{a}_1+\left(\frac{1}{2}-y_9\right)\mathbf{a}_2-z_9\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_9\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_9\right)b\,\hat{\mathbf{y}}-z_9c\,\hat{\mathbf{z}}$	(8*d*)	O VI
$\mathbf{B}_{65}$	=	$-x_9\,\mathbf{a}_1-y_9\,\mathbf{a}_2-z_9\,\mathbf{a}_3$	=	$-x_9a\,\hat{\mathbf{x}}-y_9b\,\hat{\mathbf{y}}-z_9c\,\hat{\mathbf{z}}$	(8*d*)	O VI
$\mathbf{B}_{66}$	=	$\left(\frac{1}{2}+x_9\right)\mathbf{a}_1+\left(\frac{1}{2}+y_9\right)\mathbf{a}_2+\left(\frac{1}{2}-z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_9\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_9\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_9\right)c\,\hat{\mathbf{z}}$	(8*d*)	O VI
$\mathbf{B}_{67}$	=	$x_9\,\mathbf{a}_1-y_9\,\mathbf{a}_2+\left(\frac{1}{2}+z_9\right)\mathbf{a}_3$	=	$x_9a\,\hat{\mathbf{x}}-y_9b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_9\right)c\,\hat{\mathbf{z}}$	(8*d*)	O VI
$\mathbf{B}_{68}$	=	$\left(\frac{1}{2}-x_9\right)\mathbf{a}_1+\left(\frac{1}{2}+y_9\right)\mathbf{a}_2+z_9\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_9\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_9\right)b\,\hat{\mathbf{y}}+z_9c\,\hat{\mathbf{z}}$	(8*d*)	O VI

References:

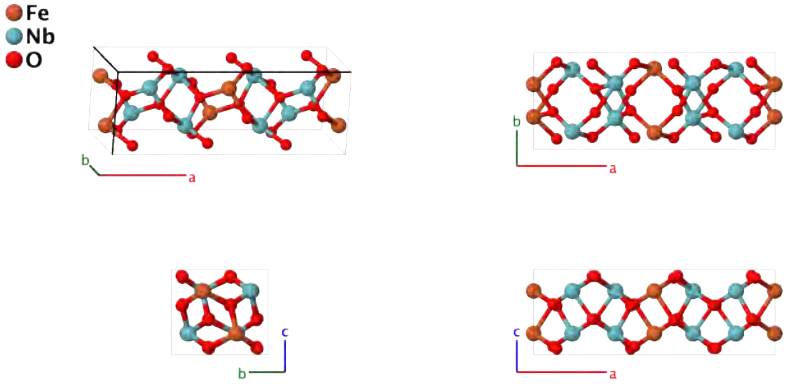
- K.-A. Wilhelmi, *The Crystal Structure of* Cr_5O_{12}, Acta Chem. Scand. **19**, 165–176 (1965), doi:10.3891/acta.chem.scand.19-0165.

Geometry files:

- CIF: pp. 1622
- POSCAR: pp. 1623

Columbite ($FeNb_2O_6$, $E5_1$) Structure: AB2C6_oP36_60_c_d_3d

http://aflow.org/prototype-encyclopedia/AB2C6_oP36_60_c_d_3d



Prototype	:	$FeNb_2O_6$
AFLOW prototype label	:	AB2C6_oP36_60_c_d_3d
***Strukturbericht* designation**	:	$E5_1$
Pearson symbol	:	oP36
Space group number	:	60
Space group symbol	:	*Pbcn*
AFLOW prototype command	:	`aflow --proto=AB2C6_oP36_60_c_d_3d --params=`$a, b/a, c/a, y_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5$

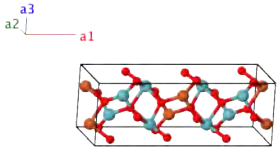
Other compounds with this structure

- $CaNb_2O_6$ (fermsmite) and $Zr_5Ti_7O_{24}$

- The original *Strukturbericht* reference uses $(Fe,Mn)Nb_2O_6$ as the prototype for columbite/$E5_1$.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$y_1\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_1 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($4c$)	Fe
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_1\right)\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_1\right)b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	($4c$)	Fe
$\mathbf{B}_3$	=	$-y_1\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_1 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	($4c$)	Fe
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_1\right)\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_1\right)b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($4c$)	Fe

$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(8*d*)	Nb
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	(8*d*)	Nb
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	(8*d*)	Nb
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right)b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(8*d*)	Nb
$\mathbf{B}_9$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(8*d*)	Nb
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	(8*d*)	Nb
$\mathbf{B}_{11}$	=	$x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	(8*d*)	Nb
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right)b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(8*d*)	Nb
$\mathbf{B}_{13}$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(8*d*)	O I
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(8*d*)	O I
$\mathbf{B}_{15}$	=	$-x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(8*d*)	O I
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(8*d*)	O I
$\mathbf{B}_{17}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(8*d*)	O I
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(8*d*)	O I
$\mathbf{B}_{19}$	=	$x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(8*d*)	O I
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(8*d*)	O I
$\mathbf{B}_{21}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{23}$	=	$-x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{25}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{27}$	=	$x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{29}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{31}$	=	$-x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right)b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{33}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{34}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	(8*d*)	O III

$$\mathbf{B}_{35} = x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_3 = x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_5\right) c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O III}$$

$$\mathbf{B}_{36} = \left(\tfrac{1}{2} - x_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_5\right)\mathbf{a}_2 + z_5\,\mathbf{a}_3 = \left(\tfrac{1}{2} - x_5\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_5\right) b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O III}$$

References:

- P. Bordet, A. McHale, A. Santoro, and R. S. Roth, *Powder neutron diffraction study of $ZrTiO_4$, $Zr_5Ti_7O_{24}$, and $FeNb_2O_6$*, J. Solid State Chem. **64**, 30–46 (1986), doi:10.1016/0022-4596(86)90119-2.
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Found in:

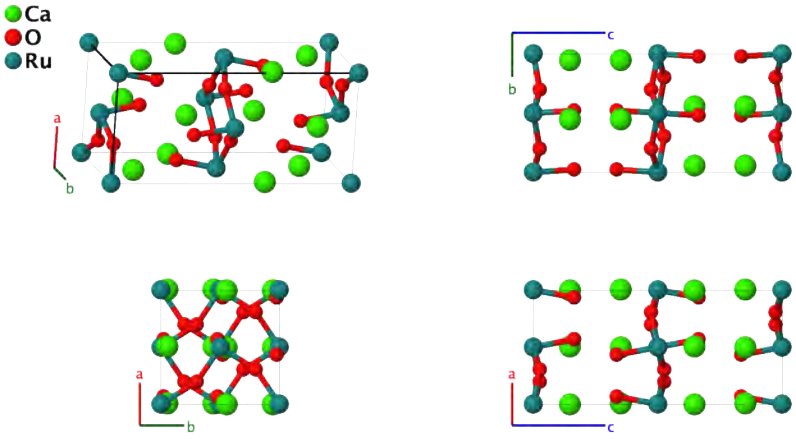
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1623
- POSCAR: pp. 1624

Ca_2RuO_4 Structure: A2B4C_oP28_61_c_2c_a

http://aflow.org/prototype-encyclopedia/A2B4C_oP28_61_c_2c_a



Prototype	:	Ca_2O_4Ru
AFLOW prototype label	:	A2B4C_oP28_61_c_2c_a
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP28
Space group number	:	61
Space group symbol	:	*Pbca*
AFLOW prototype command	:	`aflow --proto=A2B4C_oP28_61_c_2c_a --params=`$a, b/a, c/a, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4$

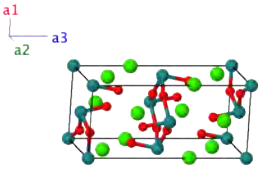
Other compounds with this structure

- $Ca_{2-x}Sr_xRuO_4$

- (Friedt, 2001) never give the positions of the ruthenium atoms in this structure, however the composition of the crystal dictates that they can only be at the (4*a*) or (4*b*) Wyckoff positions, and the Ru-O distances given are consistent with the (4*a*) site.
- The authors identify a low temperature "$S-Pbca$" phase and a high temperature "$L-Pbca$" phase, with a phase transition in the range 350-400 K. The major difference between the two phases is a 4% elongation of the c axis in the $L-Pbca$ phase, accompanied by a 4.5% contraction along the b axis, with the a axis being substantially unchanged. There is also a substantial change in the tilt of the oxygen octahedra surrounding the Ru atoms. Here we show the crystal in the $S-Pbca$ phase using data taken at 180 K.
- (Bertinshaw, 2019) identify S^* and L^* phases that have the same symmetry but stabilized by the application of an electric current.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(4*a*)	Ru
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*a*)	Ru
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*a*)	Ru
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$	(4*a*)	Ru
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2a\,\hat{\mathbf{x}} + y_2b\,\hat{\mathbf{y}} + z_2c\,\hat{\mathbf{z}}$	(8*c*)	Ca
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} - y_2b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	(8*c*)	Ca
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	(8*c*)	Ca
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right)b\,\hat{\mathbf{y}} - z_2c\,\hat{\mathbf{z}}$	(8*c*)	Ca
$\mathbf{B}_9$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}} - y_2b\,\hat{\mathbf{y}} - z_2c\,\hat{\mathbf{z}}$	(8*c*)	Ca
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + y_2b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	(8*c*)	Ca
$\mathbf{B}_{11}$	=	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$x_2a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	(8*c*)	Ca
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right)b\,\hat{\mathbf{y}} + z_2c\,\hat{\mathbf{z}}$	(8*c*)	Ca
$\mathbf{B}_{13}$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + y_3b\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	(8*c*)	O I
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} - y_3b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(8*c*)	O I
$\mathbf{B}_{15}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(8*c*)	O I
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)b\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	(8*c*)	O I
$\mathbf{B}_{17}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - y_3b\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	(8*c*)	O I
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + y_3b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(8*c*)	O I
$\mathbf{B}_{19}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(8*c*)	O I
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)b\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	(8*c*)	O I
$\mathbf{B}_{21}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + y_4b\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	(8*c*)	O II
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} - y_4b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(8*c*)	O II
$\mathbf{B}_{23}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(8*c*)	O II
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)b\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	(8*c*)	O II
$\mathbf{B}_{25}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} - y_4b\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	(8*c*)	O II
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + y_4b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(8*c*)	O II
$\mathbf{B}_{27}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(8*c*)	O II
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)b\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	(8*c*)	O II

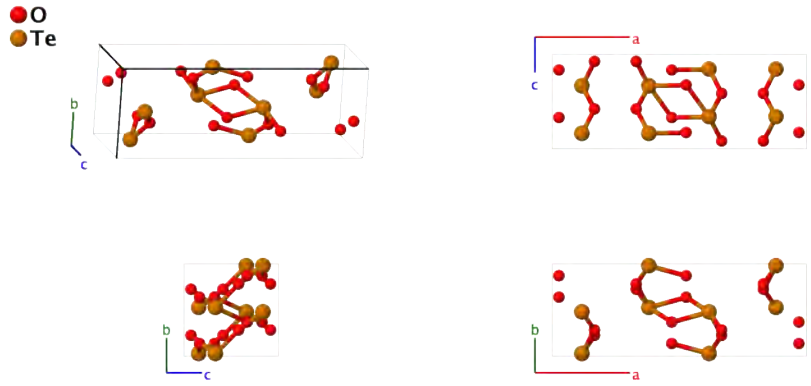
References:
- O. Friedt, M. Braden, G. André, P. Adelmann, S. Nakatsuji, and Y. Maeno, *Structural and magnetic aspects of the metal-insulator transition in $Ca_{2-x}Sr_xRuO_4$*, Phys. Rev. B **63**, 174432 (2001), doi:10.1103/PhysRevB.63.174432.

Found in:
- J. Bertinshaw, N. Gurung, P. Jorba, H. Liu, M. Schmid, D. T. Mantadakis, M. Daghofer, M. Krautloher, A. Jain, G. H. Ryu, O. Fabelo, P. Hansmann, G. Khaliullin, C. Pfleiderer, B. Keimer, and B. J. Kim, *Unique Crystal Structure of Ca_2RuO_4 in the Current Stabilized Semimetallic State*, Phys. Rev. Lett. **123**, 137204 (2019), doi:10.1103/PhysRevLett.123.137204.

Geometry files:
- CIF: pp. 1624
- POSCAR: pp. 1624

Tellurite (β-TeO_2, $C52$) Structure: A2B_oP24_61_2c_c

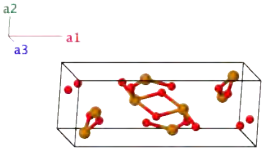
http://aflow.org/prototype-encyclopedia/A2B_oP24_61_2c_c.TeO2



Prototype	:	O_2Te
AFLOW prototype label	:	A2B_oP24_61_2c_c
***Strukturbericht* designation**	:	$C52$
Pearson symbol	:	oP24
Space group number	:	61
Space group symbol	:	*Pbca*
AFLOW prototype command	:	`aflow --proto=A2B_oP24_61_2c_c --params=`$a, b/a, c/a, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3$

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	($8c$)	O I
$\mathbf{B}_2$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right)a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\,\hat{\mathbf{z}}$	($8c$)	O I
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2} + y_1\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right)c\,\hat{\mathbf{z}}$	($8c$)	O I
$\mathbf{B}_4$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_1\right)\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_1\right)b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	($8c$)	O I
$\mathbf{B}_5$	=	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	($8c$)	O I
$\mathbf{B}_6$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + y_1\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right)a\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right)c\,\hat{\mathbf{z}}$	($8c$)	O I
$\mathbf{B}_7$	=	$x_1\,\mathbf{a}_1 + \left(\frac{1}{2} - y_1\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\,\hat{\mathbf{z}}$	($8c$)	O I
$\mathbf{B}_8$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_1\right)\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_1\right)b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	($8c$)	O I
$\mathbf{B}_9$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	($8c$)	O II

$$\mathbf{B}_{10} = \left(\tfrac{1}{2} - x_2\right)\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_2\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_2\right)c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O II}$$

$$\mathbf{B}_{11} = -x_2\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_2\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_2\right)\mathbf{a}_3 = -x_2 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_2\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_2\right)c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O II}$$

$$\mathbf{B}_{12} = \left(\tfrac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_2\right)\mathbf{a}_2 - z_2\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_2\right)b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O II}$$

$$\mathbf{B}_{13} = -x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3 = -x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O II}$$

$$\mathbf{B}_{14} = \left(\tfrac{1}{2} + x_2\right)\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_2\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_2\right)c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O II}$$

$$\mathbf{B}_{15} = x_2\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_2\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_2\right)\mathbf{a}_3 = x_2 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_2\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_2\right)c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O II}$$

$$\mathbf{B}_{16} = \left(\tfrac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_2\right)\mathbf{a}_2 + z_2\,\mathbf{a}_3 = \left(\tfrac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_2\right)b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}} \quad (8c) \quad \text{O II}$$

$$\mathbf{B}_{17} = x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3 = x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}} \quad (8c) \quad \text{Te}$$

$$\mathbf{B}_{18} = \left(\tfrac{1}{2} - x_3\right)\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_3\right)c\,\hat{\mathbf{z}} \quad (8c) \quad \text{Te}$$

$$\mathbf{B}_{19} = -x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_3\right)\mathbf{a}_3 = -x_3 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_3\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_3\right)c\,\hat{\mathbf{z}} \quad (8c) \quad \text{Te}$$

$$\mathbf{B}_{20} = \left(\tfrac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_3\right)\mathbf{a}_2 - z_3\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_3\right)b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}} \quad (8c) \quad \text{Te}$$

$$\mathbf{B}_{21} = -x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3 = -x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}} \quad (8c) \quad \text{Te}$$

$$\mathbf{B}_{22} = \left(\tfrac{1}{2} + x_3\right)\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_3\right)c\,\hat{\mathbf{z}} \quad (8c) \quad \text{Te}$$

$$\mathbf{B}_{23} = x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_3\right)\mathbf{a}_3 = x_3 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_3\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_3\right)c\,\hat{\mathbf{z}} \quad (8c) \quad \text{Te}$$

$$\mathbf{B}_{24} = \left(\tfrac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_3\right)\mathbf{a}_2 + z_3\,\mathbf{a}_3 = \left(\tfrac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_3\right)b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}} \quad (8c) \quad \text{Te}$$

References:

- H. Beyer, *Verfeinerung der Kristallstruktur von Tellurit, dem rhombischen TeO_2*, Zeitschrift für Kristallographie - Crystalline Materials **124**, 228–237 (1967), doi:10.1524/zkri.1967.124.3.228.

Found in:

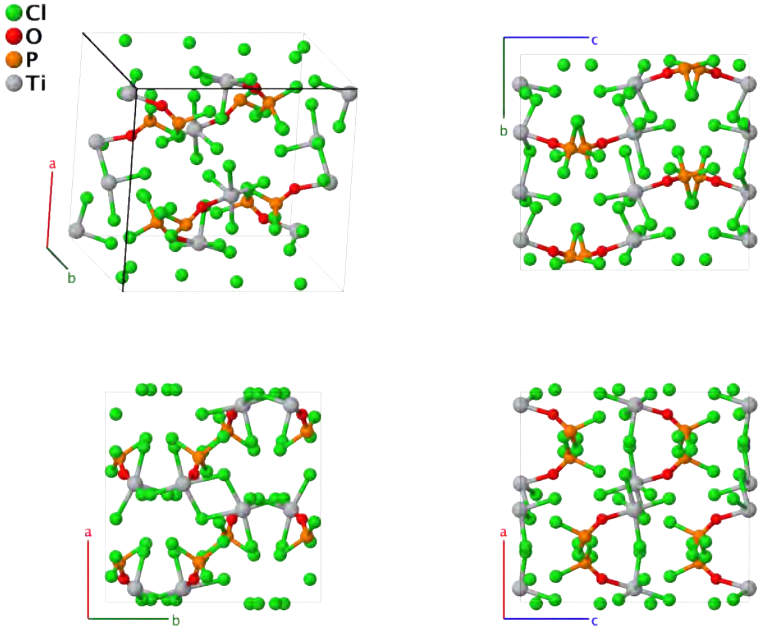
- M. Ceriotti, F. Pietrucci, and M. Bernasconi, Ab initio *study of the vibrational properties of crystalline TeO_2: The α, β, and γ phases*, Phys. Rev. B **73**, 104304 (2006), doi:10.1103/PhysRevB.73.104304.

Geometry files:

- CIF: pp. 1624
- POSCAR: pp. 1625

$(TiCl_4{\cdot}POCl_3)_2$ Structure: A7BCD_oP80_61_7c_c_c_c

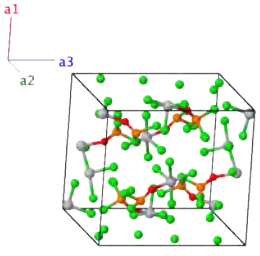
http://aflow.org/prototype-encyclopedia/A7BCD_oP80_61_7c_c_c_c



Prototype	:	Cl_7OPTi
AFLOW prototype label	:	A7BCD_oP80_61_7c_c_c_c
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP80
Space group number	:	61
Space group symbol	:	*Pbca*
AFLOW prototype command	:	`aflow --proto=A7BCD_oP80_61_7c_c_c_c --params=`$a, b/a, c/a, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}$

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$ =	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	($8c$)	Cl I

$$
\begin{aligned}
\mathbf{B}_{2} &= \left(\tfrac{1}{2}-x_1\right)\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_1\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_1\right)a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_1\right)c\,\hat{\mathbf{z}} && (8c) && \text{Cl I}\\
\mathbf{B}_{3} &= -x_1\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_1\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_1\right)\mathbf{a}_3 &&= -x_1 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_1\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_1\right)c\,\hat{\mathbf{z}} && (8c) && \text{Cl I}\\
\mathbf{B}_{4} &= \left(\tfrac{1}{2}+x_1\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_1\right)\mathbf{a}_2 - z_1\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_1\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_1\right)b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}} && (8c) && \text{Cl I}\\
\mathbf{B}_{5} &= -x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3 &&= -x_1 a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}} && (8c) && \text{Cl I}\\
\mathbf{B}_{6} &= \left(\tfrac{1}{2}+x_1\right)\mathbf{a}_1 + y_1\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_1\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_1\right)a\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_1\right)c\,\hat{\mathbf{z}} && (8c) && \text{Cl I}\\
\mathbf{B}_{7} &= x_1\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_1\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_1\right)\mathbf{a}_3 &&= x_1 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_1\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_1\right)c\,\hat{\mathbf{z}} && (8c) && \text{Cl I}\\
\mathbf{B}_{8} &= \left(\tfrac{1}{2}-x_1\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_1\right)\mathbf{a}_2 + z_1\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_1\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_1\right)b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}} && (8c) && \text{Cl I}\\
\mathbf{B}_{9} &= x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3 &&= x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}} && (8c) && \text{Cl II}\\
\mathbf{B}_{10} &= \left(\tfrac{1}{2}-x_2\right)\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_2\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_2\right)a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_2\right)c\,\hat{\mathbf{z}} && (8c) && \text{Cl II}\\
\mathbf{B}_{11} &= -x_2\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_2\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_2\right)\mathbf{a}_3 &&= -x_2 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_2\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_2\right)c\,\hat{\mathbf{z}} && (8c) && \text{Cl II}\\
\mathbf{B}_{12} &= \left(\tfrac{1}{2}+x_2\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_2\right)\mathbf{a}_2 - z_2\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_2\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_2\right)b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}} && (8c) && \text{Cl II}\\
\mathbf{B}_{13} &= -x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3 &&= -x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}} && (8c) && \text{Cl II}\\
\mathbf{B}_{14} &= \left(\tfrac{1}{2}+x_2\right)\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_2\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_2\right)a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_2\right)c\,\hat{\mathbf{z}} && (8c) && \text{Cl II}\\
\mathbf{B}_{15} &= x_2\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_2\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_2\right)\mathbf{a}_3 &&= x_2 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_2\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_2\right)c\,\hat{\mathbf{z}} && (8c) && \text{Cl II}\\
\mathbf{B}_{16} &= \left(\tfrac{1}{2}-x_2\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_2\right)\mathbf{a}_2 + z_2\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_2\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_2\right)b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}} && (8c) && \text{Cl II}\\
\mathbf{B}_{17} &= x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3 &&= x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}} && (8c) && \text{Cl III}\\
\mathbf{B}_{18} &= \left(\tfrac{1}{2}-x_3\right)\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_3\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_3\right)a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_3\right)c\,\hat{\mathbf{z}} && (8c) && \text{Cl III}\\
\mathbf{B}_{19} &= -x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_3\right)\mathbf{a}_3 &&= -x_3 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_3\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_3\right)c\,\hat{\mathbf{z}} && (8c) && \text{Cl III}\\
\mathbf{B}_{20} &= \left(\tfrac{1}{2}+x_3\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_3\right)\mathbf{a}_2 - z_3\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_3\right)b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}} && (8c) && \text{Cl III}\\
\mathbf{B}_{21} &= -x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3 &&= -x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}} && (8c) && \text{Cl III}\\
\mathbf{B}_{22} &= \left(\tfrac{1}{2}+x_3\right)\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_3\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_3\right)a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_3\right)c\,\hat{\mathbf{z}} && (8c) && \text{Cl III}\\
\mathbf{B}_{23} &= x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_3\right)\mathbf{a}_3 &&= x_3 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_3\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_3\right)c\,\hat{\mathbf{z}} && (8c) && \text{Cl III}\\
\mathbf{B}_{24} &= \left(\tfrac{1}{2}-x_3\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_3\right)\mathbf{a}_2 + z_3\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_3\right)b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}} && (8c) && \text{Cl III}\\
\mathbf{B}_{25} &= x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3 &&= x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}} && (8c) && \text{Cl IV}\\
\mathbf{B}_{26} &= \left(\tfrac{1}{2}-x_4\right)\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_4\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_4\right)a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_4\right)c\,\hat{\mathbf{z}} && (8c) && \text{Cl IV}\\
\mathbf{B}_{27} &= -x_4\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_4\right)\mathbf{a}_3 &&= -x_4 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_4\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_4\right)c\,\hat{\mathbf{z}} && (8c) && \text{Cl IV}\\
\mathbf{B}_{28} &= \left(\tfrac{1}{2}+x_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_4\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_4\right)b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}} && (8c) && \text{Cl IV}\\
\mathbf{B}_{29} &= -x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3 &&= -x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}} && (8c) && \text{Cl IV}\\
\mathbf{B}_{30} &= \left(\tfrac{1}{2}+x_4\right)\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_4\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_4\right)a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_4\right)c\,\hat{\mathbf{z}} && (8c) && \text{Cl IV}\\
\mathbf{B}_{31} &= x_4\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_4\right)\mathbf{a}_3 &&= x_4 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_4\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_4\right)c\,\hat{\mathbf{z}} && (8c) && \text{Cl IV}\\
\mathbf{B}_{32} &= \left(\tfrac{1}{2}-x_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_4\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_4\right)b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}} && (8c) && \text{Cl IV}\\
\mathbf{B}_{33} &= x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3 &&= x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}} && (8c) && \text{Cl V}\\
\mathbf{B}_{34} &= \left(\tfrac{1}{2}-x_5\right)\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_5\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_5\right)a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_5\right)c\,\hat{\mathbf{z}} && (8c) && \text{Cl V}\\
\mathbf{B}_{35} &= -x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_5\right)\mathbf{a}_3 &&= -x_5 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_5\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_5\right)c\,\hat{\mathbf{z}} && (8c) && \text{Cl V}\\
\mathbf{B}_{36} &= \left(\tfrac{1}{2}+x_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_5\right)b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}} && (8c) && \text{Cl V}\\
\mathbf{B}_{37} &= -x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3 &&= -x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}} && (8c) && \text{Cl V}
\end{aligned}
$$

$$
\begin{aligned}
\mathbf{B}_{38} &= \left(\tfrac{1}{2}+x_5\right)\mathbf{a}_1 + y_5\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_5\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_5\right)a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_5\right)c\,\hat{\mathbf{z}} && (8c) && \text{Cl V}\\
\mathbf{B}_{39} &= x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_5\right)\mathbf{a}_3 &&= x_5 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_5\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_5\right)c\,\hat{\mathbf{z}} && (8c) && \text{Cl V}\\
\mathbf{B}_{40} &= \left(\tfrac{1}{2}-x_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_5\right)\mathbf{a}_2 + z_5\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_5\right)b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}} && (8c) && \text{Cl V}\\
\mathbf{B}_{41} &= x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 &&= x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}} && (8c) && \text{Cl VI}\\
\mathbf{B}_{42} &= \left(\tfrac{1}{2}-x_6\right)\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_6\right)c\,\hat{\mathbf{z}} && (8c) && \text{Cl VI}\\
\mathbf{B}_{43} &= -x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_6\right)\mathbf{a}_3 &&= -x_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_6\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_6\right)c\,\hat{\mathbf{z}} && (8c) && \text{Cl VI}\\
\mathbf{B}_{44} &= \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_6\right)b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}} && (8c) && \text{Cl VI}\\
\mathbf{B}_{45} &= -x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3 &&= -x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}} && (8c) && \text{Cl VI}\\
\mathbf{B}_{46} &= \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_1 + y_6\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_6\right)c\,\hat{\mathbf{z}} && (8c) && \text{Cl VI}\\
\mathbf{B}_{47} &= x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_6\right)\mathbf{a}_3 &&= x_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_6\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_6\right)c\,\hat{\mathbf{z}} && (8c) && \text{Cl VI}\\
\mathbf{B}_{48} &= \left(\tfrac{1}{2}-x_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_6\right)\mathbf{a}_2 + z_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_6\right)b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}} && (8c) && \text{Cl VI}\\
\mathbf{B}_{49} &= x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 &&= x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}} && (8c) && \text{Cl VII}\\
\mathbf{B}_{50} &= \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_7\right)c\,\hat{\mathbf{z}} && (8c) && \text{Cl VII}\\
\mathbf{B}_{51} &= -x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_3 &&= -x_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_7\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_7\right)c\,\hat{\mathbf{z}} && (8c) && \text{Cl VII}\\
\mathbf{B}_{52} &= \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_7\right)\mathbf{a}_2 - z_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_7\right)b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}} && (8c) && \text{Cl VII}\\
\mathbf{B}_{53} &= -x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3 &&= -x_7 a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}} && (8c) && \text{Cl VII}\\
\mathbf{B}_{54} &= \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1 + y_7\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_7\right)c\,\hat{\mathbf{z}} && (8c) && \text{Cl VII}\\
\mathbf{B}_{55} &= x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 &&= x_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_7\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_7\right)c\,\hat{\mathbf{z}} && (8c) && \text{Cl VII}\\
\mathbf{B}_{56} &= \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_7\right)\mathbf{a}_2 + z_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_7\right)b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}} && (8c) && \text{Cl VII}\\
\mathbf{B}_{57} &= x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 &&= x_8 a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}} && (8c) && \text{O}\\
\mathbf{B}_{58} &= \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_1 - y_8\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_8\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_8\right)c\,\hat{\mathbf{z}} && (8c) && \text{O}\\
\mathbf{B}_{59} &= -x_8\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_8\right)\mathbf{a}_3 &&= -x_8 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_8\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_8\right)c\,\hat{\mathbf{z}} && (8c) && \text{O}\\
\mathbf{B}_{60} &= \left(\tfrac{1}{2}+x_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_8\right)\mathbf{a}_2 - z_8\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_8\right)b\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}} && (8c) && \text{O}\\
\mathbf{B}_{61} &= -x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3 &&= -x_8 a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}} && (8c) && \text{O}\\
\mathbf{B}_{62} &= \left(\tfrac{1}{2}+x_8\right)\mathbf{a}_1 + y_8\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_8\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_8\right)c\,\hat{\mathbf{z}} && (8c) && \text{O}\\
\mathbf{B}_{63} &= x_8\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_8\right)\mathbf{a}_3 &&= x_8 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_8\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_8\right)c\,\hat{\mathbf{z}} && (8c) && \text{O}\\
\mathbf{B}_{64} &= \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_8\right)\mathbf{a}_2 + z_8\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_8\right)b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}} && (8c) && \text{O}\\
\mathbf{B}_{65} &= x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 &&= x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}} && (8c) && \text{P}\\
\mathbf{B}_{66} &= \left(\tfrac{1}{2}-x_9\right)\mathbf{a}_1 - y_9\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_9\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_9\right)a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_9\right)c\,\hat{\mathbf{z}} && (8c) && \text{P}\\
\mathbf{B}_{67} &= -x_9\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_9\right)\mathbf{a}_3 &&= -x_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_9\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_9\right)c\,\hat{\mathbf{z}} && (8c) && \text{P}\\
\mathbf{B}_{68} &= \left(\tfrac{1}{2}+x_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_9\right)\mathbf{a}_2 - z_9\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_9\right)b\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}} && (8c) && \text{P}\\
\mathbf{B}_{69} &= -x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3 &&= -x_9 a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}} && (8c) && \text{P}\\
\mathbf{B}_{70} &= \left(\tfrac{1}{2}+x_9\right)\mathbf{a}_1 + y_9\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_9\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_9\right)a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_9\right)c\,\hat{\mathbf{z}} && (8c) && \text{P}\\
\mathbf{B}_{71} &= x_9\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_9\right)\mathbf{a}_3 &&= x_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_9\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_9\right)c\,\hat{\mathbf{z}} && (8c) && \text{P}\\
\mathbf{B}_{72} &= \left(\tfrac{1}{2}-x_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_9\right)\mathbf{a}_2 + z_9\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_9\right)b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}} && (8c) && \text{P}\\
\mathbf{B}_{73} &= x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 &&= x_{10} a\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}} && (8c) && \text{Ti}
\end{aligned}
$$

$\mathbf{B}_{74}$	=	$\left(\frac{1}{2}-x_{10}\right)\mathbf{a}_1-y_{10}\mathbf{a}_2+\left(\frac{1}{2}+z_{10}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{x}}-y_{10}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{10}\right)c\,\hat{\mathbf{z}}$	(8*c*)	Ti
$\mathbf{B}_{75}$	=	$-x_{10}\mathbf{a}_1+\left(\frac{1}{2}+y_{10}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{10}\right)\mathbf{a}_3$	=	$-x_{10}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{10}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{10}\right)c\,\hat{\mathbf{z}}$	(8*c*)	Ti
$\mathbf{B}_{76}$	=	$\left(\frac{1}{2}+x_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{10}\right)\mathbf{a}_2-z_{10}\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{10}\right)b\,\hat{\mathbf{y}}-z_{10}c\,\hat{\mathbf{z}}$	(8*c*)	Ti
$\mathbf{B}_{77}$	=	$-x_{10}\mathbf{a}_1-y_{10}\mathbf{a}_2-z_{10}\mathbf{a}_3$	=	$-x_{10}a\,\hat{\mathbf{x}}-y_{10}b\,\hat{\mathbf{y}}-z_{10}c\,\hat{\mathbf{z}}$	(8*c*)	Ti
$\mathbf{B}_{78}$	=	$\left(\frac{1}{2}+x_{10}\right)\mathbf{a}_1+y_{10}\mathbf{a}_2+\left(\frac{1}{2}-z_{10}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{x}}+y_{10}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{10}\right)c\,\hat{\mathbf{z}}$	(8*c*)	Ti
$\mathbf{B}_{79}$	=	$x_{10}\mathbf{a}_1+\left(\frac{1}{2}-y_{10}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{10}\right)\mathbf{a}_3$	=	$x_{10}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{10}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{10}\right)c\,\hat{\mathbf{z}}$	(8*c*)	Ti
$\mathbf{B}_{80}$	=	$\left(\frac{1}{2}-x_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{10}\right)\mathbf{a}_2+z_{10}\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{10}\right)b\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}}$	(8*c*)	Ti

References:

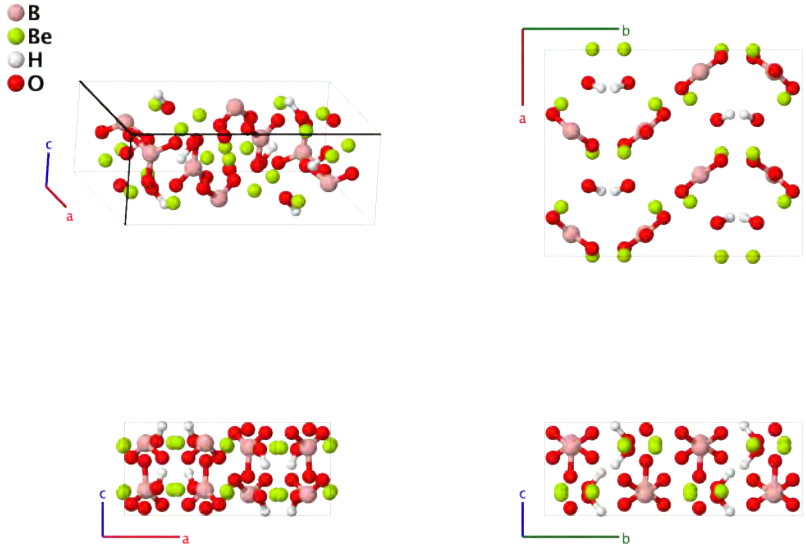
- C.-I. Brändén and I. Lindqvist, *The Crystal Structure of ($TiCl_4$·$POCl_3$)$_2$*, Acta Chem. Scand. **14**, 726–732 (1960), doi:10.3891/acta.chem.scand.14-0726.

Geometry files:

- CIF: pp. 1625
- POSCAR: pp. 1625

Hambergite [$Be_2BO_3(OH)$, $G7_2$] Structure: AB2CD4_oP64_61_c_2c_c_4c

http://aflow.org/prototype-encyclopedia/AB2CD4_oP64_61_c_2c_c_4c

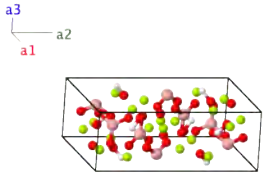


Prototype	:	BBe_2HO_4
AFLOW prototype label	:	AB2CD4_oP64_61_c_2c_c_4c
***Strukturbericht* designation**	:	$G7_2$
Pearson symbol	:	oP64
Space group number	:	61
Space group symbol	:	*Pbca*
AFLOW prototype command	:	`aflow --proto=AB2CD4_oP64_61_c_2c_c_4c --params=`$a, b/a, c/a, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8$

- The original description of this structure in (Hermann, 1937) did not include the positions of the hydrogen atoms, instead using OH for the O(IV)-H radical.
- The sample actually measured by (Gatta, 2012) has 4% of the OH radicals replaced by fluorine.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(8*c*)	B
$\mathbf{B}_2$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right)a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\,\hat{\mathbf{z}}$	(8*c*)	B
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2} + y_1\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right)c\,\hat{\mathbf{z}}$	(8*c*)	B
$\mathbf{B}_4$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_1\right)\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_1\right)b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	(8*c*)	B
$\mathbf{B}_5$	=	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	(8*c*)	B
$\mathbf{B}_6$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + y_1\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right)a\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right)c\,\hat{\mathbf{z}}$	(8*c*)	B
$\mathbf{B}_7$	=	$x_1\,\mathbf{a}_1 + \left(\frac{1}{2} - y_1\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\,\hat{\mathbf{z}}$	(8*c*)	B
$\mathbf{B}_8$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_1\right)\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_1\right)b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(8*c*)	B
$\mathbf{B}_9$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(8*c*)	Be I
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	(8*c*)	Be I
$\mathbf{B}_{11}$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	(8*c*)	Be I
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right)b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(8*c*)	Be I
$\mathbf{B}_{13}$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(8*c*)	Be I
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	(8*c*)	Be I
$\mathbf{B}_{15}$	=	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	(8*c*)	Be I
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right)b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(8*c*)	Be I
$\mathbf{B}_{17}$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(8*c*)	Be II
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(8*c*)	Be II
$\mathbf{B}_{19}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(8*c*)	Be II
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(8*c*)	Be II
$\mathbf{B}_{21}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(8*c*)	Be II
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(8*c*)	Be II
$\mathbf{B}_{23}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(8*c*)	Be II
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(8*c*)	Be II
$\mathbf{B}_{25}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*c*)	H
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(8*c*)	H
$\mathbf{B}_{27}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(8*c*)	H
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*c*)	H
$\mathbf{B}_{29}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*c*)	H
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(8*c*)	H
$\mathbf{B}_{31}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(8*c*)	H
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*c*)	H
$\mathbf{B}_{33}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(8*c*)	O I
$\mathbf{B}_{34}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	(8*c*)	O I
$\mathbf{B}_{35}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	(8*c*)	O I

$$
\begin{aligned}
\mathbf{B}_{36} &= \left(\tfrac{1}{2}+x_5\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_5\right)\mathbf{a}_2-z_5\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_5\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_5\right)b\,\hat{\mathbf{y}}-z_5c\,\hat{\mathbf{z}} && (8c) && \text{O I}\\
\mathbf{B}_{37} &= -x_5\,\mathbf{a}_1-y_5\,\mathbf{a}_2-z_5\,\mathbf{a}_3 &&= -x_5a\,\hat{\mathbf{x}}-y_5b\,\hat{\mathbf{y}}-z_5c\,\hat{\mathbf{z}} && (8c) && \text{O I}\\
\mathbf{B}_{38} &= \left(\tfrac{1}{2}+x_5\right)\mathbf{a}_1+y_5\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_5\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_5\right)a\,\hat{\mathbf{x}}+y_5b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_5\right)c\,\hat{\mathbf{z}} && (8c) && \text{O I}\\
\mathbf{B}_{39} &= x_5\,\mathbf{a}_1+\left(\tfrac{1}{2}-y_5\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_5\right)\mathbf{a}_3 &&= x_5a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_5\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_5\right)c\,\hat{\mathbf{z}} && (8c) && \text{O I}\\
\mathbf{B}_{40} &= \left(\tfrac{1}{2}-x_5\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_5\right)\mathbf{a}_2+z_5\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_5\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_5\right)b\,\hat{\mathbf{y}}+z_5c\,\hat{\mathbf{z}} && (8c) && \text{O I}\\
\mathbf{B}_{41} &= x_6\,\mathbf{a}_1+y_6\,\mathbf{a}_2+z_6\,\mathbf{a}_3 &&= x_6a\,\hat{\mathbf{x}}+y_6b\,\hat{\mathbf{y}}+z_6c\,\hat{\mathbf{z}} && (8c) && \text{O II}\\
\mathbf{B}_{42} &= \left(\tfrac{1}{2}-x_6\right)\mathbf{a}_1-y_6\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{x}}-y_6b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_6\right)c\,\hat{\mathbf{z}} && (8c) && \text{O II}\\
\mathbf{B}_{43} &= -x_6\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_6\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_6\right)\mathbf{a}_3 &&= -x_6a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_6\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_6\right)c\,\hat{\mathbf{z}} && (8c) && \text{O II}\\
\mathbf{B}_{44} &= \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_6\right)\mathbf{a}_2-z_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_6\right)b\,\hat{\mathbf{y}}-z_6c\,\hat{\mathbf{z}} && (8c) && \text{O II}\\
\mathbf{B}_{45} &= -x_6\,\mathbf{a}_1-y_6\,\mathbf{a}_2-z_6\,\mathbf{a}_3 &&= -x_6a\,\hat{\mathbf{x}}-y_6b\,\hat{\mathbf{y}}-z_6c\,\hat{\mathbf{z}} && (8c) && \text{O II}\\
\mathbf{B}_{46} &= \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_1+y_6\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{x}}+y_6b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_6\right)c\,\hat{\mathbf{z}} && (8c) && \text{O II}\\
\mathbf{B}_{47} &= x_6\,\mathbf{a}_1+\left(\tfrac{1}{2}-y_6\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_6\right)\mathbf{a}_3 &&= x_6a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_6\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_6\right)c\,\hat{\mathbf{z}} && (8c) && \text{O II}\\
\mathbf{B}_{48} &= \left(\tfrac{1}{2}-x_6\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_6\right)\mathbf{a}_2+z_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_6\right)b\,\hat{\mathbf{y}}+z_6c\,\hat{\mathbf{z}} && (8c) && \text{O II}\\
\mathbf{B}_{49} &= x_7\,\mathbf{a}_1+y_7\,\mathbf{a}_2+z_7\,\mathbf{a}_3 &&= x_7a\,\hat{\mathbf{x}}+y_7b\,\hat{\mathbf{y}}+z_7c\,\hat{\mathbf{z}} && (8c) && \text{O III}\\
\mathbf{B}_{50} &= \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_1-y_7\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{x}}-y_7b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_7\right)c\,\hat{\mathbf{z}} && (8c) && \text{O III}\\
\mathbf{B}_{51} &= -x_7\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_7\right)\mathbf{a}_3 &&= -x_7a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_7\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_7\right)c\,\hat{\mathbf{z}} && (8c) && \text{O III}\\
\mathbf{B}_{52} &= \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_7\right)\mathbf{a}_2-z_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_7\right)b\,\hat{\mathbf{y}}-z_7c\,\hat{\mathbf{z}} && (8c) && \text{O III}\\
\mathbf{B}_{53} &= -x_7\,\mathbf{a}_1-y_7\,\mathbf{a}_2-z_7\,\mathbf{a}_3 &&= -x_7a\,\hat{\mathbf{x}}-y_7b\,\hat{\mathbf{y}}-z_7c\,\hat{\mathbf{z}} && (8c) && \text{O III}\\
\mathbf{B}_{54} &= \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1+y_7\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{x}}+y_7b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_7\right)c\,\hat{\mathbf{z}} && (8c) && \text{O III}\\
\mathbf{B}_{55} &= x_7\,\mathbf{a}_1+\left(\tfrac{1}{2}-y_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 &&= x_7a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_7\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_7\right)c\,\hat{\mathbf{z}} && (8c) && \text{O III}\\
\mathbf{B}_{56} &= \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_7\right)\mathbf{a}_2+z_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_7\right)b\,\hat{\mathbf{y}}+z_7c\,\hat{\mathbf{z}} && (8c) && \text{O III}\\
\mathbf{B}_{57} &= x_8\,\mathbf{a}_1+y_8\,\mathbf{a}_2+z_8\,\mathbf{a}_3 &&= x_8a\,\hat{\mathbf{x}}+y_8b\,\hat{\mathbf{y}}+z_8c\,\hat{\mathbf{z}} && (8c) && \text{O IV}\\
\mathbf{B}_{58} &= \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_1-y_8\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_8\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{x}}-y_8b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_8\right)c\,\hat{\mathbf{z}} && (8c) && \text{O IV}\\
\mathbf{B}_{59} &= -x_8\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_8\right)\mathbf{a}_3 &&= -x_8a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_8\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_8\right)c\,\hat{\mathbf{z}} && (8c) && \text{O IV}\\
\mathbf{B}_{60} &= \left(\tfrac{1}{2}+x_8\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_8\right)\mathbf{a}_2-z_8\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_8\right)b\,\hat{\mathbf{y}}-z_8c\,\hat{\mathbf{z}} && (8c) && \text{O IV}\\
\mathbf{B}_{61} &= -x_8\,\mathbf{a}_1-y_8\,\mathbf{a}_2-z_8\,\mathbf{a}_3 &&= -x_8a\,\hat{\mathbf{x}}-y_8b\,\hat{\mathbf{y}}-z_8c\,\hat{\mathbf{z}} && (8c) && \text{O IV}\\
\mathbf{B}_{62} &= \left(\tfrac{1}{2}+x_8\right)\mathbf{a}_1+y_8\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_8\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{x}}+y_8b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_8\right)c\,\hat{\mathbf{z}} && (8c) && \text{O IV}\\
\mathbf{B}_{63} &= x_8\,\mathbf{a}_1+\left(\tfrac{1}{2}-y_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_8\right)\mathbf{a}_3 &&= x_8a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_8\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_8\right)c\,\hat{\mathbf{z}} && (8c) && \text{O IV}\\
\mathbf{B}_{64} &= \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_8\right)\mathbf{a}_2+z_8\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_8\right)b\,\hat{\mathbf{y}}+z_8c\,\hat{\mathbf{z}} && (8c) && \text{O IV}
\end{aligned}
$$

References:

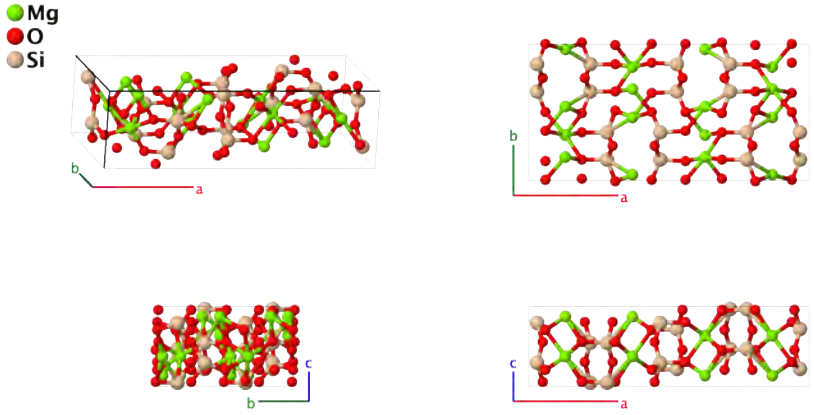
- G. D. Gatta, G. J. McIntyre, G. Bromiley, A. Guastoni, and F. Nestola, *A single-crystal neutron diffraction study of hambergite, $Be_2BO_3(OH,F)$*, Am. Mineral. **97**, 1891–1897 (2012).
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1626
- POSCAR: pp. 1626

Enstatite ($MgSiO_3$, $S4_3$) Structure: AB3C_oP80_61_2c_6c_2c

http://aflow.org/prototype-encyclopedia/AB3C_oP80_61_2c_6c_2c



Prototype	:	MgO_3Si
AFLOW prototype label	:	AB3C_oP80_61_2c_6c_2c
***Strukturbericht* designation**	:	$S4_3$
Pearson symbol	:	oP80
Space group number	:	61
Space group symbol	:	*Pbca*
AFLOW prototype command	:	`aflow --proto=AB3C_oP80_61_2c_6c_2c --params=`$a, b/a, c/a, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}$

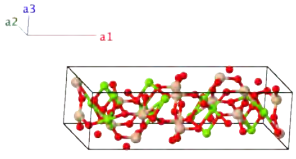
Other compounds with this structure

- $FeSiO_3$ (ferrosilite) and $CoSiO_3$ (Co-pyroxene)

- Enstatite (or properly orthoenstatite) is an end point of orthopyroxene, $XYSi_2O_6$, where X and Y are small-radius divalent cations. If either X or Y is replaced by a transition metal the system becomes a monoclinic clinopyroxene, in space group $C2/c$ #15.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	($8c$)	Mg I

$$
\begin{aligned}
\mathbf{B}_{2} &= \left(\tfrac{1}{2}-x_1\right)\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_1\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_1\right)a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_1\right)c\,\hat{\mathbf{z}} && (8c) && \text{Mg I}\\
\mathbf{B}_{3} &= -x_1\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_1\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_1\right)\mathbf{a}_3 &&= -x_1 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_1\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_1\right)c\,\hat{\mathbf{z}} && (8c) && \text{Mg I}\\
\mathbf{B}_{4} &= \left(\tfrac{1}{2}+x_1\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_1\right)\mathbf{a}_2 - z_1\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_1\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_1\right)b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}} && (8c) && \text{Mg I}\\
\mathbf{B}_{5} &= -x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3 &&= -x_1 a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}} && (8c) && \text{Mg I}\\
\mathbf{B}_{6} &= \left(\tfrac{1}{2}+x_1\right)\mathbf{a}_1 + y_1\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_1\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_1\right)a\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_1\right)c\,\hat{\mathbf{z}} && (8c) && \text{Mg I}\\
\mathbf{B}_{7} &= x_1\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_1\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_1\right)\mathbf{a}_3 &&= x_1 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_1\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_1\right)c\,\hat{\mathbf{z}} && (8c) && \text{Mg I}\\
\mathbf{B}_{8} &= \left(\tfrac{1}{2}-x_1\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_1\right)\mathbf{a}_2 + z_1\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_1\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_1\right)b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}} && (8c) && \text{Mg I}\\
\mathbf{B}_{9} &= x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3 &&= x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}} && (8c) && \text{Mg II}\\
\mathbf{B}_{10} &= \left(\tfrac{1}{2}-x_2\right)\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_2\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_2\right)a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_2\right)c\,\hat{\mathbf{z}} && (8c) && \text{Mg II}\\
\mathbf{B}_{11} &= -x_2\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_2\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_2\right)\mathbf{a}_3 &&= -x_2 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_2\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_2\right)c\,\hat{\mathbf{z}} && (8c) && \text{Mg II}\\
\mathbf{B}_{12} &= \left(\tfrac{1}{2}+x_2\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_2\right)\mathbf{a}_2 - z_2\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_2\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_2\right)b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}} && (8c) && \text{Mg II}\\
\mathbf{B}_{13} &= -x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3 &&= -x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}} && (8c) && \text{Mg II}\\
\mathbf{B}_{14} &= \left(\tfrac{1}{2}+x_2\right)\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_2\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_2\right)a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_2\right)c\,\hat{\mathbf{z}} && (8c) && \text{Mg II}\\
\mathbf{B}_{15} &= x_2\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_2\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_2\right)\mathbf{a}_3 &&= x_2 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_2\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_2\right)c\,\hat{\mathbf{z}} && (8c) && \text{Mg II}\\
\mathbf{B}_{16} &= \left(\tfrac{1}{2}-x_2\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_2\right)\mathbf{a}_2 + z_2\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_2\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_2\right)b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}} && (8c) && \text{Mg II}\\
\mathbf{B}_{17} &= x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3 &&= x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}} && (8c) && \text{O I}\\
\mathbf{B}_{18} &= \left(\tfrac{1}{2}-x_3\right)\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_3\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_3\right)a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_3\right)c\,\hat{\mathbf{z}} && (8c) && \text{O I}\\
\mathbf{B}_{19} &= -x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_3\right)\mathbf{a}_3 &&= -x_3 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_3\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_3\right)c\,\hat{\mathbf{z}} && (8c) && \text{O I}\\
\mathbf{B}_{20} &= \left(\tfrac{1}{2}+x_3\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_3\right)\mathbf{a}_2 - z_3\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_3\right)b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}} && (8c) && \text{O I}\\
\mathbf{B}_{21} &= -x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3 &&= -x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}} && (8c) && \text{O I}\\
\mathbf{B}_{22} &= \left(\tfrac{1}{2}+x_3\right)\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_3\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_3\right)a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_3\right)c\,\hat{\mathbf{z}} && (8c) && \text{O I}\\
\mathbf{B}_{23} &= x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_3\right)\mathbf{a}_3 &&= x_3 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_3\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_3\right)c\,\hat{\mathbf{z}} && (8c) && \text{O I}\\
\mathbf{B}_{24} &= \left(\tfrac{1}{2}-x_3\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_3\right)\mathbf{a}_2 + z_3\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_3\right)b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}} && (8c) && \text{O I}\\
\mathbf{B}_{25} &= x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3 &&= x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}} && (8c) && \text{O II}\\
\mathbf{B}_{26} &= \left(\tfrac{1}{2}-x_4\right)\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_4\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_4\right)a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_4\right)c\,\hat{\mathbf{z}} && (8c) && \text{O II}\\
\mathbf{B}_{27} &= -x_4\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_4\right)\mathbf{a}_3 &&= -x_4 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_4\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_4\right)c\,\hat{\mathbf{z}} && (8c) && \text{O II}\\
\mathbf{B}_{28} &= \left(\tfrac{1}{2}+x_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_4\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_4\right)b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}} && (8c) && \text{O II}\\
\mathbf{B}_{29} &= -x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3 &&= -x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}} && (8c) && \text{O II}\\
\mathbf{B}_{30} &= \left(\tfrac{1}{2}+x_4\right)\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_4\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_4\right)a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_4\right)c\,\hat{\mathbf{z}} && (8c) && \text{O II}\\
\mathbf{B}_{31} &= x_4\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_4\right)\mathbf{a}_3 &&= x_4 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_4\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_4\right)c\,\hat{\mathbf{z}} && (8c) && \text{O II}\\
\mathbf{B}_{32} &= \left(\tfrac{1}{2}-x_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_4\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_4\right)b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}} && (8c) && \text{O II}\\
\mathbf{B}_{33} &= x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3 &&= x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}} && (8c) && \text{O III}\\
\mathbf{B}_{34} &= \left(\tfrac{1}{2}-x_5\right)\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_5\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_5\right)a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_5\right)c\,\hat{\mathbf{z}} && (8c) && \text{O III}\\
\mathbf{B}_{35} &= -x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_5\right)\mathbf{a}_3 &&= -x_5 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_5\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_5\right)c\,\hat{\mathbf{z}} && (8c) && \text{O III}\\
\mathbf{B}_{36} &= \left(\tfrac{1}{2}+x_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_5\right)b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}} && (8c) && \text{O III}\\
\mathbf{B}_{37} &= -x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3 &&= -x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}} && (8c) && \text{O III}
\end{aligned}
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\begin{aligned}
\mathbf{B}_{38} &= \left(\tfrac{1}{2}+x_5\right)\mathbf{a}_1 + y_5\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_5\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_5\right)a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_5\right)c\,\hat{\mathbf{z}} && (8c) && \text{O III}\\
\mathbf{B}_{39} &= x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_5\right)\mathbf{a}_3 &&= x_5 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_5\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_5\right)c\,\hat{\mathbf{z}} && (8c) && \text{O III}\\
\mathbf{B}_{40} &= \left(\tfrac{1}{2}-x_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_5\right)\mathbf{a}_2 + z_5\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_5\right)b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}} && (8c) && \text{O III}\\
\mathbf{B}_{41} &= x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 &&= x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}} && (8c) && \text{O IV}\\
\mathbf{B}_{42} &= \left(\tfrac{1}{2}-x_6\right)\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_6\right)c\,\hat{\mathbf{z}} && (8c) && \text{O IV}\\
\mathbf{B}_{43} &= -x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_6\right)\mathbf{a}_3 &&= -x_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_6\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_6\right)c\,\hat{\mathbf{z}} && (8c) && \text{O IV}\\
\mathbf{B}_{44} &= \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_6\right)b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}} && (8c) && \text{O IV}\\
\mathbf{B}_{45} &= -x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3 &&= -x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}} && (8c) && \text{O IV}\\
\mathbf{B}_{46} &= \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_1 + y_6\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_6\right)c\,\hat{\mathbf{z}} && (8c) && \text{O IV}\\
\mathbf{B}_{47} &= x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_6\right)\mathbf{a}_3 &&= x_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_6\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_6\right)c\,\hat{\mathbf{z}} && (8c) && \text{O IV}\\
\mathbf{B}_{48} &= \left(\tfrac{1}{2}-x_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_6\right)\mathbf{a}_2 + z_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_6\right)b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}} && (8c) && \text{O IV}\\
\mathbf{B}_{49} &= x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 &&= x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}} && (8c) && \text{O V}\\
\mathbf{B}_{50} &= \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_7\right)c\,\hat{\mathbf{z}} && (8c) && \text{O V}\\
\mathbf{B}_{51} &= -x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_3 &&= -x_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_7\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_7\right)c\,\hat{\mathbf{z}} && (8c) && \text{O V}\\
\mathbf{B}_{52} &= \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_7\right)\mathbf{a}_2 - z_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_7\right)b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}} && (8c) && \text{O V}\\
\mathbf{B}_{53} &= -x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3 &&= -x_7 a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}} && (8c) && \text{O V}\\
\mathbf{B}_{54} &= \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1 + y_7\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_7\right)c\,\hat{\mathbf{z}} && (8c) && \text{O V}\\
\mathbf{B}_{55} &= x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 &&= x_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_7\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_7\right)c\,\hat{\mathbf{z}} && (8c) && \text{O V}\\
\mathbf{B}_{56} &= \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_7\right)\mathbf{a}_2 + z_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_7\right)b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}} && (8c) && \text{O V}\\
\mathbf{B}_{57} &= x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 &&= x_8 a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}} && (8c) && \text{O VI}\\
\mathbf{B}_{58} &= \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_1 - y_8\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_8\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_8\right)c\,\hat{\mathbf{z}} && (8c) && \text{O VI}\\
\mathbf{B}_{59} &= -x_8\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_8\right)\mathbf{a}_3 &&= -x_8 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_8\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_8\right)c\,\hat{\mathbf{z}} && (8c) && \text{O VI}\\
\mathbf{B}_{60} &= \left(\tfrac{1}{2}+x_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_8\right)\mathbf{a}_2 - z_8\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_8\right)b\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}} && (8c) && \text{O VI}\\
\mathbf{B}_{61} &= -x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3 &&= -x_8 a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}} && (8c) && \text{O VI}\\
\mathbf{B}_{62} &= \left(\tfrac{1}{2}+x_8\right)\mathbf{a}_1 + y_8\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_8\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_8\right)c\,\hat{\mathbf{z}} && (8c) && \text{O VI}\\
\mathbf{B}_{63} &= x_8\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_8\right)\mathbf{a}_3 &&= x_8 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_8\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_8\right)c\,\hat{\mathbf{z}} && (8c) && \text{O VI}\\
\mathbf{B}_{64} &= \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_8\right)\mathbf{a}_2 + z_8\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_8\right)b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}} && (8c) && \text{O VI}\\
\mathbf{B}_{65} &= x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 &&= x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}} && (8c) && \text{Si I}\\
\mathbf{B}_{66} &= \left(\tfrac{1}{2}-x_9\right)\mathbf{a}_1 - y_9\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_9\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_9\right)a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_9\right)c\,\hat{\mathbf{z}} && (8c) && \text{Si I}\\
\mathbf{B}_{67} &= -x_9\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_9\right)\mathbf{a}_3 &&= -x_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_9\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_9\right)c\,\hat{\mathbf{z}} && (8c) && \text{Si I}\\
\mathbf{B}_{68} &= \left(\tfrac{1}{2}+x_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_9\right)\mathbf{a}_2 - z_9\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_9\right)b\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}} && (8c) && \text{Si I}\\
\mathbf{B}_{69} &= -x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3 &&= -x_9 a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}} && (8c) && \text{Si I}\\
\mathbf{B}_{70} &= \left(\tfrac{1}{2}+x_9\right)\mathbf{a}_1 + y_9\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_9\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_9\right)a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_9\right)c\,\hat{\mathbf{z}} && (8c) && \text{Si I}\\
\mathbf{B}_{71} &= x_9\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_9\right)\mathbf{a}_3 &&= x_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_9\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_9\right)c\,\hat{\mathbf{z}} && (8c) && \text{Si I}\\
\mathbf{B}_{72} &= \left(\tfrac{1}{2}-x_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_9\right)\mathbf{a}_2 + z_9\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_9\right)b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}} && (8c) && \text{Si I}\\
\mathbf{B}_{73} &= x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 &&= x_{10} a\,\hat{\mathbf{x}} + y_{10} b\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}} && (8c) && \text{Si II}
\end{aligned}
$$

$\mathbf{B}_{74}$	=	$\left(\frac{1}{2}-x_{10}\right)\mathbf{a}_1-y_{10}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{10}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{x}}-y_{10}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{10}\right)c\,\hat{\mathbf{z}}$	(8*c*)	Si II
$\mathbf{B}_{75}$	=	$-x_{10}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{10}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{10}\right)\mathbf{a}_3$	=	$-x_{10}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{10}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{10}\right)c\,\hat{\mathbf{z}}$	(8*c*)	Si II
$\mathbf{B}_{76}$	=	$\left(\frac{1}{2}+x_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{10}\right)\mathbf{a}_2-z_{10}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{10}\right)b\,\hat{\mathbf{y}}-z_{10}c\,\hat{\mathbf{z}}$	(8*c*)	Si II
$\mathbf{B}_{77}$	=	$-x_{10}\,\mathbf{a}_1-y_{10}\,\mathbf{a}_2-z_{10}\,\mathbf{a}_3$	=	$-x_{10}a\,\hat{\mathbf{x}}-y_{10}b\,\hat{\mathbf{y}}-z_{10}c\,\hat{\mathbf{z}}$	(8*c*)	Si II
$\mathbf{B}_{78}$	=	$\left(\frac{1}{2}+x_{10}\right)\mathbf{a}_1+y_{10}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{10}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{x}}+y_{10}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{10}\right)c\,\hat{\mathbf{z}}$	(8*c*)	Si II
$\mathbf{B}_{79}$	=	$x_{10}\,\mathbf{a}_1+\left(\frac{1}{2}-y_{10}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{10}\right)\mathbf{a}_3$	=	$x_{10}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{10}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{10}\right)c\,\hat{\mathbf{z}}$	(8*c*)	Si II
$\mathbf{B}_{80}$	=	$\left(\frac{1}{2}-x_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{10}\right)\mathbf{a}_2+z_{10}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{10}\right)b\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}}$	(8*c*)	Si II

References:

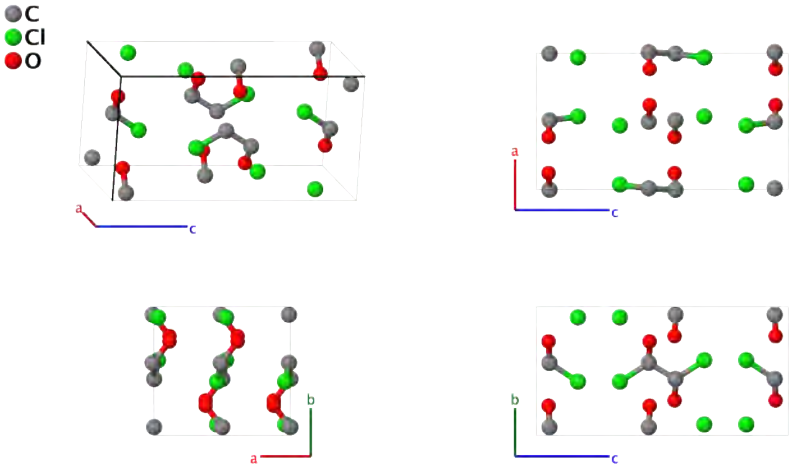
- B. E. Warren and D. I. Modell, *The Structure of Enstatite $MgSiO_3$*, Zeitschrift für Kristallographie - Crystalline Materials **75**, 1–14 (1930), doi:10.1515/zkri-1930-0102.

Geometry files:

- CIF: pp. 1627
- POSCAR: pp. 1627

COCl Structure: ABC_oP24_61_c_c_c

http://aflow.org/prototype-encyclopedia/ABC_oP24_61_c_c_c

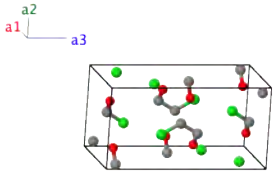


Prototype	:	CClO
AFLOW prototype label	:	ABC_oP24_61_c_c_c
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP24
Space group number	:	61
Space group symbol	:	*Pbca*
AFLOW prototype command	:	`aflow --proto=ABC_oP24_61_c_c_c --params=`$a, b/a, c/a, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3$

- (Groth, 1962) are not confident about the atomic positions, saying that "The computed values for the oxalyl chloride molecule . . . [are probably] less reliable than those listed for the bromide molecule [COBr]." In fact, we have not been able to reconcile their crystal structure for COBr with the distances between the atoms, so we have left that structure out of the encyclopedia at this time.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$ =	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$ =	$x_1 a\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(8*c*)	C
$\mathbf{B}_2$ =	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$ =	$\left(\frac{1}{2} - x_1\right)a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\,\hat{\mathbf{z}}$	(8*c*)	C
$\mathbf{B}_3$ =	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2} + y_1\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$ =	$-x_1 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_1\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right)c\,\hat{\mathbf{z}}$	(8*c*)	C
$\mathbf{B}_4$ =	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_1\right)\mathbf{a}_2 - z_1\,\mathbf{a}_3$ =	$\left(\frac{1}{2} + x_1\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_1\right)b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	(8*c*)	C

$$
\begin{array}{lcccccc}
\mathbf{B}_5 & = & -x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3 & = & -x_1 a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}} & (8c) & \mathrm{C} \\
\mathbf{B}_6 & = & \left(\tfrac{1}{2}+x_1\right)\mathbf{a}_1 + y_1\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_1\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}+x_1\right)a\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_1\right)c\,\hat{\mathbf{z}} & (8c) & \mathrm{C} \\
\mathbf{B}_7 & = & x_1\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_1\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_1\right)\mathbf{a}_3 & = & x_1 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_1\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_1\right)c\,\hat{\mathbf{z}} & (8c) & \mathrm{C} \\
\mathbf{B}_8 & = & \left(\tfrac{1}{2}-x_1\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_1\right)\mathbf{a}_2 + z_1\,\mathbf{a}_3 & = & \left(\tfrac{1}{2}-x_1\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_1\right)b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}} & (8c) & \mathrm{C} \\
\mathbf{B}_9 & = & x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3 & = & x_2 a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}} & (8c) & \mathrm{Cl} \\
\mathbf{B}_{10} & = & \left(\tfrac{1}{2}-x_2\right)\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_2\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}-x_2\right)a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_2\right)c\,\hat{\mathbf{z}} & (8c) & \mathrm{Cl} \\
\mathbf{B}_{11} & = & -x_2\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_2\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_2\right)\mathbf{a}_3 & = & -x_2 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_2\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_2\right)c\,\hat{\mathbf{z}} & (8c) & \mathrm{Cl} \\
\mathbf{B}_{12} & = & \left(\tfrac{1}{2}+x_2\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_2\right)\mathbf{a}_2 - z_2\,\mathbf{a}_3 & = & \left(\tfrac{1}{2}+x_2\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_2\right)b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}} & (8c) & \mathrm{Cl} \\
\mathbf{B}_{13} & = & -x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3 & = & -x_2 a\,\hat{\mathbf{x}} - y_2 b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}} & (8c) & \mathrm{Cl} \\
\mathbf{B}_{14} & = & \left(\tfrac{1}{2}+x_2\right)\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_2\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}+x_2\right)a\,\hat{\mathbf{x}} + y_2 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_2\right)c\,\hat{\mathbf{z}} & (8c) & \mathrm{Cl} \\
\mathbf{B}_{15} & = & x_2\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_2\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_2\right)\mathbf{a}_3 & = & x_2 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_2\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_2\right)c\,\hat{\mathbf{z}} & (8c) & \mathrm{Cl} \\
\mathbf{B}_{16} & = & \left(\tfrac{1}{2}-x_2\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_2\right)\mathbf{a}_2 + z_2\,\mathbf{a}_3 & = & \left(\tfrac{1}{2}-x_2\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_2\right)b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}} & (8c) & \mathrm{Cl} \\
\mathbf{B}_{17} & = & x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3 & = & x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}} & (8c) & \mathrm{O} \\
\mathbf{B}_{18} & = & \left(\tfrac{1}{2}-x_3\right)\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_3\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}-x_3\right)a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_3\right)c\,\hat{\mathbf{z}} & (8c) & \mathrm{O} \\
\mathbf{B}_{19} & = & -x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_3\right)\mathbf{a}_3 & = & -x_3 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_3\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_3\right)c\,\hat{\mathbf{z}} & (8c) & \mathrm{O} \\
\mathbf{B}_{20} & = & \left(\tfrac{1}{2}+x_3\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_3\right)\mathbf{a}_2 - z_3\,\mathbf{a}_3 & = & \left(\tfrac{1}{2}+x_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_3\right)b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}} & (8c) & \mathrm{O} \\
\mathbf{B}_{21} & = & -x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3 & = & -x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}} & (8c) & \mathrm{O} \\
\mathbf{B}_{22} & = & \left(\tfrac{1}{2}+x_3\right)\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_3\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2}+x_3\right)a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_3\right)c\,\hat{\mathbf{z}} & (8c) & \mathrm{O} \\
\mathbf{B}_{23} & = & x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_3\right)\mathbf{a}_3 & = & x_3 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_3\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_3\right)c\,\hat{\mathbf{z}} & (8c) & \mathrm{O} \\
\mathbf{B}_{24} & = & \left(\tfrac{1}{2}-x_3\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_3\right)\mathbf{a}_2 + z_3\,\mathbf{a}_3 & = & \left(\tfrac{1}{2}-x_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_3\right)b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}} & (8c) & \mathrm{O}
\end{array}
$$

References:

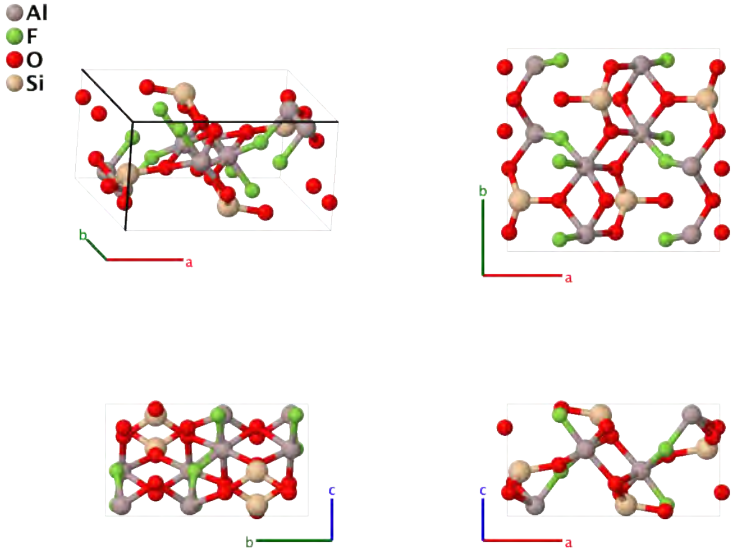
- P. Groth and O. Hassel, *Crystal Structures of Oxalyl Bromide and Oxalyl Chloride*, Acta Chem. Scand. **16**, 2311–2317 (1962), doi:10.3891/acta.chem.scand.16-2311.

Geometry files:

- CIF: pp. 1627
- POSCAR: pp. 1628

Topaz ($Al_2SiO_4F_2$, $S0_5$) Structure: A2B2C4D_oP36_62_d_d_2cd_c

http://aflow.org/prototype-encyclopedia/A2B2C4D_oP36_62_d_d_2cd_c



Prototype	:	$Al_2F_2O_4Si$
AFLOW prototype label	:	A2B2C4D_oP36_62_d_d_2cd_c
***Strukturbericht* designation**	:	$S0_5$
Pearson symbol	:	oP36
Space group number	:	62
Space group symbol	:	*Pnma*
AFLOW prototype command	:	`aflow --proto=A2B2C4D_oP36_62_d_d_2cd_c --params=`$a, b/a, c/a, x_1, z_1, x_2, z_2, x_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6$

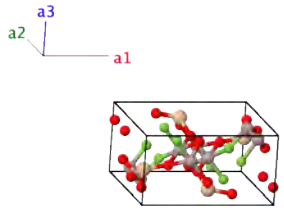
Other compounds with this structure

- $Al_2SiO_4(F,OH)_2$

- The fluorine (8*d*) site in topaz can have a composition $F_{1-x}(OH)_x$, where $x < 0.3$.
- We use the data at $T = 298$ K from the "Pax08" sample in (Komatsu, 2003).
- The data was presented in the *Pbnm* setting of space group #62. We used FINDSYM to transform this into the standard *Pnma* setting.
- (Herman, 1937) gave this the $S0_5$ designation, but co-listed it as $H5_5$ in the index.

Simple Orthorhombic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= a\,\hat{\mathbf{x}}\\ \mathbf{a}_2 &= b\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	$x_1 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_2$	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_3$	$-x_1\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	$-x_1 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_4$	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_1\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_5$	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	$x_2 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_6$	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_7$	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	$-x_2 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_8$	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right) c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_9$	$x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$x_3 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*c*)	Si
$\mathbf{B}_{10}$	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_3\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	(4*c*)	Si
$\mathbf{B}_{11}$	$-x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	$-x_3 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(4*c*)	Si
$\mathbf{B}_{12}$	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_3\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right) c\,\hat{\mathbf{z}}$	(4*c*)	Si
$\mathbf{B}_{13}$	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*d*)	Al
$\mathbf{B}_{14}$	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	(8*d*)	Al
$\mathbf{B}_{15}$	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3$	$-x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right) b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*d*)	Al
$\mathbf{B}_{16}$	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\,\hat{\mathbf{z}}$	(8*d*)	Al
$\mathbf{B}_{17}$	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*d*)	Al
$\mathbf{B}_{18}$	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\,\hat{\mathbf{z}}$	(8*d*)	Al
$\mathbf{B}_{19}$	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right) b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*d*)	Al
$\mathbf{B}_{20}$	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	(8*d*)	Al
$\mathbf{B}_{21}$	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(8*d*)	F
$\mathbf{B}_{22}$	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_5\right) a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	(8*d*)	F
$\mathbf{B}_{23}$	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3$	$-x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right) b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(8*d*)	F
$\mathbf{B}_{24}$	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right) c\,\hat{\mathbf{z}}$	(8*d*)	F
$\mathbf{B}_{25}$	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(8*d*)	F

$\mathbf{B}_{26}$	=	$\left(\frac{1}{2}+x_5\right)\mathbf{a}_1+y_5\,\mathbf{a}_2+\left(\frac{1}{2}-z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{x}}+y_5 b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_5\right)c\,\hat{\mathbf{z}}$	(8*d*)	F
$\mathbf{B}_{27}$	=	$x_5\,\mathbf{a}_1+\left(\frac{1}{2}-y_5\right)\mathbf{a}_2+z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_5\right)b\,\hat{\mathbf{y}}+z_5 c\,\hat{\mathbf{z}}$	(8*d*)	F
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2}-x_5\right)\mathbf{a}_1+\left(\frac{1}{2}+y_5\right)\mathbf{a}_2+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_5\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	(8*d*)	F
$\mathbf{B}_{29}$	=	$x_6\,\mathbf{a}_1+y_6\,\mathbf{a}_2+z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}}+y_6 b\,\hat{\mathbf{y}}+z_6 c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2}-x_6\right)\mathbf{a}_1-y_6\,\mathbf{a}_2+\left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{x}}-y_6 b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{31}$	=	$-x_6\,\mathbf{a}_1+\left(\frac{1}{2}+y_6\right)\mathbf{a}_2-z_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_6\right)b\,\hat{\mathbf{y}}-z_6 c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2}+x_6\right)\mathbf{a}_1+\left(\frac{1}{2}-y_6\right)\mathbf{a}_2+\left(\frac{1}{2}-z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_6\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_6\right)c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{33}$	=	$-x_6\,\mathbf{a}_1-y_6\,\mathbf{a}_2-z_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}}-y_6 b\,\hat{\mathbf{y}}-z_6 c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{34}$	=	$\left(\frac{1}{2}+x_6\right)\mathbf{a}_1+y_6\,\mathbf{a}_2+\left(\frac{1}{2}-z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{x}}+y_6 b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_6\right)c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{35}$	=	$x_6\,\mathbf{a}_1+\left(\frac{1}{2}-y_6\right)\mathbf{a}_2+z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_6\right)b\,\hat{\mathbf{y}}+z_6 c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{36}$	=	$\left(\frac{1}{2}-x_6\right)\mathbf{a}_1+\left(\frac{1}{2}+y_6\right)\mathbf{a}_2+\left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_6\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	(8*d*)	O III

References:

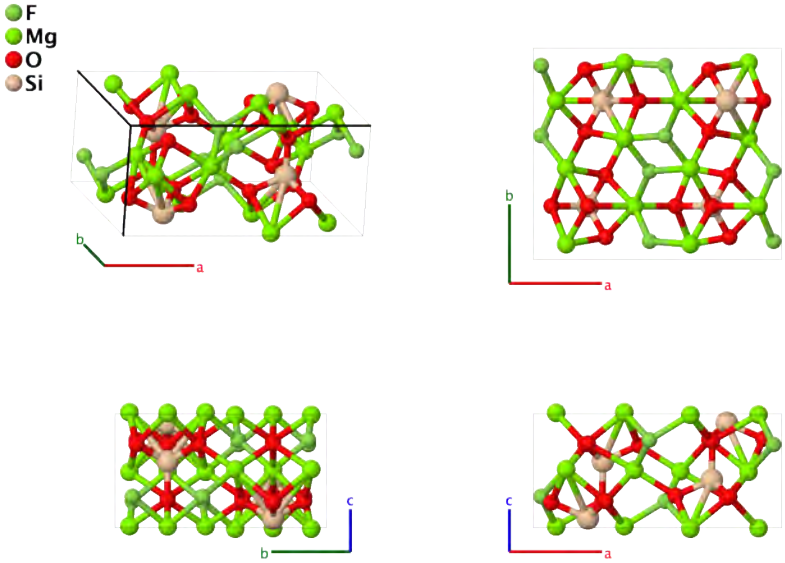
- K. Komatsu, T. Kuribayashi, and Y. Kudoh, *Effect of temperature and pressure on the crystal structure of topaz, $Al_2SiO_4(OH,F)_2$*, J. Miner. Petrol. Sci. **98**, 167–180 (2003), doi:10.2465/jmps.98.167.
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1628
- POSCAR: pp. 1628

Norbergite [$Mg(F,OH)_2 \cdot Mg_2SiO_4$, $S0_7$] Structure: A2B3C4D_oP40_62_d_cd_2cd_c

http://aflow.org/prototype-encyclopedia/A2B3C4D_oP40_62_d_cd_2cd_c

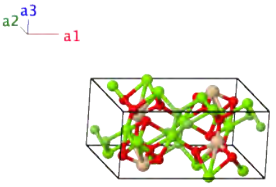


Prototype	:	$F_2Mg_3O_4Si$
AFLOW prototype label	:	A2B3C4D_oP40_62_d_cd_2cd_c
***Strukturbericht* designation**	:	$S0_7$
Pearson symbol	:	oP40
Space group number	:	62
Space group symbol	:	*Pnma*
AFLOW prototype command	:	`aflow --proto=A2B3C4D_oP40_62_d_cd_2cd_c --params=`$a, b/a, c/a, x_1, z_1, x_2, z_2, x_3, z_3, x_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7$

- (Hermann, 1937) define *Strukturbericht* designation $S0_7$ as the general class of humite materials, $Mg(F,OH)_2 \cdot nMg_2SiO_4$. Their major reference for this is (Taylor, 1929), who gave structural data for the $n = 1$ structure, norbergite.
- We use the data from the fluorine-rich sample studied by (Gibbs, 1969), and assume that the (F, OH) site is pure fluorine. (Camara, 1997) studied OH-rich samples, and were able to locate the hydrogen atoms.
- (Gibbs, 1969) give the structure in the *Pbnm* setting of space group #62. We used FINDSYM to change this to the standard *Pnma* setting. This involves a rotation of the axis so that $\hat{x} \to \hat{z} \to \hat{y} \to \hat{x}$.

Simple Orthorhombic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= a\,\hat{\mathbf{x}}\\ \mathbf{a}_2 &= b\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	$x_1 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	$(4c)$	Mg I
$\mathbf{B}_2$	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	$(4c)$	Mg I
$\mathbf{B}_3$	$-x_1\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	$-x_1 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	$(4c)$	Mg I
$\mathbf{B}_4$	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_1\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right) c\,\hat{\mathbf{z}}$	$(4c)$	Mg I
$\mathbf{B}_5$	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	$x_2 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(4c)$	O I
$\mathbf{B}_6$	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	$(4c)$	O I
$\mathbf{B}_7$	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	$-x_2 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	$(4c)$	O I
$\mathbf{B}_8$	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right) c\,\hat{\mathbf{z}}$	$(4c)$	O I
$\mathbf{B}_9$	$x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$x_3 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(4c)$	O II
$\mathbf{B}_{10}$	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_3\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	$(4c)$	O II
$\mathbf{B}_{11}$	$-x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	$-x_3 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(4c)$	O II
$\mathbf{B}_{12}$	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_3\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right) c\,\hat{\mathbf{z}}$	$(4c)$	O II
$\mathbf{B}_{13}$	$x_4\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$x_4 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(4c)$	Si
$\mathbf{B}_{14}$	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	$(4c)$	Si
$\mathbf{B}_{15}$	$-x_4\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	$-x_4 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	$(4c)$	Si
$\mathbf{B}_{16}$	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\,\hat{\mathbf{z}}$	$(4c)$	Si
$\mathbf{B}_{17}$	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(8d)$	F
$\mathbf{B}_{18}$	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_5\right) a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	$(8d)$	F
$\mathbf{B}_{19}$	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3$	$-x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right) b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	$(8d)$	F
$\mathbf{B}_{20}$	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right) c\,\hat{\mathbf{z}}$	$(8d)$	F
$\mathbf{B}_{21}$	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	$(8d)$	F
$\mathbf{B}_{22}$	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + y_5\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right) c\,\hat{\mathbf{z}}$	$(8d)$	F
$\mathbf{B}_{23}$	$x_5\,\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + z_5\,\mathbf{a}_3$	$x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right) b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(8d)$	F
$\mathbf{B}_{24}$	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_5\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	$(8d)$	F
$\mathbf{B}_{25}$	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(8d)$	Mg II
$\mathbf{B}_{26}$	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	$(8d)$	Mg II

$\mathbf{B}_{27}$	=	$-x_6\,\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right) b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(8*d*)	Mg II
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right) c\,\hat{\mathbf{z}}$	(8*d*)	Mg II
$\mathbf{B}_{29}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(8*d*)	Mg II
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + y_6\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right) c\,\hat{\mathbf{z}}$	(8*d*)	Mg II
$\mathbf{B}_{31}$	=	$x_6\,\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right) b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(8*d*)	Mg II
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	(8*d*)	Mg II
$\mathbf{B}_{33}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{34}$	=	$\left(\frac{1}{2} - x_7\right)\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_7\right) a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right) c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{35}$	=	$-x_7\,\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right) b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{36}$	=	$\left(\frac{1}{2} + x_7\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_7\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_7\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_7\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right) c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{37}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{38}$	=	$\left(\frac{1}{2} + x_7\right)\mathbf{a}_1 + y_7\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_7\right) a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right) c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{39}$	=	$x_7\,\mathbf{a}_1 + \left(\frac{1}{2} - y_7\right)\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_7\right) b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{40}$	=	$\left(\frac{1}{2} - x_7\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_7\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right) c\,\hat{\mathbf{z}}$	(8*d*)	O III

References:

- G. V. Gibbs and P. H. Ribbe, *The crystal structures of the humite minerals: I. Norbergite*, Am. Mineral. **54**, 376–390 (1969).
- F. Cámara, *New data on the structure of norbergite; location of hydrogen by X-ray diffraction*, Can. Mineral. **35**, 1523–1530 (1997).
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).
- W. H. Taylor and J. West, *The structure of norbergite*, Zeitschrift für Kristallographie - Crystalline Materials **70**, 461–474 (1929), doi:10.1524/zkri.1929.70.1.461.

Found in:

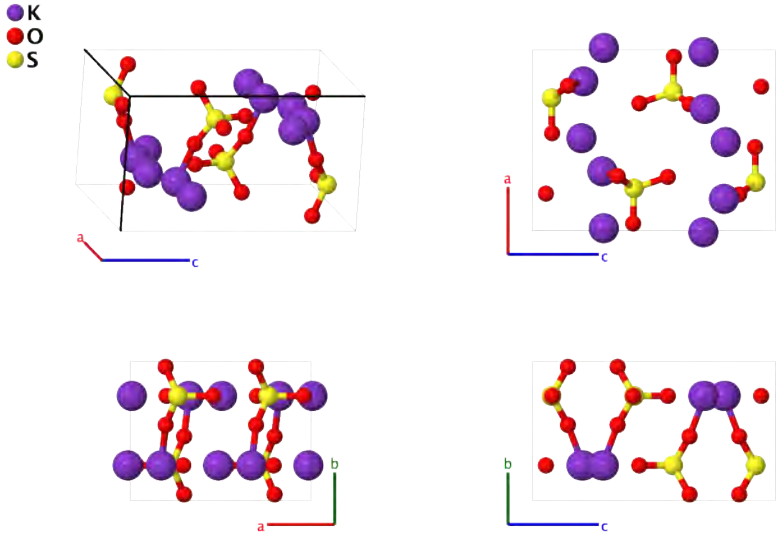
- B. Downs and H. Yang, *The RRUFF^{TM} Project*, http://rruff.info/Norbergite. Norbergite, RRUFF ID: R050280.

Geometry files:

- CIF: pp. 1629
- POSCAR: pp. 1629

Arcanite (K_2SO_4, $H1_6$) Structure: A2B4C_oP28_62_2c_2cd_c

http://aflow.org/prototype-encyclopedia/A2B4C_oP28_62_2c_2cd_c



Prototype	:	K_2O_4S
AFLOW prototype label	:	A2B4C_oP28_62_2c_2cd_c
***Strukturbericht* designation**	:	$H1_6$
Pearson symbol	:	oP28
Space group number	:	62
Space group symbol	:	*Pnma*
AFLOW prototype command	:	`aflow --proto=A2B4C_oP28_62_2c_2cd_c --params=`$a, b/a, c/a, x_1, z_1, x_2, z_2, x_3, z_3, x_4, z_4, x_5, z_5, x_6, y_6, z_6$

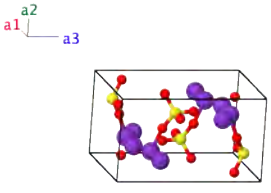
Other compounds with this structure

- K_2CrO_4, K_2MnO_4, K_2SeO_4, and Cs_2CuCl_4

- (McGinnety, 1972) gives the atomic positions using the *Pnam* setting of space group #62. We have switched the y- and z-axes to put the crystal in the *Pnma* setting.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(4*c*)	K I
$\mathbf{B}_2$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	K I
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	(4*c*)	K I
$\mathbf{B}_4$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	K I
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*c*)	K II
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	(4*c*)	K II
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(4*c*)	K II
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right) c\,\hat{\mathbf{z}}$	(4*c*)	K II
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right) c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(4*c*)	S
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	(4*c*)	S
$\mathbf{B}_{19}$	=	$-x_5\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(4*c*)	S
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right) c\,\hat{\mathbf{z}}$	(4*c*)	S
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{23}$	=	$-x_6\,\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right) b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right) c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{25}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + y_6\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right) c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{27}$	=	$x_6\,\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right) b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	(8*d*)	O III

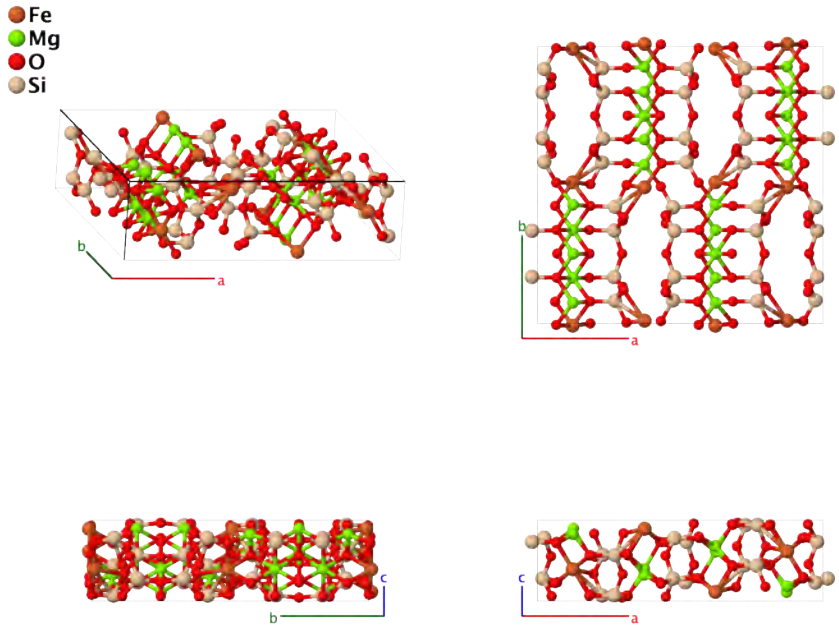
References:
- J. A. McGinnety, *Redetermination of the structures of potassium sulphate and potassium chromate: the effect of electrostatic crystal forces upon observed bond lengths*, Acta Crystallogr. Sect. B Struct. Sci. **28**, 2845–2852 (1972), doi:10.1107/S0567740872007022.

Found in:
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:
- CIF: pp. 1629
- POSCAR: pp. 1630

Anthophyllite ($Mg_5Fe_2Si_8O_{22}(OH)_2$, $S4_4$) Structure: A2B5C22D2E8_oP156_62_d_c2d_2c10d_2c_4d

http://aflow.org/prototype-encyclopedia/A2B5C22D2E8_oP156_62_d_c2d_2c10d_2c_4d



Prototype	:	$Fe_2Mg_5O_{22}(OH)_2Si_8$
AFLOW prototype label	:	A2B5C22D2E8_oP156_62_d_c2d_2c10d_2c_4d
***Strukturbericht* designation**	:	$S4_4$
Pearson symbol	:	oP156
Space group number	:	62
Space group symbol	:	*Pnma*
AFLOW prototype command	:	`aflow --proto=A2B5C22D2E8_oP156_62_d_c2d_2c10d_2c_4d --params=`$a, b/a, c/a, x_1, z_1, x_2, z_2, x_3, z_3, x_4, z_4, x_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}, x_{17}, y_{17}, z_{17}, x_{18}, y_{18}, z_{18}, x_{19}, y_{19}, z_{19}, x_{20}, y_{20}, z_{20}, x_{21}, y_{21}, z_{21}, x_{22}, y_{22}, z_{22}$

Other compounds with this structure

- $H_2Mg_7Si_8O_{24}$

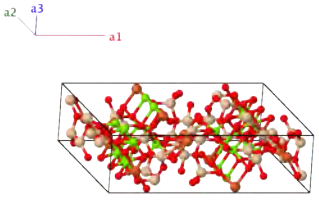
- (Warren, 1930) analyzed their sample of anthophyllite under the assumption that it was free of iron, and this structure was given the $S4_4$ designation by (Hermann, 1937).
- (Walitzi, 1989) found a more accurate determination of the structure, but their sample included substantial amounts of iron. The Mg-I (4*c*) site is actually ($Mg_{0.99}Fe_{0.01}$), Mg-II (8*d*) is ($Mg_{0.98}Fe_{0.02}$), and Fe (8*d*) is ($Mg_{0.38}Fe_{0.62}$). Trace amounts of calcium, manganese and sodium were found on the magnesium/iron sites, and trace amounts of aluminum

were found on the silicon sites. Neither paper was able to determine the positions of the hydrogen atoms, which are included in the OH radicals.

- Anthophyllite can be thought of as an approximation to a doubled unit cell of protoanthophyllite, where the hydrogen atoms have been located.

Simple Orthorhombic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= a\,\hat{\mathbf{x}}\\ \mathbf{a}_2 &= b\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(4*c*)	Mg I
$\mathbf{B}_2$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	Mg I
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	(4*c*)	Mg I
$\mathbf{B}_4$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	Mg I
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right) c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right) c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(4*c*)	OH I
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	(4*c*)	OH I
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(4*c*)	OH I
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\,\hat{\mathbf{z}}$	(4*c*)	OH I
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(4*c*)	OH II
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	(4*c*)	OH II
$\mathbf{B}_{19}$	=	$-x_5\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(4*c*)	OH II
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right) c\,\hat{\mathbf{z}}$	(4*c*)	OH II
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(8*d*)	Fe
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	(8*d*)	Fe
$\mathbf{B}_{23}$	=	$-x_6\,\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right) b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(8*d*)	Fe

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\begin{aligned}
\mathbf{B}_{24} &= \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_6\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_6\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_6\right)c\,\hat{\mathbf{z}} && (8d) && \text{Fe}\\
\mathbf{B}_{25} &= -x_6\,\mathbf{a}_1-y_6\,\mathbf{a}_2-z_6\,\mathbf{a}_3 &&= -x_6a\,\hat{\mathbf{x}}-y_6b\,\hat{\mathbf{y}}-z_6c\,\hat{\mathbf{z}} && (8d) && \text{Fe}\\
\mathbf{B}_{26} &= \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_1+y_6\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{x}}+y_6b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_6\right)c\,\hat{\mathbf{z}} && (8d) && \text{Fe}\\
\mathbf{B}_{27} &= x_6\,\mathbf{a}_1+\left(\tfrac{1}{2}-y_6\right)\mathbf{a}_2+z_6\,\mathbf{a}_3 &&= x_6a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_6\right)b\,\hat{\mathbf{y}}+z_6c\,\hat{\mathbf{z}} && (8d) && \text{Fe}\\
\mathbf{B}_{28} &= \left(\tfrac{1}{2}-x_6\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_6\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_6\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_6\right)c\,\hat{\mathbf{z}} && (8d) && \text{Fe}\\
\mathbf{B}_{29} &= x_7\,\mathbf{a}_1+y_7\,\mathbf{a}_2+z_7\,\mathbf{a}_3 &&= x_7a\,\hat{\mathbf{x}}+y_7b\,\hat{\mathbf{y}}+z_7c\,\hat{\mathbf{z}} && (8d) && \text{Mg II}\\
\mathbf{B}_{30} &= \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_1-y_7\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{x}}-y_7b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_7\right)c\,\hat{\mathbf{z}} && (8d) && \text{Mg II}\\
\mathbf{B}_{31} &= -x_7\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_7\right)\mathbf{a}_2-z_7\,\mathbf{a}_3 &&= -x_7a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_7\right)b\,\hat{\mathbf{y}}-z_7c\,\hat{\mathbf{z}} && (8d) && \text{Mg II}\\
\mathbf{B}_{32} &= \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_7\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_7\right)c\,\hat{\mathbf{z}} && (8d) && \text{Mg II}\\
\mathbf{B}_{33} &= -x_7\,\mathbf{a}_1-y_7\,\mathbf{a}_2-z_7\,\mathbf{a}_3 &&= -x_7a\,\hat{\mathbf{x}}-y_7b\,\hat{\mathbf{y}}-z_7c\,\hat{\mathbf{z}} && (8d) && \text{Mg II}\\
\mathbf{B}_{34} &= \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1+y_7\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{x}}+y_7b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_7\right)c\,\hat{\mathbf{z}} && (8d) && \text{Mg II}\\
\mathbf{B}_{35} &= x_7\,\mathbf{a}_1+\left(\tfrac{1}{2}-y_7\right)\mathbf{a}_2+z_7\,\mathbf{a}_3 &&= x_7a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_7\right)b\,\hat{\mathbf{y}}+z_7c\,\hat{\mathbf{z}} && (8d) && \text{Mg II}\\
\mathbf{B}_{36} &= \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_7\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_7\right)c\,\hat{\mathbf{z}} && (8d) && \text{Mg II}\\
\mathbf{B}_{37} &= x_8\,\mathbf{a}_1+y_8\,\mathbf{a}_2+z_8\,\mathbf{a}_3 &&= x_8a\,\hat{\mathbf{x}}+y_8b\,\hat{\mathbf{y}}+z_8c\,\hat{\mathbf{z}} && (8d) && \text{Mg III}\\
\mathbf{B}_{38} &= \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_1-y_8\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_8\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{x}}-y_8b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_8\right)c\,\hat{\mathbf{z}} && (8d) && \text{Mg III}\\
\mathbf{B}_{39} &= -x_8\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_8\right)\mathbf{a}_2-z_8\,\mathbf{a}_3 &&= -x_8a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_8\right)b\,\hat{\mathbf{y}}-z_8c\,\hat{\mathbf{z}} && (8d) && \text{Mg III}\\
\mathbf{B}_{40} &= \left(\tfrac{1}{2}+x_8\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_8\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_8\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_8\right)c\,\hat{\mathbf{z}} && (8d) && \text{Mg III}\\
\mathbf{B}_{41} &= -x_8\,\mathbf{a}_1-y_8\,\mathbf{a}_2-z_8\,\mathbf{a}_3 &&= -x_8a\,\hat{\mathbf{x}}-y_8b\,\hat{\mathbf{y}}-z_8c\,\hat{\mathbf{z}} && (8d) && \text{Mg III}\\
\mathbf{B}_{42} &= \left(\tfrac{1}{2}+x_8\right)\mathbf{a}_1+y_8\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_8\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{x}}+y_8b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_8\right)c\,\hat{\mathbf{z}} && (8d) && \text{Mg III}\\
\mathbf{B}_{43} &= x_8\,\mathbf{a}_1+\left(\tfrac{1}{2}-y_8\right)\mathbf{a}_2+z_8\,\mathbf{a}_3 &&= x_8a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_8\right)b\,\hat{\mathbf{y}}+z_8c\,\hat{\mathbf{z}} && (8d) && \text{Mg III}\\
\mathbf{B}_{44} &= \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_8\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_8\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_8\right)c\,\hat{\mathbf{z}} && (8d) && \text{Mg III}\\
\mathbf{B}_{45} &= x_9\,\mathbf{a}_1+y_9\,\mathbf{a}_2+z_9\,\mathbf{a}_3 &&= x_9a\,\hat{\mathbf{x}}+y_9b\,\hat{\mathbf{y}}+z_9c\,\hat{\mathbf{z}} && (8d) && \text{O III}\\
\mathbf{B}_{46} &= \left(\tfrac{1}{2}-x_9\right)\mathbf{a}_1-y_9\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_9\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_9\right)a\,\hat{\mathbf{x}}-y_9b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_9\right)c\,\hat{\mathbf{z}} && (8d) && \text{O III}\\
\mathbf{B}_{47} &= -x_9\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_9\right)\mathbf{a}_2-z_9\,\mathbf{a}_3 &&= -x_9a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_9\right)b\,\hat{\mathbf{y}}-z_9c\,\hat{\mathbf{z}} && (8d) && \text{O III}\\
\mathbf{B}_{48} &= \left(\tfrac{1}{2}+x_9\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_9\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_9\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_9\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_9\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_9\right)c\,\hat{\mathbf{z}} && (8d) && \text{O III}\\
\mathbf{B}_{49} &= -x_9\,\mathbf{a}_1-y_9\,\mathbf{a}_2-z_9\,\mathbf{a}_3 &&= -x_9a\,\hat{\mathbf{x}}-y_9b\,\hat{\mathbf{y}}-z_9c\,\hat{\mathbf{z}} && (8d) && \text{O III}\\
\mathbf{B}_{50} &= \left(\tfrac{1}{2}+x_9\right)\mathbf{a}_1+y_9\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_9\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_9\right)a\,\hat{\mathbf{x}}+y_9b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_9\right)c\,\hat{\mathbf{z}} && (8d) && \text{O III}\\
\mathbf{B}_{51} &= x_9\,\mathbf{a}_1+\left(\tfrac{1}{2}-y_9\right)\mathbf{a}_2+z_9\,\mathbf{a}_3 &&= x_9a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_9\right)b\,\hat{\mathbf{y}}+z_9c\,\hat{\mathbf{z}} && (8d) && \text{O III}\\
\mathbf{B}_{52} &= \left(\tfrac{1}{2}-x_9\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_9\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_9\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_9\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_9\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_9\right)c\,\hat{\mathbf{z}} && (8d) && \text{O III}\\
\mathbf{B}_{53} &= x_{10}\,\mathbf{a}_1+y_{10}\,\mathbf{a}_2+z_{10}\,\mathbf{a}_3 &&= x_{10}a\,\hat{\mathbf{x}}+y_{10}b\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}} && (8d) && \text{O IV}
\end{aligned}
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\begin{array}{lcccccll}
\mathbf{B}_{54} & = & \left(\frac{1}{2}-x_{10}\right)\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 + \left(\frac{1}{2}+z_{10}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{x}} - y_{10}b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_{10}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O IV} \\
\mathbf{B}_{55} & = & -x_{10}\,\mathbf{a}_1 + \left(\frac{1}{2}+y_{10}\right)\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 & = & -x_{10}a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_{10}\right)b\,\hat{\mathbf{y}} - z_{10}c\,\hat{\mathbf{z}} & (8d) & \text{O IV} \\
\mathbf{B}_{56} & = & \left(\frac{1}{2}+x_{10}\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_{10}\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_{10}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_{10}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_{10}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O IV} \\
\mathbf{B}_{57} & = & -x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 & = & -x_{10}a\,\hat{\mathbf{x}} - y_{10}b\,\hat{\mathbf{y}} - z_{10}c\,\hat{\mathbf{z}} & (8d) & \text{O IV} \\
\mathbf{B}_{58} & = & \left(\frac{1}{2}+x_{10}\right)\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + \left(\frac{1}{2}-z_{10}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{x}} + y_{10}b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_{10}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O IV} \\
\mathbf{B}_{59} & = & x_{10}\,\mathbf{a}_1 + \left(\frac{1}{2}-y_{10}\right)\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 & = & x_{10}a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_{10}\right)b\,\hat{\mathbf{y}} + z_{10}c\,\hat{\mathbf{z}} & (8d) & \text{O IV} \\
\mathbf{B}_{60} & = & \left(\frac{1}{2}-x_{10}\right)\mathbf{a}_1 + \left(\frac{1}{2}+y_{10}\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_{10}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_{10}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_{10}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O IV} \\
\mathbf{B}_{61} & = & x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 & = & x_{11}a\,\hat{\mathbf{x}} + y_{11}b\,\hat{\mathbf{y}} + z_{11}c\,\hat{\mathbf{z}} & (8d) & \text{O V} \\
\mathbf{B}_{62} & = & \left(\frac{1}{2}-x_{11}\right)\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 + \left(\frac{1}{2}+z_{11}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{11}\right)a\,\hat{\mathbf{x}} - y_{11}b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_{11}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O V} \\
\mathbf{B}_{63} & = & -x_{11}\,\mathbf{a}_1 + \left(\frac{1}{2}+y_{11}\right)\mathbf{a}_2 - z_{11}\,\mathbf{a}_3 & = & -x_{11}a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_{11}\right)b\,\hat{\mathbf{y}} - z_{11}c\,\hat{\mathbf{z}} & (8d) & \text{O V} \\
\mathbf{B}_{64} & = & \left(\frac{1}{2}+x_{11}\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_{11}\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_{11}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{11}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_{11}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_{11}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O V} \\
\mathbf{B}_{65} & = & -x_{11}\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3 & = & -x_{11}a\,\hat{\mathbf{x}} - y_{11}b\,\hat{\mathbf{y}} - z_{11}c\,\hat{\mathbf{z}} & (8d) & \text{O V} \\
\mathbf{B}_{66} & = & \left(\frac{1}{2}+x_{11}\right)\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + \left(\frac{1}{2}-z_{11}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{11}\right)a\,\hat{\mathbf{x}} + y_{11}b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_{11}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O V} \\
\mathbf{B}_{67} & = & x_{11}\,\mathbf{a}_1 + \left(\frac{1}{2}-y_{11}\right)\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 & = & x_{11}a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_{11}\right)b\,\hat{\mathbf{y}} + z_{11}c\,\hat{\mathbf{z}} & (8d) & \text{O V} \\
\mathbf{B}_{68} & = & \left(\frac{1}{2}-x_{11}\right)\mathbf{a}_1 + \left(\frac{1}{2}+y_{11}\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_{11}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{11}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_{11}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_{11}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O V} \\
\mathbf{B}_{69} & = & x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 & = & x_{12}a\,\hat{\mathbf{x}} + y_{12}b\,\hat{\mathbf{y}} + z_{12}c\,\hat{\mathbf{z}} & (8d) & \text{O VI} \\
\mathbf{B}_{70} & = & \left(\frac{1}{2}-x_{12}\right)\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 + \left(\frac{1}{2}+z_{12}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{12}\right)a\,\hat{\mathbf{x}} - y_{12}b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_{12}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O VI} \\
\mathbf{B}_{71} & = & -x_{12}\,\mathbf{a}_1 + \left(\frac{1}{2}+y_{12}\right)\mathbf{a}_2 - z_{12}\,\mathbf{a}_3 & = & -x_{12}a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_{12}\right)b\,\hat{\mathbf{y}} - z_{12}c\,\hat{\mathbf{z}} & (8d) & \text{O VI} \\
\mathbf{B}_{72} & = & \left(\frac{1}{2}+x_{12}\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_{12}\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_{12}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{12}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_{12}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_{12}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O VI} \\
\mathbf{B}_{73} & = & -x_{12}\,\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3 & = & -x_{12}a\,\hat{\mathbf{x}} - y_{12}b\,\hat{\mathbf{y}} - z_{12}c\,\hat{\mathbf{z}} & (8d) & \text{O VI} \\
\mathbf{B}_{74} & = & \left(\frac{1}{2}+x_{12}\right)\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + \left(\frac{1}{2}-z_{12}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{12}\right)a\,\hat{\mathbf{x}} + y_{12}b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_{12}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O VI} \\
\mathbf{B}_{75} & = & x_{12}\,\mathbf{a}_1 + \left(\frac{1}{2}-y_{12}\right)\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 & = & x_{12}a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_{12}\right)b\,\hat{\mathbf{y}} + z_{12}c\,\hat{\mathbf{z}} & (8d) & \text{O VI} \\
\mathbf{B}_{76} & = & \left(\frac{1}{2}-x_{12}\right)\mathbf{a}_1 + \left(\frac{1}{2}+y_{12}\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_{12}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{12}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_{12}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_{12}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O VI} \\
\mathbf{B}_{77} & = & x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3 & = & x_{13}a\,\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} + z_{13}c\,\hat{\mathbf{z}} & (8d) & \text{O VII} \\
\mathbf{B}_{78} & = & \left(\frac{1}{2}-x_{13}\right)\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 + \left(\frac{1}{2}+z_{13}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{13}\right)a\,\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_{13}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O VII} \\
\mathbf{B}_{79} & = & -x_{13}\,\mathbf{a}_1 + \left(\frac{1}{2}+y_{13}\right)\mathbf{a}_2 - z_{13}\,\mathbf{a}_3 & = & -x_{13}a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_{13}\right)b\,\hat{\mathbf{y}} - z_{13}c\,\hat{\mathbf{z}} & (8d) & \text{O VII} \\
\mathbf{B}_{80} & = & \left(\frac{1}{2}+x_{13}\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_{13}\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_{13}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{13}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_{13}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_{13}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O VII} \\
\mathbf{B}_{81} & = & -x_{13}\,\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 - z_{13}\,\mathbf{a}_3 & = & -x_{13}a\,\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} - z_{13}c\,\hat{\mathbf{z}} & (8d) & \text{O VII} \\
\mathbf{B}_{82} & = & \left(\frac{1}{2}+x_{13}\right)\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + \left(\frac{1}{2}-z_{13}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{13}\right)a\,\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_{13}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O VII} \\
\mathbf{B}_{83} & = & x_{13}\,\mathbf{a}_1 + \left(\frac{1}{2}-y_{13}\right)\mathbf{a}_2 + z_{13}\,\mathbf{a}_3 & = & x_{13}a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_{13}\right)b\,\hat{\mathbf{y}} + z_{13}c\,\hat{\mathbf{z}} & (8d) & \text{O VII}
\end{array}
$$

$$\mathbf{B}_{84} = \left(\tfrac{1}{2} - x_{13}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{13}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{13}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{13}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{13}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O VII}$$

$$\mathbf{B}_{85} = x_{14}\,\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3 = x_{14}a\,\hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} + z_{14}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O VIII}$$

$$\mathbf{B}_{86} = \left(\tfrac{1}{2} - x_{14}\right)\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{14}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{14}\right)a\,\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{14}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O VIII}$$

$$\mathbf{B}_{87} = -x_{14}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{14}\right)\mathbf{a}_2 - z_{14}\,\mathbf{a}_3 = -x_{14}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{14}\right)b\,\hat{\mathbf{y}} - z_{14}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O VIII}$$

$$\mathbf{B}_{88} = \left(\tfrac{1}{2} + x_{14}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{14}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{14}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{14}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{14}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{14}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O VIII}$$

$$\mathbf{B}_{89} = -x_{14}\,\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 - z_{14}\,\mathbf{a}_3 = -x_{14}a\,\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} - z_{14}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O VIII}$$

$$\mathbf{B}_{90} = \left(\tfrac{1}{2} + x_{14}\right)\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{14}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{14}\right)a\,\hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{14}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O VIII}$$

$$\mathbf{B}_{91} = x_{14}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{14}\right)\mathbf{a}_2 + z_{14}\,\mathbf{a}_3 = x_{14}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{14}\right)b\,\hat{\mathbf{y}} + z_{14}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O VIII}$$

$$\mathbf{B}_{92} = \left(\tfrac{1}{2} - x_{14}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{14}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{14}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{14}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{14}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{14}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O VIII}$$

$$\mathbf{B}_{93} = x_{15}\,\mathbf{a}_1 + y_{15}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 = x_{15}a\,\hat{\mathbf{x}} + y_{15}b\,\hat{\mathbf{y}} + z_{15}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O IX}$$

$$\mathbf{B}_{94} = \left(\tfrac{1}{2} - x_{15}\right)\mathbf{a}_1 - y_{15}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{15}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{15}\right)a\,\hat{\mathbf{x}} - y_{15}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{15}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O IX}$$

$$\mathbf{B}_{95} = -x_{15}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{15}\right)\mathbf{a}_2 - z_{15}\,\mathbf{a}_3 = -x_{15}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{15}\right)b\,\hat{\mathbf{y}} - z_{15}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O IX}$$

$$\mathbf{B}_{96} = \left(\tfrac{1}{2} + x_{15}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{15}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{15}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{15}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{15}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{15}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O IX}$$

$$\mathbf{B}_{97} = -x_{15}\,\mathbf{a}_1 - y_{15}\,\mathbf{a}_2 - z_{15}\,\mathbf{a}_3 = -x_{15}a\,\hat{\mathbf{x}} - y_{15}b\,\hat{\mathbf{y}} - z_{15}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O IX}$$

$$\mathbf{B}_{98} = \left(\tfrac{1}{2} + x_{15}\right)\mathbf{a}_1 + y_{15}\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{15}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{15}\right)a\,\hat{\mathbf{x}} + y_{15}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{15}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O IX}$$

$$\mathbf{B}_{99} = x_{15}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{15}\right)\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 = x_{15}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{15}\right)b\,\hat{\mathbf{y}} + z_{15}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O IX}$$

$$\mathbf{B}_{100} = \left(\tfrac{1}{2} - x_{15}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{15}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{15}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{15}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{15}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{15}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O IX}$$

$$\mathbf{B}_{101} = x_{16}\,\mathbf{a}_1 + y_{16}\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 = x_{16}a\,\hat{\mathbf{x}} + y_{16}b\,\hat{\mathbf{y}} + z_{16}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O X}$$

$$\mathbf{B}_{102} = \left(\tfrac{1}{2} - x_{16}\right)\mathbf{a}_1 - y_{16}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{16}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{16}\right)a\,\hat{\mathbf{x}} - y_{16}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{16}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O X}$$

$$\mathbf{B}_{103} = -x_{16}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{16}\right)\mathbf{a}_2 - z_{16}\,\mathbf{a}_3 = -x_{16}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{16}\right)b\,\hat{\mathbf{y}} - z_{16}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O X}$$

$$\mathbf{B}_{104} = \left(\tfrac{1}{2} + x_{16}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{16}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{16}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{16}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{16}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{16}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O X}$$

$$\mathbf{B}_{105} = -x_{16}\,\mathbf{a}_1 - y_{16}\,\mathbf{a}_2 - z_{16}\,\mathbf{a}_3 = -x_{16}a\,\hat{\mathbf{x}} - y_{16}b\,\hat{\mathbf{y}} - z_{16}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O X}$$

$$\mathbf{B}_{106} = \left(\tfrac{1}{2} + x_{16}\right)\mathbf{a}_1 + y_{16}\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{16}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{16}\right)a\,\hat{\mathbf{x}} + y_{16}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{16}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O X}$$

$$\mathbf{B}_{107} = x_{16}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{16}\right)\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 = x_{16}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{16}\right)b\,\hat{\mathbf{y}} + z_{16}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O X}$$

$$\mathbf{B}_{108} = \left(\tfrac{1}{2} - x_{16}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{16}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{16}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{16}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{16}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{16}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O X}$$

$$\mathbf{B}_{109} = x_{17}\,\mathbf{a}_1 + y_{17}\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 = x_{17}a\,\hat{\mathbf{x}} + y_{17}b\,\hat{\mathbf{y}} + z_{17}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O XI}$$

$$\mathbf{B}_{110} = \left(\tfrac{1}{2} - x_{17}\right)\mathbf{a}_1 - y_{17}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{17}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{17}\right)a\,\hat{\mathbf{x}} - y_{17}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{17}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O XI}$$

$$\mathbf{B}_{111} = -x_{17}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{17}\right)\mathbf{a}_2 - z_{17}\,\mathbf{a}_3 = -x_{17}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{17}\right)b\,\hat{\mathbf{y}} - z_{17}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O XI}$$

$$\mathbf{B}_{112} = \left(\tfrac{1}{2} + x_{17}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{17}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{17}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{17}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{17}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{17}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O XI}$$

$$\mathbf{B}_{113} = -x_{17}\,\mathbf{a}_1 - y_{17}\,\mathbf{a}_2 - z_{17}\,\mathbf{a}_3 = -x_{17}a\,\hat{\mathbf{x}} - y_{17}b\,\hat{\mathbf{y}} - z_{17}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O XI}$$

$$\mathbf{B}_{114} = \left(\tfrac{1}{2}+x_{17}\right)\mathbf{a}_1+y_{17}\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_{17}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{17}\right)a\,\hat{\mathbf{x}}+y_{17}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{17}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O XI}$$

$$\mathbf{B}_{115} = x_{17}\,\mathbf{a}_1+\left(\tfrac{1}{2}-y_{17}\right)\mathbf{a}_2+z_{17}\,\mathbf{a}_3 = x_{17}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{17}\right)b\,\hat{\mathbf{y}}+z_{17}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O XI}$$

$$\mathbf{B}_{116} = \left(\tfrac{1}{2}-x_{17}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{17}\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_{17}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{17}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{17}\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{17}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O XI}$$

$$\mathbf{B}_{117} = x_{18}\,\mathbf{a}_1+y_{18}\,\mathbf{a}_2+z_{18}\,\mathbf{a}_3 = x_{18}a\,\hat{\mathbf{x}}+y_{18}b\,\hat{\mathbf{y}}+z_{18}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O XII}$$

$$\mathbf{B}_{118} = \left(\tfrac{1}{2}-x_{18}\right)\mathbf{a}_1-y_{18}\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_{18}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{18}\right)a\,\hat{\mathbf{x}}-y_{18}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{18}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O XII}$$

$$\mathbf{B}_{119} = -x_{18}\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_{18}\right)\mathbf{a}_2-z_{18}\,\mathbf{a}_3 = -x_{18}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{18}\right)b\,\hat{\mathbf{y}}-z_{18}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O XII}$$

$$\mathbf{B}_{120} = \left(\tfrac{1}{2}+x_{18}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{18}\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_{18}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{18}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{18}\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{18}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O XII}$$

$$\mathbf{B}_{121} = -x_{18}\,\mathbf{a}_1-y_{18}\,\mathbf{a}_2-z_{18}\,\mathbf{a}_3 = -x_{18}a\,\hat{\mathbf{x}}-y_{18}b\,\hat{\mathbf{y}}-z_{18}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O XII}$$

$$\mathbf{B}_{122} = \left(\tfrac{1}{2}+x_{18}\right)\mathbf{a}_1+y_{18}\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_{18}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{18}\right)a\,\hat{\mathbf{x}}+y_{18}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{18}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O XII}$$

$$\mathbf{B}_{123} = x_{18}\,\mathbf{a}_1+\left(\tfrac{1}{2}-y_{18}\right)\mathbf{a}_2+z_{18}\,\mathbf{a}_3 = x_{18}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{18}\right)b\,\hat{\mathbf{y}}+z_{18}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O XII}$$

$$\mathbf{B}_{124} = \left(\tfrac{1}{2}-x_{18}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{18}\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_{18}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{18}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{18}\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{18}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{O XII}$$

$$\mathbf{B}_{125} = x_{19}\,\mathbf{a}_1+y_{19}\,\mathbf{a}_2+z_{19}\,\mathbf{a}_3 = x_{19}a\,\hat{\mathbf{x}}+y_{19}b\,\hat{\mathbf{y}}+z_{19}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{Si I}$$

$$\mathbf{B}_{126} = \left(\tfrac{1}{2}-x_{19}\right)\mathbf{a}_1-y_{19}\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_{19}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{19}\right)a\,\hat{\mathbf{x}}-y_{19}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{19}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{Si I}$$

$$\mathbf{B}_{127} = -x_{19}\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_{19}\right)\mathbf{a}_2-z_{19}\,\mathbf{a}_3 = -x_{19}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{19}\right)b\,\hat{\mathbf{y}}-z_{19}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{Si I}$$

$$\mathbf{B}_{128} = \left(\tfrac{1}{2}+x_{19}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{19}\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_{19}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{19}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{19}\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{19}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{Si I}$$

$$\mathbf{B}_{129} = -x_{19}\,\mathbf{a}_1-y_{19}\,\mathbf{a}_2-z_{19}\,\mathbf{a}_3 = -x_{19}a\,\hat{\mathbf{x}}-y_{19}b\,\hat{\mathbf{y}}-z_{19}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{Si I}$$

$$\mathbf{B}_{130} = \left(\tfrac{1}{2}+x_{19}\right)\mathbf{a}_1+y_{19}\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_{19}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{19}\right)a\,\hat{\mathbf{x}}+y_{19}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{19}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{Si I}$$

$$\mathbf{B}_{131} = x_{19}\,\mathbf{a}_1+\left(\tfrac{1}{2}-y_{19}\right)\mathbf{a}_2+z_{19}\,\mathbf{a}_3 = x_{19}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{19}\right)b\,\hat{\mathbf{y}}+z_{19}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{Si I}$$

$$\mathbf{B}_{132} = \left(\tfrac{1}{2}-x_{19}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{19}\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_{19}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{19}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{19}\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{19}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{Si I}$$

$$\mathbf{B}_{133} = x_{20}\,\mathbf{a}_1+y_{20}\,\mathbf{a}_2+z_{20}\,\mathbf{a}_3 = x_{20}a\,\hat{\mathbf{x}}+y_{20}b\,\hat{\mathbf{y}}+z_{20}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{Si II}$$

$$\mathbf{B}_{134} = \left(\tfrac{1}{2}-x_{20}\right)\mathbf{a}_1-y_{20}\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_{20}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{20}\right)a\,\hat{\mathbf{x}}-y_{20}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{20}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{Si II}$$

$$\mathbf{B}_{135} = -x_{20}\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_{20}\right)\mathbf{a}_2-z_{20}\,\mathbf{a}_3 = -x_{20}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{20}\right)b\,\hat{\mathbf{y}}-z_{20}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{Si II}$$

$$\mathbf{B}_{136} = \left(\tfrac{1}{2}+x_{20}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{20}\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_{20}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{20}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{20}\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{20}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{Si II}$$

$$\mathbf{B}_{137} = -x_{20}\,\mathbf{a}_1-y_{20}\,\mathbf{a}_2-z_{20}\,\mathbf{a}_3 = -x_{20}a\,\hat{\mathbf{x}}-y_{20}b\,\hat{\mathbf{y}}-z_{20}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{Si II}$$

$$\mathbf{B}_{138} = \left(\tfrac{1}{2}+x_{20}\right)\mathbf{a}_1+y_{20}\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_{20}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{20}\right)a\,\hat{\mathbf{x}}+y_{20}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{20}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{Si II}$$

$$\mathbf{B}_{139} = x_{20}\,\mathbf{a}_1+\left(\tfrac{1}{2}-y_{20}\right)\mathbf{a}_2+z_{20}\,\mathbf{a}_3 = x_{20}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{20}\right)b\,\hat{\mathbf{y}}+z_{20}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{Si II}$$

$$\mathbf{B}_{140} = \left(\tfrac{1}{2}-x_{20}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{20}\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_{20}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{20}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{20}\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{20}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{Si II}$$

$$\mathbf{B}_{141} = x_{21}\,\mathbf{a}_1+y_{21}\,\mathbf{a}_2+z_{21}\,\mathbf{a}_3 = x_{21}a\,\hat{\mathbf{x}}+y_{21}b\,\hat{\mathbf{y}}+z_{21}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{Si III}$$

$$\mathbf{B}_{142} = \left(\tfrac{1}{2}-x_{21}\right)\mathbf{a}_1-y_{21}\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_{21}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{21}\right)a\,\hat{\mathbf{x}}-y_{21}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{21}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{Si III}$$

$$\mathbf{B}_{143} = -x_{21}\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_{21}\right)\mathbf{a}_2-z_{21}\,\mathbf{a}_3 = -x_{21}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{21}\right)b\,\hat{\mathbf{y}}-z_{21}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{Si III}$$

$\mathbf{B}_{144}$	=	$\left(\frac{1}{2}+x_{21}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{21}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{21}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{21}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{21}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{21}\right)c\,\hat{\mathbf{z}}$	(8*d*)	Si III
$\mathbf{B}_{145}$	=	$-x_{21}\,\mathbf{a}_1-y_{21}\,\mathbf{a}_2-z_{21}\,\mathbf{a}_3$	=	$-x_{21}a\,\hat{\mathbf{x}}-y_{21}b\,\hat{\mathbf{y}}-z_{21}c\,\hat{\mathbf{z}}$	(8*d*)	Si III
$\mathbf{B}_{146}$	=	$\left(\frac{1}{2}+x_{21}\right)\mathbf{a}_1+y_{21}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{21}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{21}\right)a\,\hat{\mathbf{x}}+y_{21}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{21}\right)c\,\hat{\mathbf{z}}$	(8*d*)	Si III
$\mathbf{B}_{147}$	=	$x_{21}\,\mathbf{a}_1+\left(\frac{1}{2}-y_{21}\right)\mathbf{a}_2+z_{21}\,\mathbf{a}_3$	=	$x_{21}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{21}\right)b\,\hat{\mathbf{y}}+z_{21}c\,\hat{\mathbf{z}}$	(8*d*)	Si III
$\mathbf{B}_{148}$	=	$\left(\frac{1}{2}-x_{21}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{21}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{21}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{21}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{21}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{21}\right)c\,\hat{\mathbf{z}}$	(8*d*)	Si III
$\mathbf{B}_{149}$	=	$x_{22}\,\mathbf{a}_1+y_{22}\,\mathbf{a}_2+z_{22}\,\mathbf{a}_3$	=	$x_{22}a\,\hat{\mathbf{x}}+y_{22}b\,\hat{\mathbf{y}}+z_{22}c\,\hat{\mathbf{z}}$	(8*d*)	Si IV
$\mathbf{B}_{150}$	=	$\left(\frac{1}{2}-x_{22}\right)\mathbf{a}_1-y_{22}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{22}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{22}\right)a\,\hat{\mathbf{x}}-y_{22}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{22}\right)c\,\hat{\mathbf{z}}$	(8*d*)	Si IV
$\mathbf{B}_{151}$	=	$-x_{22}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{22}\right)\mathbf{a}_2-z_{22}\,\mathbf{a}_3$	=	$-x_{22}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{22}\right)b\,\hat{\mathbf{y}}-z_{22}c\,\hat{\mathbf{z}}$	(8*d*)	Si IV
$\mathbf{B}_{152}$	=	$\left(\frac{1}{2}+x_{22}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{22}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{22}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{22}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{22}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{22}\right)c\,\hat{\mathbf{z}}$	(8*d*)	Si IV
$\mathbf{B}_{153}$	=	$-x_{22}\,\mathbf{a}_1-y_{22}\,\mathbf{a}_2-z_{22}\,\mathbf{a}_3$	=	$-x_{22}a\,\hat{\mathbf{x}}-y_{22}b\,\hat{\mathbf{y}}-z_{22}c\,\hat{\mathbf{z}}$	(8*d*)	Si IV
$\mathbf{B}_{154}$	=	$\left(\frac{1}{2}+x_{22}\right)\mathbf{a}_1+y_{22}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{22}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{22}\right)a\,\hat{\mathbf{x}}+y_{22}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{22}\right)c\,\hat{\mathbf{z}}$	(8*d*)	Si IV
$\mathbf{B}_{155}$	=	$x_{22}\,\mathbf{a}_1+\left(\frac{1}{2}-y_{22}\right)\mathbf{a}_2+z_{22}\,\mathbf{a}_3$	=	$x_{22}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{22}\right)b\,\hat{\mathbf{y}}+z_{22}c\,\hat{\mathbf{z}}$	(8*d*)	Si IV
$\mathbf{B}_{156}$	=	$\left(\frac{1}{2}-x_{22}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{22}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{22}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{22}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{22}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{22}\right)c\,\hat{\mathbf{z}}$	(8*d*)	Si IV

References:

- E. M. Walitzi, F. Walter, and K. Ettinger, *Verfeinerung der Kristallstruktur von Anthophyllit vom Ochsenkogel/Gleinalpe, Österreich*, Zeitschrift für Kristallographie - Crystalline Materials **188**, 237–244 (1989), doi:10.1524/zkri.1989.188.14.237.
- B. E. Warren and D. I. Modell, *The Structure of Anthophyllite $H_2Mg_7(SiO_3)_8$*, Zeitschrift für Kristallographie - Crystalline Materials **75**, 161–178 (1930), doi:10.1515/zkri-1930-0112.
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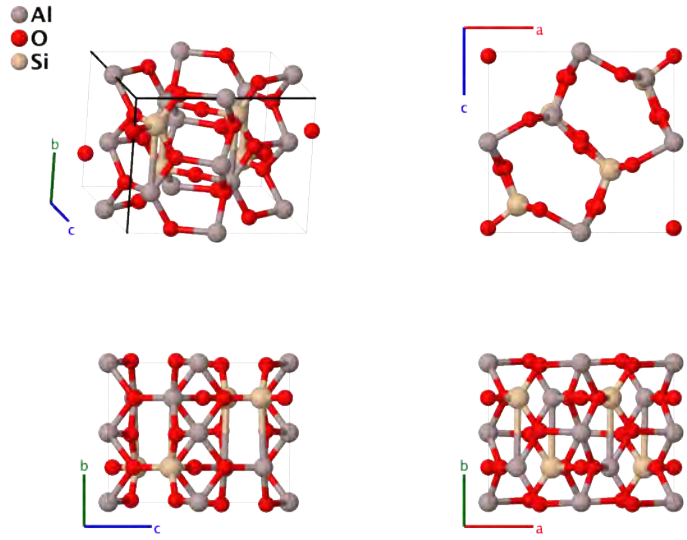
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1630
- POSCAR: pp. 1630

Sillimanite (Al_2SiO_5, $S0_3$) Structure: A2B5C_oP32_62_bc_3cd_c

http://aflow.org/prototype-encyclopedia/A2B5C_oP32_62_bc_3cd_c

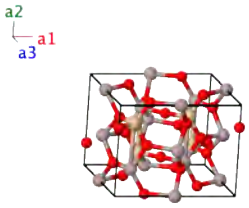


Prototype	:	Al_2O_5Si
AFLOW prototype label	:	A2B5C_oP32_62_bc_3cd_c
***Strukturbericht* designation**	:	$S0_3$
Pearson symbol	:	oP32
Space group number	:	62
Space group symbol	:	*Pnma*
AFLOW prototype command	:	`aflow --proto=A2B5C_oP32_62_bc_3cd_c --params=`$a, b/a, c/a, x_2, z_2, x_3, z_3, x_4, z_4, x_5, z_5, x_6, z_6, x_7, y_7, z_7$

- Three crystal polymorphs of Al_2SiO_5 have been characterized: kyanite ($S0_1$), space group $P\bar{1}$ #2, andalusite ($S0_2$), space group *Pnnm* #58, and sillimanite ($S0_3$), space group *Pnma* #62. All are characterized chains of edge-sharing SiO_6 tetrahedra and Al octahedra.
- We use the ambient pressure data of (Yang, 1997).
- (Yang, 1997) give the Wyckoff positions in the *Pbnm* setting of space group #62. We use FINDSYM to transform this to the standard *Pnma* setting. This involves a rotation of the principle axes and a shift of the set of aluminum atoms from the (4*a*) to the (4*b*) Wyckoff position.
- (Hermann, 1937) defined this as $S0_3$, but also listed it as $H5_2$ in the index.

Simple Orthorhombic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= a\,\hat{\mathbf{x}}\\ \mathbf{a}_2 &= b\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$\frac{1}{2}\,\mathbf{a}_3$	$\frac{1}{2}c\,\hat{\mathbf{z}}$	$(4b)$	Al I
$\mathbf{B}_2$	$\frac{1}{2}\,\mathbf{a}_1$	$\frac{1}{2}a\,\hat{\mathbf{x}}$	$(4b)$	Al I
$\mathbf{B}_3$	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	$\frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(4b)$	Al I
$\mathbf{B}_4$	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$	$(4b)$	Al I
$\mathbf{B}_5$	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	$x_2a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_2c\,\hat{\mathbf{z}}$	$(4c)$	Al II
$\mathbf{B}_6$	$\left(\frac{1}{2} - x_2\right)\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\,\mathbf{a}_3$	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	$(4c)$	Al II
$\mathbf{B}_7$	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	$-x_2a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_2c\,\hat{\mathbf{z}}$	$(4c)$	Al II
$\mathbf{B}_8$	$\left(\frac{1}{2} + x_2\right)\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	$(4c)$	Al II
$\mathbf{B}_9$	$x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$x_3a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	$(4c)$	O I
$\mathbf{B}_{10}$	$\left(\frac{1}{2} - x_3\right)\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\,\mathbf{a}_3$	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	$(4c)$	O I
$\mathbf{B}_{11}$	$-x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	$-x_3a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	$(4c)$	O I
$\mathbf{B}_{12}$	$\left(\frac{1}{2} + x_3\right)\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	$(4c)$	O I
$\mathbf{B}_{13}$	$x_4\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$x_4a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	$(4c)$	O II
$\mathbf{B}_{14}$	$\left(\frac{1}{2} - x_4\right)\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\,\mathbf{a}_3$	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	$(4c)$	O II
$\mathbf{B}_{15}$	$-x_4\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	$-x_4a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	$(4c)$	O II
$\mathbf{B}_{16}$	$\left(\frac{1}{2} + x_4\right)\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	$(4c)$	O II
$\mathbf{B}_{17}$	$x_5\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	$x_5a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_5c\,\hat{\mathbf{z}}$	$(4c)$	O III
$\mathbf{B}_{18}$	$\left(\frac{1}{2} - x_5\right)\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\,\mathbf{a}_3$	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	$(4c)$	O III
$\mathbf{B}_{19}$	$-x_5\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	$-x_5a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_5c\,\hat{\mathbf{z}}$	$(4c)$	O III
$\mathbf{B}_{20}$	$\left(\frac{1}{2} + x_5\right)\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	$(4c)$	O III
$\mathbf{B}_{21}$	$x_6\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	$x_6a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_6c\,\hat{\mathbf{z}}$	$(4c)$	Si
$\mathbf{B}_{22}$	$\left(\frac{1}{2} - x_6\right)\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\,\mathbf{a}_3$	$\left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\,\hat{\mathbf{z}}$	$(4c)$	Si
$\mathbf{B}_{23}$	$-x_6\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	$-x_6a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_6c\,\hat{\mathbf{z}}$	$(4c)$	Si
$\mathbf{B}_{24}$	$\left(\frac{1}{2} + x_6\right)\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\,\hat{\mathbf{z}}$	$(4c)$	Si
$\mathbf{B}_{25}$	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	$x_7a\,\hat{\mathbf{x}} + y_7b\,\hat{\mathbf{y}} + z_7c\,\hat{\mathbf{z}}$	$(8d)$	O IV
$\mathbf{B}_{26}$	$\left(\frac{1}{2} - x_7\right)\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\,\mathbf{a}_3$	$\left(\frac{1}{2} - x_7\right)a\,\hat{\mathbf{x}} - y_7b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right)c\,\hat{\mathbf{z}}$	$(8d)$	O IV
$\mathbf{B}_{27}$	$-x_7\,\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	$-x_7a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right)b\,\hat{\mathbf{y}} - z_7c\,\hat{\mathbf{z}}$	$(8d)$	O IV

$\mathbf{B}_{28}$	=	$\left(\frac{1}{2}+x_7\right)\mathbf{a}_1+\left(\frac{1}{2}-y_7\right)\mathbf{a}_2+\left(\frac{1}{2}-z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_7\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_7\right)c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{29}$	=	$-x_7\,\mathbf{a}_1-y_7\,\mathbf{a}_2-z_7\,\mathbf{a}_3$	=	$-x_7a\,\hat{\mathbf{x}}-y_7b\,\hat{\mathbf{y}}-z_7c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2}+x_7\right)\mathbf{a}_1+y_7\,\mathbf{a}_2+\left(\frac{1}{2}-z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{x}}+y_7b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_7\right)c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{31}$	=	$x_7\,\mathbf{a}_1+\left(\frac{1}{2}-y_7\right)\mathbf{a}_2+z_7\,\mathbf{a}_3$	=	$x_7a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_7\right)b\,\hat{\mathbf{y}}+z_7c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2}-x_7\right)\mathbf{a}_1+\left(\frac{1}{2}+y_7\right)\mathbf{a}_2+\left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_7\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_7\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	(8*d*)	O IV

References:

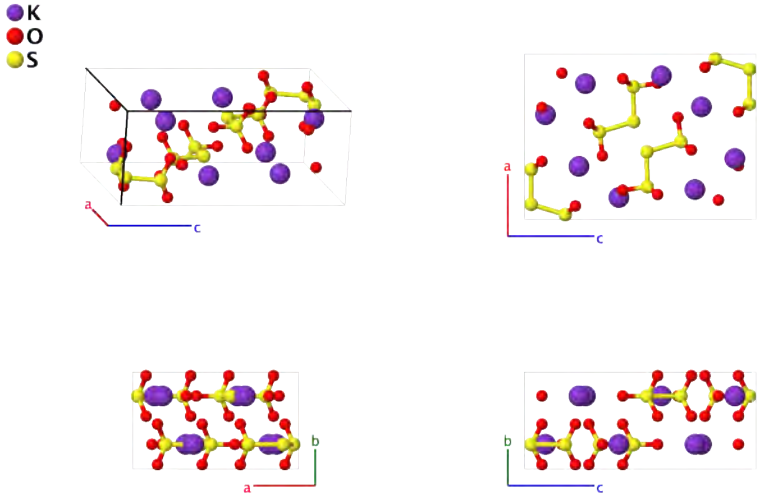
- H. Yang, R. M. Hazen, L. W. Finger, C. T. Prewitt, and R. T. Downs, *Compressibility and crystal structure of sillimanite, Al_2SiO_5, at high pressure*, Phys. Chem. Miner. **25**, 39–47 (1997), doi:10.1007/s002690050084.
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1631
- POSCAR: pp. 1632

$K_2S_3O_6$ ($K5_1$) Structure: A2B6C3_oP44_62_2c_2c2d_3c

http://aflow.org/prototype-encyclopedia/A2B6C3_oP44_62_2c_2c2d_3c



Prototype	:	$K_2O_6S_3$
AFLOW prototype label	:	A2B6C3_oP44_62_2c_2c2d_3c
***Strukturbericht* designation**	:	$K5_1$
Pearson symbol	:	oP44
Space group number	:	62
Space group symbol	:	*Pnma*
AFLOW prototype command	:	`aflow --proto=A2B6C3_oP44_62_2c_2c2d_3c --params=`$a, b/a, c/a, x_1, z_1, x_2, z_2, x_3, z_3, x_4, z_4, x_5, z_5, x_6, z_6, x_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9$

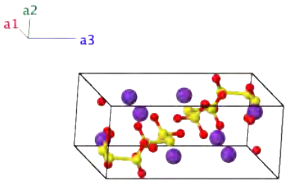
Other compounds with this structure

- $Rb_2S_3O_6$

- (Stewart, 1979) give the Wyckoff positions of this structure using the *Pnam* orientation of space group #62. We have used FINDSYM to change this to our standard *Pnma* orientation.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(4*c*)	K I
$\mathbf{B}_2$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	K I
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	(4*c*)	K I
$\mathbf{B}_4$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	K I
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*c*)	K II
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	(4*c*)	K II
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(4*c*)	K II
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right) c\,\hat{\mathbf{z}}$	(4*c*)	K II
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right) c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(4*c*)	S I
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	(4*c*)	S I
$\mathbf{B}_{19}$	=	$-x_5\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(4*c*)	S I
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right) c\,\hat{\mathbf{z}}$	(4*c*)	S I
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(4*c*)	S II
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	(4*c*)	S II
$\mathbf{B}_{23}$	=	$-x_6\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(4*c*)	S II
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right) c\,\hat{\mathbf{z}}$	(4*c*)	S II
$\mathbf{B}_{25}$	=	$x_7\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(4*c*)	S III
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} - x_7\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_7\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right) c\,\hat{\mathbf{z}}$	(4*c*)	S III
$\mathbf{B}_{27}$	=	$-x_7\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	(4*c*)	S III
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} + x_7\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_7\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right) c\,\hat{\mathbf{z}}$	(4*c*)	S III
$\mathbf{B}_{29}$	=	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2} - x_8\right)\mathbf{a}_1 - y_8\,\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_8\right) a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right) c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{31}$	=	$-x_8\,\mathbf{a}_1 + \left(\frac{1}{2} + y_8\right)\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_8\right) b\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2} + x_8\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_8\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_8\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_8\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right) c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{33}$	=	$-x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{34}$	=	$\left(\frac{1}{2} + x_8\right)\mathbf{a}_1 + y_8\,\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_8\right) a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right) c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{35}$	=	$x_8\,\mathbf{a}_1 + \left(\frac{1}{2} - y_8\right)\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_8\right) b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(8*d*)	O III

$\mathbf{B}_{36}$	=	$\left(\frac{1}{2}-x_8\right)\mathbf{a}_1+\left(\frac{1}{2}+y_8\right)\mathbf{a}_2+\left(\frac{1}{2}+z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_8\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_8\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_8\right)c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{37}$	=	$x_9\,\mathbf{a}_1+y_9\,\mathbf{a}_2+z_9\,\mathbf{a}_3$	=	$x_9a\,\hat{\mathbf{x}}+y_9b\,\hat{\mathbf{y}}+z_9c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{38}$	=	$\left(\frac{1}{2}-x_9\right)\mathbf{a}_1-y_9\,\mathbf{a}_2+\left(\frac{1}{2}+z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_9\right)a\,\hat{\mathbf{x}}-y_9b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_9\right)c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{39}$	=	$-x_9\,\mathbf{a}_1+\left(\frac{1}{2}+y_9\right)\mathbf{a}_2-z_9\,\mathbf{a}_3$	=	$-x_9a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_9\right)b\,\hat{\mathbf{y}}-z_9c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{40}$	=	$\left(\frac{1}{2}+x_9\right)\mathbf{a}_1+\left(\frac{1}{2}-y_9\right)\mathbf{a}_2+\left(\frac{1}{2}-z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_9\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_9\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_9\right)c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{41}$	=	$-x_9\,\mathbf{a}_1-y_9\,\mathbf{a}_2-z_9\,\mathbf{a}_3$	=	$-x_9a\,\hat{\mathbf{x}}-y_9b\,\hat{\mathbf{y}}-z_9c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{42}$	=	$\left(\frac{1}{2}+x_9\right)\mathbf{a}_1+y_9\,\mathbf{a}_2+\left(\frac{1}{2}-z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_9\right)a\,\hat{\mathbf{x}}+y_9b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_9\right)c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{43}$	=	$x_9\,\mathbf{a}_1+\left(\frac{1}{2}-y_9\right)\mathbf{a}_2+z_9\,\mathbf{a}_3$	=	$x_9a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_9\right)b\,\hat{\mathbf{y}}+z_9c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{44}$	=	$\left(\frac{1}{2}-x_9\right)\mathbf{a}_1+\left(\frac{1}{2}+y_9\right)\mathbf{a}_2+\left(\frac{1}{2}+z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_9\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_9\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_9\right)c\,\hat{\mathbf{z}}$	(8*d*)	O IV

References:

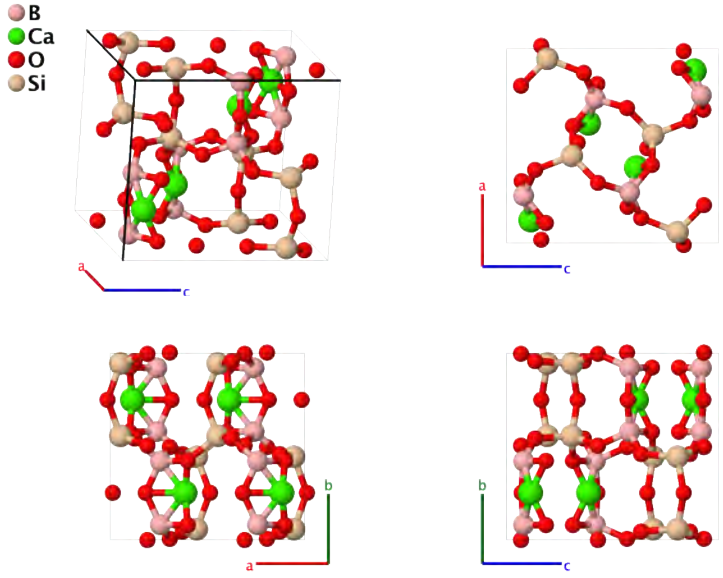
- J. M. Stewart and J. T. Szymański, *A redetermination of the crystal structure of potassium trithionate, $K_2S_3O_6$*, Acta Crystallogr. Sect. B Struct. Sci. **35**, 1967–1970 (1979), doi:10.1107/S0567740879008268.

Geometry files:

- CIF: pp. 1632
- POSCAR: pp. 1632

Danburite ($CaB_2Si_2O_8$, $S6_3$) Structure: A2BC8D2_oP52_62_d_c_2c3d_d

http://aflow.org/prototype-encyclopedia/A2BC8D2_oP52_62_d_c_2c3d_d

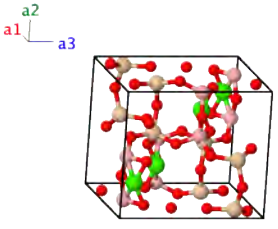


Prototype	:	$B_2CaO_8Si_2$
AFLOW prototype label	:	A2BC8D2_oP52_62_d_c_2c3d_d
***Strukturbericht* designation**	:	$S6_3$
Pearson symbol	:	oP52
Space group number	:	62
Space group symbol	:	*Pnma*
AFLOW prototype command	:	`aflow --proto=A2BC8D2_oP52_62_d_c_2c3d_d --params=`$a, b/a, c/a, x_1, z_1, x_2, z_2, x_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8$

- We use the 25 °C data from (Sugiyama, 1985). They presented the structure in the *Pnam* orientation of space group #62. We used FINDSYM to transform this to the standard *Pnma* orientation.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(4*c*)	Ca
$\mathbf{B}_2$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	Ca
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	(4*c*)	Ca
$\mathbf{B}_4$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	Ca
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right) c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right) c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*d*)	B
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	(8*d*)	B
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right) b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*d*)	B
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\,\hat{\mathbf{z}}$	(8*d*)	B
$\mathbf{B}_{17}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*d*)	B
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\,\hat{\mathbf{z}}$	(8*d*)	B
$\mathbf{B}_{19}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right) b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*d*)	B
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	(8*d*)	B
$\mathbf{B}_{21}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right) a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{23}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right) b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right) c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{25}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + y_5\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right) c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{27}$	=	$x_5\,\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right) b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{29}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{31}$	=	$-x_6\,\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right) b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right) c\,\hat{\mathbf{z}}$	(8*d*)	O IV

$\mathbf{B}_{33}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{34}$	=	$\left(\frac{1}{2}+x_6\right)\mathbf{a}_1 + y_6\,\mathbf{a}_2 + \left(\frac{1}{2}-z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_6\right)c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{35}$	=	$x_6\,\mathbf{a}_1 + \left(\frac{1}{2}-y_6\right)\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_6\right)b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{36}$	=	$\left(\frac{1}{2}-x_6\right)\mathbf{a}_1 + \left(\frac{1}{2}+y_6\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_6\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{37}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(8*d*)	O V
$\mathbf{B}_{38}$	=	$\left(\frac{1}{2}-x_7\right)\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_7\right)a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	(8*d*)	O V
$\mathbf{B}_{39}$	=	$-x_7\,\mathbf{a}_1 + \left(\frac{1}{2}+y_7\right)\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_7\right)b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	(8*d*)	O V
$\mathbf{B}_{40}$	=	$\left(\frac{1}{2}+x_7\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_7\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_7\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_7\right)c\,\hat{\mathbf{z}}$	(8*d*)	O V
$\mathbf{B}_{41}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	(8*d*)	O V
$\mathbf{B}_{42}$	=	$\left(\frac{1}{2}+x_7\right)\mathbf{a}_1 + y_7\,\mathbf{a}_2 + \left(\frac{1}{2}-z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_7\right)c\,\hat{\mathbf{z}}$	(8*d*)	O V
$\mathbf{B}_{43}$	=	$x_7\,\mathbf{a}_1 + \left(\frac{1}{2}-y_7\right)\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_7\right)b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(8*d*)	O V
$\mathbf{B}_{44}$	=	$\left(\frac{1}{2}-x_7\right)\mathbf{a}_1 + \left(\frac{1}{2}+y_7\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_7\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	(8*d*)	O V
$\mathbf{B}_{45}$	=	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(8*d*)	Si
$\mathbf{B}_{46}$	=	$\left(\frac{1}{2}-x_8\right)\mathbf{a}_1 - y_8\,\mathbf{a}_2 + \left(\frac{1}{2}+z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_8\right)a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_8\right)c\,\hat{\mathbf{z}}$	(8*d*)	Si
$\mathbf{B}_{47}$	=	$-x_8\,\mathbf{a}_1 + \left(\frac{1}{2}+y_8\right)\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_8\right)b\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(8*d*)	Si
$\mathbf{B}_{48}$	=	$\left(\frac{1}{2}+x_8\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_8\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_8\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_8\right)c\,\hat{\mathbf{z}}$	(8*d*)	Si
$\mathbf{B}_{49}$	=	$-x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(8*d*)	Si
$\mathbf{B}_{50}$	=	$\left(\frac{1}{2}+x_8\right)\mathbf{a}_1 + y_8\,\mathbf{a}_2 + \left(\frac{1}{2}-z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_8\right)a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_8\right)c\,\hat{\mathbf{z}}$	(8*d*)	Si
$\mathbf{B}_{51}$	=	$x_8\,\mathbf{a}_1 + \left(\frac{1}{2}-y_8\right)\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_8\right)b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(8*d*)	Si
$\mathbf{B}_{52}$	=	$\left(\frac{1}{2}-x_8\right)\mathbf{a}_1 + \left(\frac{1}{2}+y_8\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_8\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_8\right)c\,\hat{\mathbf{z}}$	(8*d*)	Si

References:

- K. Sugiyama and Y. Takéuchi, *Unusual thermal expansion of a B-O bond in the structure of danburite* $CaB_2Si_2O_8$, Zeitschrift für Kristallographie - Crystalline Materials **173**, 293–304 (1985), doi:10.1524/zkri.1985.173.14.293.

Found in:

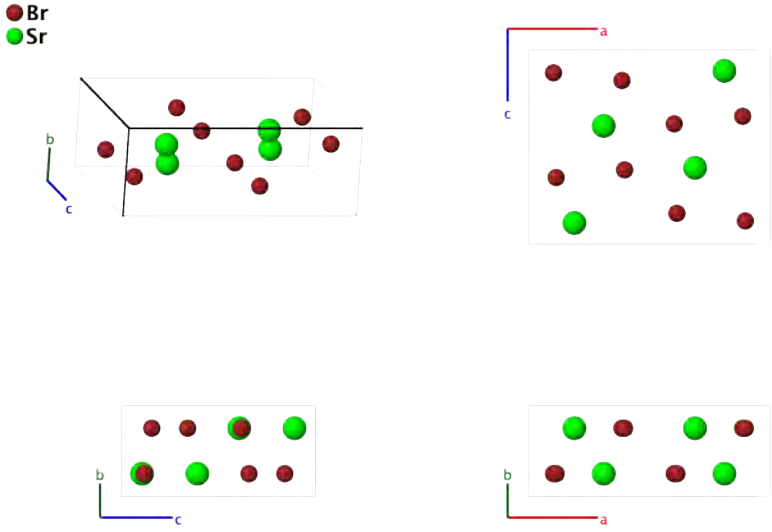
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1632
- POSCAR: pp. 1633

*C*53 ($SrBr_2$) (*obsolete*) Structure: A2B_oP12_62_2c_c

http://aflow.org/prototype-encyclopedia/A2B_oP12_62_2c_c.SrBr2



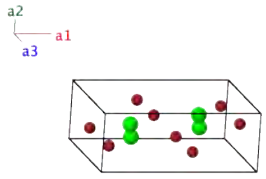
Prototype	:	Br_2Sr
AFLOW prototype label	:	A2B_oP12_62_2c_c
***Strukturbericht* designation**	:	*C*53
Pearson symbol	:	oP12
Space group number	:	62
Space group symbol	:	*Pnma*
AFLOW prototype command	:	`aflow --proto=A2B_oP12_62_2c_c --params=`$a, b/a, c/a, x_1, z_1, x_2, z_2, x_3, z_3$

- (Kamermans, 1939) determined that $SrBr_2$ had a distorted $PbCl_2$ structure, and assigned it to space group *Pnma* #62. (Hermann, 1939) gave this the *Strukturbericht* designation *C*53.
- Some years later, (Sass, 1963) showed that this structure was incorrect, and that α-$SrBr_2$ was actually in space group $P4/n$ #85, rendering the *C*53 symbol obsolete. We include it here for historical interest.
- (Kamermans, 1939) gave the atomic positions in terms of the *Pbnm* setting of space group #62. We have used FINDSYM to translate this to the standard *Pnma* setting.
- Note that
  - $PbCl_2$ *C*23,
  - $HgCl_2$ *C*25,
  - SrH_2 *C*29,
  - Co_2Si *C*37, and
  - $SrBr_2$ *C*53

all have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters (`--params`) specified in their corresponding CIF files.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	$x_1 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	($4c$)	Br I
$\mathbf{B}_2$	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	($4c$)	Br I
$\mathbf{B}_3$	$-x_1\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	$-x_1 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	($4c$)	Br I
$\mathbf{B}_4$	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_1\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right) c\,\hat{\mathbf{z}}$	($4c$)	Br I
$\mathbf{B}_5$	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	$x_2 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	($4c$)	Br II
$\mathbf{B}_6$	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	($4c$)	Br II
$\mathbf{B}_7$	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	$-x_2 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	($4c$)	Br II
$\mathbf{B}_8$	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right) c\,\hat{\mathbf{z}}$	($4c$)	Br II
$\mathbf{B}_9$	$x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$x_3 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	($4c$)	Sr
$\mathbf{B}_{10}$	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_3\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	($4c$)	Sr
$\mathbf{B}_{11}$	$-x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	$-x_3 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	($4c$)	Sr
$\mathbf{B}_{12}$	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_3\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right) c\,\hat{\mathbf{z}}$	($4c$)	Sr

References:

- M. A. Kamermans, *The Crystal Structure of $SrBr_2$*, Zeitschrift für Kristallographie - Crystalline Materials **101**, 406–411 (1939), doi:10.1524/zkri.1939.101.1.406. C53 (obsolete) $SrBr_2$ Structure.
- R. L. Sass, T. Brackett, and E. Brackett, *The Crystal Structure Of Strontium Bromide*, J. Phys. Chem. **67**, 2862–2863 (1963), doi:10.1021/j100806a516.

Found in:

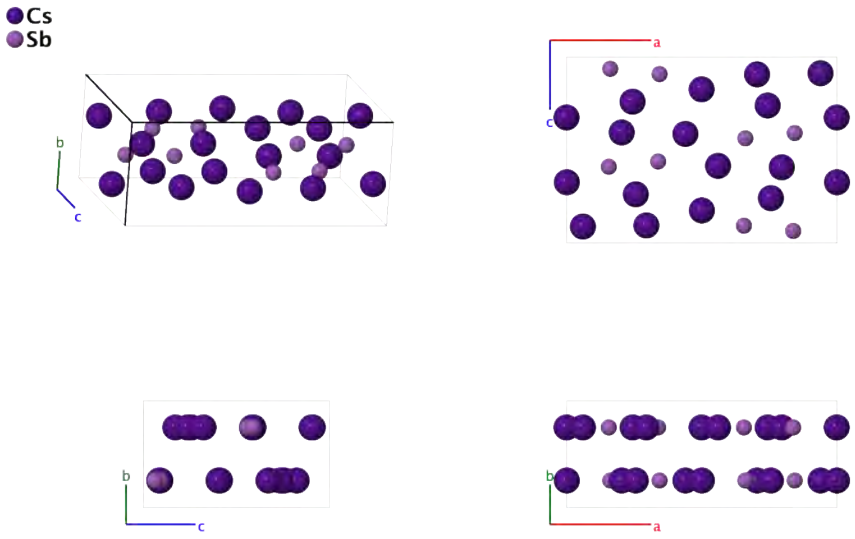
- K. Herrmann, ed., *Strukturbericht Band VII 1939* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1943).

Geometry files:

- CIF: pp. 1633
- POSCAR: pp. 1633

Cs_2Sb Structure: A2B_oP24_62_4c_2c

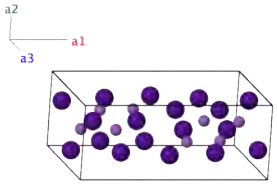
http://aflow.org/prototype-encyclopedia/A2B_oP24_62_4c_2c



Prototype	:	Cs_2Sb
AFLOW prototype label	:	A2B_oP24_62_4c_2c
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP24
Space group number	:	62
Space group symbol	:	*Pnma*
AFLOW prototype command	:	`aflow --proto=A2B_oP24_62_4c_2c --params=`$a, b/a, c/a, x_1, z_1, x_2, z_2, x_3, z_3, x_4, z_4, x_5, z_5, x_6, z_6$

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(4*c*)	Cs I
$\mathbf{B}_2$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	Cs I
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	(4*c*)	Cs I
$\mathbf{B}_4$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	Cs I
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*c*)	Cs II

$\mathbf{B}_6$	=	$\left(\frac{1}{2}-x_2\right)\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2+\left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_2\right)a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_2\right)c\,\hat{\mathbf{z}}$	(4*c*)	Cs II
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2-z_2\,\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}-z_2c\,\hat{\mathbf{z}}$	(4*c*)	Cs II
$\mathbf{B}_8$	=	$\left(\frac{1}{2}+x_2\right)\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+\left(\frac{1}{2}-z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_2\right)a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_2\right)c\,\hat{\mathbf{z}}$	(4*c*)	Cs II
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+z_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+z_3c\,\hat{\mathbf{z}}$	(4*c*)	Cs III
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2}-x_3\right)\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2+\left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_3\right)c\,\hat{\mathbf{z}}$	(4*c*)	Cs III
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2-z_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}-z_3c\,\hat{\mathbf{z}}$	(4*c*)	Cs III
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2}+x_3\right)\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+\left(\frac{1}{2}-z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_3\right)c\,\hat{\mathbf{z}}$	(4*c*)	Cs III
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+z_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+z_4c\,\hat{\mathbf{z}}$	(4*c*)	Cs IV
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2}-x_4\right)\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2+\left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_4\right)c\,\hat{\mathbf{z}}$	(4*c*)	Cs IV
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2-z_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}-z_4c\,\hat{\mathbf{z}}$	(4*c*)	Cs IV
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2}+x_4\right)\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+\left(\frac{1}{2}-z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_4\right)c\,\hat{\mathbf{z}}$	(4*c*)	Cs IV
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+z_5\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+z_5c\,\hat{\mathbf{z}}$	(4*c*)	Sb I
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2}-x_5\right)\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	(4*c*)	Sb I
$\mathbf{B}_{19}$	=	$-x_5\,\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2-z_5\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}-z_5c\,\hat{\mathbf{z}}$	(4*c*)	Sb I
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2}+x_5\right)\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+\left(\frac{1}{2}-z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_5\right)c\,\hat{\mathbf{z}}$	(4*c*)	Sb I
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+z_6\,\mathbf{a}_3$	=	$x_6a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+z_6c\,\hat{\mathbf{z}}$	(4*c*)	Sb II
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2}-x_6\right)\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2+\left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	(4*c*)	Sb II
$\mathbf{B}_{23}$	=	$-x_6\,\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2-z_6\,\mathbf{a}_3$	=	$-x_6a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}-z_6c\,\hat{\mathbf{z}}$	(4*c*)	Sb II
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2}+x_6\right)\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+\left(\frac{1}{2}-z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_6\right)c\,\hat{\mathbf{z}}$	(4*c*)	Sb II

References:

- C. Hirschle and C. Röhr, *Darstellung und Kristallstruktur der bekannten Zintl-Phasen Cs_3Sb_7 und Cs_4Sb_2*, Z. Anorg. Allg. Chem. **626**, 1992–1998 (2000), doi:10.1002/1521-3749(200009)626:9%3C1992::AID-ZAAC1992%3E3.0.CO;2-G.

Found in:

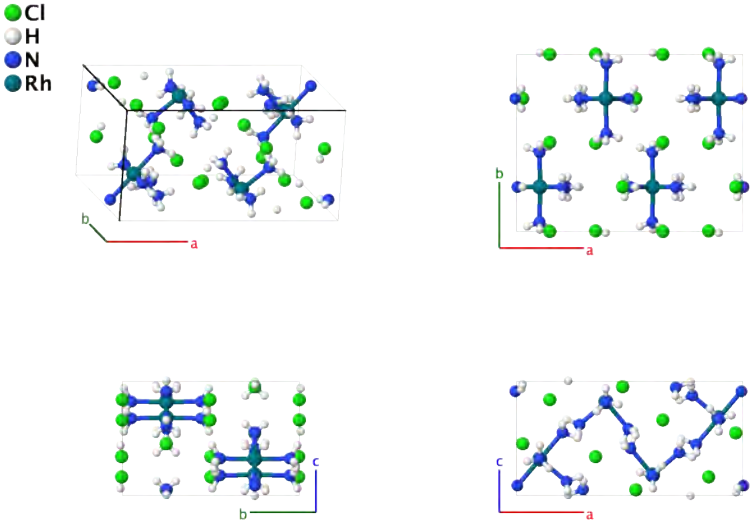
- P. Villars, ed., *PAULING FILE in: Inorganic Solid Phases (online database)* (Springer Materials, Heidelberg, 2016). Cs_4Sb_2 (Cs_2Sb) Crystal Structure.

Geometry files:

- CIF: pp. 1634
- POSCAR: pp. 1634

$RhCl_2(NH_3)_5Cl$ ($J1_8$) Structure: A3B15C5D_oP96_62_cd_3c6d_3cd_c

http://aflow.org/prototype-encyclopedia/A3B15C5D_oP96_62_cd_3c6d_3cd_c



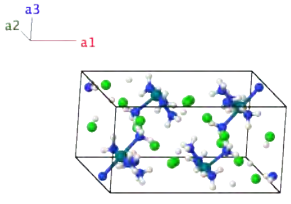
Prototype	:	$Cl_3H_{15}N_5Rh$
AFLOW prototype label	:	A3B15C5D_oP96_62_cd_3c6d_3cd_c
***Strukturbericht* designation**	:	$J1_8$
Pearson symbol	:	oP96
Space group number	:	62
Space group symbol	:	*Pnma*
AFLOW prototype command	:	`aflow --proto=A3B15C5D_oP96_62_cd_3c6d_3cd_c --params=`$a, b/a, c/a, x_1, z_1, x_2, z_2, x_3, z_3, x_4, z_4, x_5, z_5, x_6, z_6, x_7, z_7, x_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}$

Other compounds with this structure

- $CoCl_2(NH_3)_5Cl$, $CoI_2(NH_3)_5Cl$, $CrCl_2(NH_3)_5Cl$, $IrCl_2(NH_3)_5Cl$, $OsCl_2(NH_3)_5Cl$, $RhBr_2(NH_3)_5Br$, $RhCl_2(NH_3)_5Cl$, $RhI_2(NH_3)_5Cl$, and $RuCl_2(NH_3)_5Cl$

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(4*c*)	Cl I
$\mathbf{B}_2$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	Cl I
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	(4*c*)	Cl I
$\mathbf{B}_4$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	Cl I
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*c*)	H I
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	(4*c*)	H I
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(4*c*)	H I
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right) c\,\hat{\mathbf{z}}$	(4*c*)	H I
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*c*)	H II
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	(4*c*)	H II
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(4*c*)	H II
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right) c\,\hat{\mathbf{z}}$	(4*c*)	H II
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(4*c*)	H III
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	(4*c*)	H III
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(4*c*)	H III
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\,\hat{\mathbf{z}}$	(4*c*)	H III
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(4*c*)	N I
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	(4*c*)	N I
$\mathbf{B}_{19}$	=	$-x_5\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(4*c*)	N I
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right) c\,\hat{\mathbf{z}}$	(4*c*)	N I
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(4*c*)	N II
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	(4*c*)	N II
$\mathbf{B}_{23}$	=	$-x_6\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(4*c*)	N II
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right) c\,\hat{\mathbf{z}}$	(4*c*)	N II
$\mathbf{B}_{25}$	=	$x_7\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(4*c*)	N III
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} - x_7\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_7\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right) c\,\hat{\mathbf{z}}$	(4*c*)	N III
$\mathbf{B}_{27}$	=	$-x_7\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	(4*c*)	N III
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} + x_7\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_7\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right) c\,\hat{\mathbf{z}}$	(4*c*)	N III
$\mathbf{B}_{29}$	=	$x_8\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(4*c*)	Rh
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2} - x_8\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_8\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right) c\,\hat{\mathbf{z}}$	(4*c*)	Rh
$\mathbf{B}_{31}$	=	$-x_8\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(4*c*)	Rh
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2} + x_8\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_8\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right) c\,\hat{\mathbf{z}}$	(4*c*)	Rh
$\mathbf{B}_{33}$	=	$x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(8*d*)	Cl II
$\mathbf{B}_{34}$	=	$\left(\frac{1}{2} - x_9\right)\mathbf{a}_1 - y_9\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_9\right) a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right) c\,\hat{\mathbf{z}}$	(8*d*)	Cl II
$\mathbf{B}_{35}$	=	$-x_9\,\mathbf{a}_1 + \left(\frac{1}{2} + y_9\right)\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_9\right) b\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}}$	(8*d*)	Cl II

$$\mathbf{B}_{36} = \left(\tfrac{1}{2}+x_9\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_9\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_9\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_9\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_9\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{Cl II}$$

$$\mathbf{B}_{37} = -x_9\,\mathbf{a}_1-y_9\,\mathbf{a}_2-z_9\,\mathbf{a}_3 = -x_9a\,\hat{\mathbf{x}}-y_9b\,\hat{\mathbf{y}}-z_9c\,\hat{\mathbf{z}} \quad (8d) \quad \text{Cl II}$$

$$\mathbf{B}_{38} = \left(\tfrac{1}{2}+x_9\right)\mathbf{a}_1+y_9\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_9\right)a\,\hat{\mathbf{x}}+y_9b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_9\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{Cl II}$$

$$\mathbf{B}_{39} = x_9\,\mathbf{a}_1+\left(\tfrac{1}{2}-y_9\right)\mathbf{a}_2+z_9\,\mathbf{a}_3 = x_9a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_9\right)b\,\hat{\mathbf{y}}+z_9c\,\hat{\mathbf{z}} \quad (8d) \quad \text{Cl II}$$

$$\mathbf{B}_{40} = \left(\tfrac{1}{2}-x_9\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_9\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_9\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_9\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_9\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{Cl II}$$

$$\mathbf{B}_{41} = x_{10}\,\mathbf{a}_1+y_{10}\,\mathbf{a}_2+z_{10}\,\mathbf{a}_3 = x_{10}a\,\hat{\mathbf{x}}+y_{10}b\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H IV}$$

$$\mathbf{B}_{42} = \left(\tfrac{1}{2}-x_{10}\right)\mathbf{a}_1-y_{10}\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{10}\right)a\,\hat{\mathbf{x}}-y_{10}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{10}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H IV}$$

$$\mathbf{B}_{43} = -x_{10}\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_{10}\right)\mathbf{a}_2-z_{10}\,\mathbf{a}_3 = -x_{10}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{10}\right)b\,\hat{\mathbf{y}}-z_{10}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H IV}$$

$$\mathbf{B}_{44} = \left(\tfrac{1}{2}+x_{10}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{10}\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{10}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{10}\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{10}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H IV}$$

$$\mathbf{B}_{45} = -x_{10}\,\mathbf{a}_1-y_{10}\,\mathbf{a}_2-z_{10}\,\mathbf{a}_3 = -x_{10}a\,\hat{\mathbf{x}}-y_{10}b\,\hat{\mathbf{y}}-z_{10}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H IV}$$

$$\mathbf{B}_{46} = \left(\tfrac{1}{2}+x_{10}\right)\mathbf{a}_1+y_{10}\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{10}\right)a\,\hat{\mathbf{x}}+y_{10}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{10}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H IV}$$

$$\mathbf{B}_{47} = x_{10}\,\mathbf{a}_1+\left(\tfrac{1}{2}-y_{10}\right)\mathbf{a}_2+z_{10}\,\mathbf{a}_3 = x_{10}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{10}\right)b\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H IV}$$

$$\mathbf{B}_{48} = \left(\tfrac{1}{2}-x_{10}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{10}\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{10}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{10}\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{10}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H IV}$$

$$\mathbf{B}_{49} = x_{11}\,\mathbf{a}_1+y_{11}\,\mathbf{a}_2+z_{11}\,\mathbf{a}_3 = x_{11}a\,\hat{\mathbf{x}}+y_{11}b\,\hat{\mathbf{y}}+z_{11}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H V}$$

$$\mathbf{B}_{50} = \left(\tfrac{1}{2}-x_{11}\right)\mathbf{a}_1-y_{11}\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{11}\right)a\,\hat{\mathbf{x}}-y_{11}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{11}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H V}$$

$$\mathbf{B}_{51} = -x_{11}\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_{11}\right)\mathbf{a}_2-z_{11}\,\mathbf{a}_3 = -x_{11}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{11}\right)b\,\hat{\mathbf{y}}-z_{11}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H V}$$

$$\mathbf{B}_{52} = \left(\tfrac{1}{2}+x_{11}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{11}\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{11}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{11}\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{11}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H V}$$

$$\mathbf{B}_{53} = -x_{11}\,\mathbf{a}_1-y_{11}\,\mathbf{a}_2-z_{11}\,\mathbf{a}_3 = -x_{11}a\,\hat{\mathbf{x}}-y_{11}b\,\hat{\mathbf{y}}-z_{11}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H V}$$

$$\mathbf{B}_{54} = \left(\tfrac{1}{2}+x_{11}\right)\mathbf{a}_1+y_{11}\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{11}\right)a\,\hat{\mathbf{x}}+y_{11}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{11}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H V}$$

$$\mathbf{B}_{55} = x_{11}\,\mathbf{a}_1+\left(\tfrac{1}{2}-y_{11}\right)\mathbf{a}_2+z_{11}\,\mathbf{a}_3 = x_{11}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{11}\right)b\,\hat{\mathbf{y}}+z_{11}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H V}$$

$$\mathbf{B}_{56} = \left(\tfrac{1}{2}-x_{11}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{11}\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{11}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{11}\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{11}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H V}$$

$$\mathbf{B}_{57} = x_{12}\,\mathbf{a}_1+y_{12}\,\mathbf{a}_2+z_{12}\,\mathbf{a}_3 = x_{12}a\,\hat{\mathbf{x}}+y_{12}b\,\hat{\mathbf{y}}+z_{12}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H VI}$$

$$\mathbf{B}_{58} = \left(\tfrac{1}{2}-x_{12}\right)\mathbf{a}_1-y_{12}\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{12}\right)a\,\hat{\mathbf{x}}-y_{12}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{12}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H VI}$$

$$\mathbf{B}_{59} = -x_{12}\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_{12}\right)\mathbf{a}_2-z_{12}\,\mathbf{a}_3 = -x_{12}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{12}\right)b\,\hat{\mathbf{y}}-z_{12}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H VI}$$

$$\mathbf{B}_{60} = \left(\tfrac{1}{2}+x_{12}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{12}\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{12}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{12}\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{12}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H VI}$$

$$\mathbf{B}_{61} = -x_{12}\,\mathbf{a}_1-y_{12}\,\mathbf{a}_2-z_{12}\,\mathbf{a}_3 = -x_{12}a\,\hat{\mathbf{x}}-y_{12}b\,\hat{\mathbf{y}}-z_{12}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H VI}$$

$$\mathbf{B}_{62} = \left(\tfrac{1}{2}+x_{12}\right)\mathbf{a}_1+y_{12}\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{12}\right)a\,\hat{\mathbf{x}}+y_{12}b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{12}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H VI}$$

$$\mathbf{B}_{63} = x_{12}\,\mathbf{a}_1+\left(\tfrac{1}{2}-y_{12}\right)\mathbf{a}_2+z_{12}\,\mathbf{a}_3 = x_{12}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{12}\right)b\,\hat{\mathbf{y}}+z_{12}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H VI}$$

$$\mathbf{B}_{64} = \left(\tfrac{1}{2}-x_{12}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{12}\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{12}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{12}\right)b\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{12}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H VI}$$

$$\mathbf{B}_{65} = x_{13}\,\mathbf{a}_1+y_{13}\,\mathbf{a}_2+z_{13}\,\mathbf{a}_3 = x_{13}a\,\hat{\mathbf{x}}+y_{13}b\,\hat{\mathbf{y}}+z_{13}c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H VII}$$

$\mathbf{B}_{66}$	=	$\left(\frac{1}{2}-x_{13}\right)\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 + \left(\frac{1}{2}+z_{13}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{13}\right)a\,\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_{13}\right)c\,\hat{\mathbf{z}}$	(8*d*)	H VII
$\mathbf{B}_{67}$	=	$-x_{13}\,\mathbf{a}_1 + \left(\frac{1}{2}+y_{13}\right)\mathbf{a}_2 - z_{13}\,\mathbf{a}_3$	=	$-x_{13}a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_{13}\right)b\,\hat{\mathbf{y}} - z_{13}c\,\hat{\mathbf{z}}$	(8*d*)	H VII
$\mathbf{B}_{68}$	=	$\left(\frac{1}{2}+x_{13}\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_{13}\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_{13}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{13}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_{13}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_{13}\right)c\,\hat{\mathbf{z}}$	(8*d*)	H VII
$\mathbf{B}_{69}$	=	$-x_{13}\,\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 - z_{13}\,\mathbf{a}_3$	=	$-x_{13}a\,\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} - z_{13}c\,\hat{\mathbf{z}}$	(8*d*)	H VII
$\mathbf{B}_{70}$	=	$\left(\frac{1}{2}+x_{13}\right)\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + \left(\frac{1}{2}-z_{13}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{13}\right)a\,\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_{13}\right)c\,\hat{\mathbf{z}}$	(8*d*)	H VII
$\mathbf{B}_{71}$	=	$x_{13}\,\mathbf{a}_1 + \left(\frac{1}{2}-y_{13}\right)\mathbf{a}_2 + z_{13}\,\mathbf{a}_3$	=	$x_{13}a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_{13}\right)b\,\hat{\mathbf{y}} + z_{13}c\,\hat{\mathbf{z}}$	(8*d*)	H VII
$\mathbf{B}_{72}$	=	$\left(\frac{1}{2}-x_{13}\right)\mathbf{a}_1 + \left(\frac{1}{2}+y_{13}\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_{13}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{13}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_{13}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_{13}\right)c\,\hat{\mathbf{z}}$	(8*d*)	H VII
$\mathbf{B}_{73}$	=	$x_{14}\,\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3$	=	$x_{14}a\,\hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} + z_{14}c\,\hat{\mathbf{z}}$	(8*d*)	H VIII
$\mathbf{B}_{74}$	=	$\left(\frac{1}{2}-x_{14}\right)\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 + \left(\frac{1}{2}+z_{14}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{14}\right)a\,\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_{14}\right)c\,\hat{\mathbf{z}}$	(8*d*)	H VIII
$\mathbf{B}_{75}$	=	$-x_{14}\,\mathbf{a}_1 + \left(\frac{1}{2}+y_{14}\right)\mathbf{a}_2 - z_{14}\,\mathbf{a}_3$	=	$-x_{14}a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_{14}\right)b\,\hat{\mathbf{y}} - z_{14}c\,\hat{\mathbf{z}}$	(8*d*)	H VIII
$\mathbf{B}_{76}$	=	$\left(\frac{1}{2}+x_{14}\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_{14}\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_{14}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{14}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_{14}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_{14}\right)c\,\hat{\mathbf{z}}$	(8*d*)	H VIII
$\mathbf{B}_{77}$	=	$-x_{14}\,\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 - z_{14}\,\mathbf{a}_3$	=	$-x_{14}a\,\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} - z_{14}c\,\hat{\mathbf{z}}$	(8*d*)	H VIII
$\mathbf{B}_{78}$	=	$\left(\frac{1}{2}+x_{14}\right)\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + \left(\frac{1}{2}-z_{14}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{14}\right)a\,\hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_{14}\right)c\,\hat{\mathbf{z}}$	(8*d*)	H VIII
$\mathbf{B}_{79}$	=	$x_{14}\,\mathbf{a}_1 + \left(\frac{1}{2}-y_{14}\right)\mathbf{a}_2 + z_{14}\,\mathbf{a}_3$	=	$x_{14}a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_{14}\right)b\,\hat{\mathbf{y}} + z_{14}c\,\hat{\mathbf{z}}$	(8*d*)	H VIII
$\mathbf{B}_{80}$	=	$\left(\frac{1}{2}-x_{14}\right)\mathbf{a}_1 + \left(\frac{1}{2}+y_{14}\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_{14}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{14}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_{14}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_{14}\right)c\,\hat{\mathbf{z}}$	(8*d*)	H VIII
$\mathbf{B}_{81}$	=	$x_{15}\,\mathbf{a}_1 + y_{15}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3$	=	$x_{15}a\,\hat{\mathbf{x}} + y_{15}b\,\hat{\mathbf{y}} + z_{15}c\,\hat{\mathbf{z}}$	(8*d*)	H IX
$\mathbf{B}_{82}$	=	$\left(\frac{1}{2}-x_{15}\right)\mathbf{a}_1 - y_{15}\,\mathbf{a}_2 + \left(\frac{1}{2}+z_{15}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{15}\right)a\,\hat{\mathbf{x}} - y_{15}b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_{15}\right)c\,\hat{\mathbf{z}}$	(8*d*)	H IX
$\mathbf{B}_{83}$	=	$-x_{15}\,\mathbf{a}_1 + \left(\frac{1}{2}+y_{15}\right)\mathbf{a}_2 - z_{15}\,\mathbf{a}_3$	=	$-x_{15}a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_{15}\right)b\,\hat{\mathbf{y}} - z_{15}c\,\hat{\mathbf{z}}$	(8*d*)	H IX
$\mathbf{B}_{84}$	=	$\left(\frac{1}{2}+x_{15}\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_{15}\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_{15}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{15}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_{15}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_{15}\right)c\,\hat{\mathbf{z}}$	(8*d*)	H IX
$\mathbf{B}_{85}$	=	$-x_{15}\,\mathbf{a}_1 - y_{15}\,\mathbf{a}_2 - z_{15}\,\mathbf{a}_3$	=	$-x_{15}a\,\hat{\mathbf{x}} - y_{15}b\,\hat{\mathbf{y}} - z_{15}c\,\hat{\mathbf{z}}$	(8*d*)	H IX
$\mathbf{B}_{86}$	=	$\left(\frac{1}{2}+x_{15}\right)\mathbf{a}_1 + y_{15}\,\mathbf{a}_2 + \left(\frac{1}{2}-z_{15}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{15}\right)a\,\hat{\mathbf{x}} + y_{15}b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_{15}\right)c\,\hat{\mathbf{z}}$	(8*d*)	H IX
$\mathbf{B}_{87}$	=	$x_{15}\,\mathbf{a}_1 + \left(\frac{1}{2}-y_{15}\right)\mathbf{a}_2 + z_{15}\,\mathbf{a}_3$	=	$x_{15}a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_{15}\right)b\,\hat{\mathbf{y}} + z_{15}c\,\hat{\mathbf{z}}$	(8*d*)	H IX
$\mathbf{B}_{88}$	=	$\left(\frac{1}{2}-x_{15}\right)\mathbf{a}_1 + \left(\frac{1}{2}+y_{15}\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_{15}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{15}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_{15}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_{15}\right)c\,\hat{\mathbf{z}}$	(8*d*)	H IX
$\mathbf{B}_{89}$	=	$x_{16}\,\mathbf{a}_1 + y_{16}\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3$	=	$x_{16}a\,\hat{\mathbf{x}} + y_{16}b\,\hat{\mathbf{y}} + z_{16}c\,\hat{\mathbf{z}}$	(8*d*)	N IV
$\mathbf{B}_{90}$	=	$\left(\frac{1}{2}-x_{16}\right)\mathbf{a}_1 - y_{16}\,\mathbf{a}_2 + \left(\frac{1}{2}+z_{16}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{16}\right)a\,\hat{\mathbf{x}} - y_{16}b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_{16}\right)c\,\hat{\mathbf{z}}$	(8*d*)	N IV
$\mathbf{B}_{91}$	=	$-x_{16}\,\mathbf{a}_1 + \left(\frac{1}{2}+y_{16}\right)\mathbf{a}_2 - z_{16}\,\mathbf{a}_3$	=	$-x_{16}a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_{16}\right)b\,\hat{\mathbf{y}} - z_{16}c\,\hat{\mathbf{z}}$	(8*d*)	N IV
$\mathbf{B}_{92}$	=	$\left(\frac{1}{2}+x_{16}\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_{16}\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_{16}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{16}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_{16}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_{16}\right)c\,\hat{\mathbf{z}}$	(8*d*)	N IV
$\mathbf{B}_{93}$	=	$-x_{16}\,\mathbf{a}_1 - y_{16}\,\mathbf{a}_2 - z_{16}\,\mathbf{a}_3$	=	$-x_{16}a\,\hat{\mathbf{x}} - y_{16}b\,\hat{\mathbf{y}} - z_{16}c\,\hat{\mathbf{z}}$	(8*d*)	N IV
$\mathbf{B}_{94}$	=	$\left(\frac{1}{2}+x_{16}\right)\mathbf{a}_1 + y_{16}\,\mathbf{a}_2 + \left(\frac{1}{2}-z_{16}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{16}\right)a\,\hat{\mathbf{x}} + y_{16}b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_{16}\right)c\,\hat{\mathbf{z}}$	(8*d*)	N IV
$\mathbf{B}_{95}$	=	$x_{16}\,\mathbf{a}_1 + \left(\frac{1}{2}-y_{16}\right)\mathbf{a}_2 + z_{16}\,\mathbf{a}_3$	=	$x_{16}a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_{16}\right)b\,\hat{\mathbf{y}} + z_{16}c\,\hat{\mathbf{z}}$	(8*d*)	N IV

$$\mathbf{B}_{96} = \left(\tfrac{1}{2} - x_{16}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{16}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{16}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{16}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{16}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{16}\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{N IV}$$

References:

- R. S. Evans, E. A. Hopcus, J. Bordner, and A. F. Schreiner, *Molecular and crystal structures of halopentaamminerhodium-(III) complexes, [Rh(NH$_3$)$_5$Cl]Cl$_2$ and [Rh(NH$_3$)$_5$Br]Br$_2$*, J. Cryst. Mol. Struct. **3**, 235–245 (1973), doi:10.1007/BF01236644.

Found in:

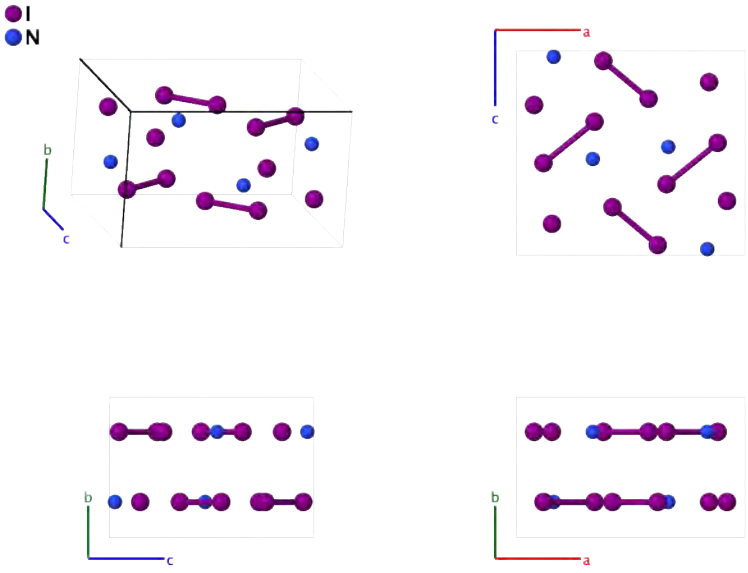
- T. W. Hambley and P. A. Lay, *Comparisons of π-bonding and hydrogen bonding in isomorphous compounds: [M(NH$_3$)$_5$Cl]Cl$_2$ (M=Cr, Co, Rh, Ir, Ru, Os)*, Inorg. Chem. **25**, 4553–4558 (1986), doi:10.1021/ic00245a020.

Geometry files:

- CIF: pp. 1634
- POSCAR: pp. 1635

NH_4I_3 ($D0_{16}$) Structure: A3B_oP16_62_3c_c

http://aflow.org/prototype-encyclopedia/A3B_oP16_62_3c_c



Prototype	:	$I_3(NH_4)$
AFLOW prototype label	:	A3B_oP16_62_3c_c
***Strukturbericht* designation**	:	$D0_{16}$
Pearson symbol	:	oP16
Space group number	:	62
Space group symbol	:	*Pnma*
AFLOW prototype command	:	`aflow --proto=A3B_oP16_62_3c_c --params=`$a, b/a, c/a, x_1, z_1, x_2, z_2, x_3, z_3, x_4, z_4$

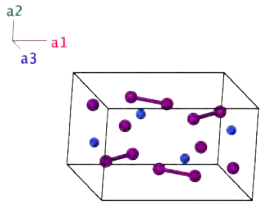
Other compounds with this structure

- CsI_3 and $CsBr_3$

- (Cheesman, 1970) gives the positions of the atoms in Cartesian coordinates, which we have translated into the standard Wyckoff notation.
- The hydrogen positions are not given, so the NH_4 molecules are centered on the ($4c$) Wyckoff position.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(4*c*)	I I
$\mathbf{B}_2$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	I I
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	(4*c*)	I I
$\mathbf{B}_4$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	I I
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*c*)	I II
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	(4*c*)	I II
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(4*c*)	I II
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right) c\,\hat{\mathbf{z}}$	(4*c*)	I II
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*c*)	I III
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	(4*c*)	I III
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(4*c*)	I III
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right) c\,\hat{\mathbf{z}}$	(4*c*)	I III
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(4*c*)	NH_4
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	(4*c*)	NH_4
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(4*c*)	NH_4
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\,\hat{\mathbf{z}}$	(4*c*)	NH_4

References:

- G. H. Cheesman and A. J. T. Finney, *Refinement of the structure of ammonium triiodide, NH_4I_3*, Acta Crystallogr. Sect. B Struct. Sci. **26**, 904–906 (1970), doi:10.1107/S0567740870003357.

Found in:

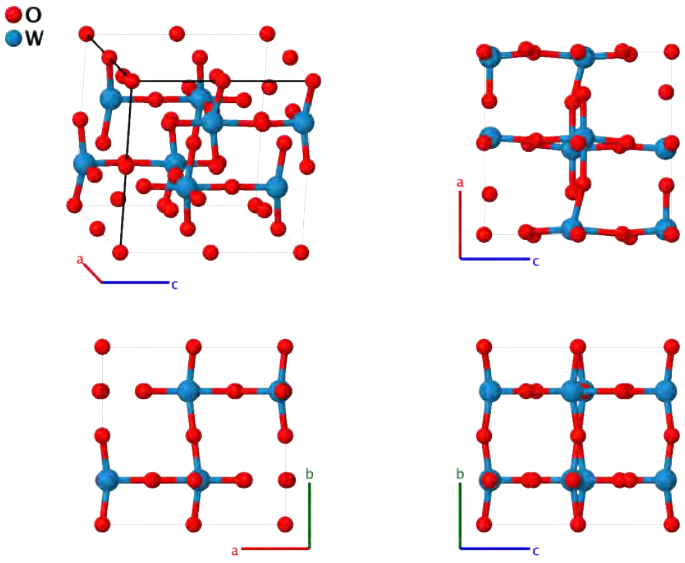
- A. J. T. Finney, *The Structure and Stability of Simple Tri-Iodides*, http://eprints.utas.edu.au/19441/1/whole_FinneyAnthonyJohnThompson1974_thesis.pdf (1973). Ph.D. Thesis, University of Tasmania.

Geometry files:

- CIF: pp. 1635
- POSCAR: pp. 1636

Original β-WO_3 (*obsolete*) Structure: A3B_oP32_62_ab4c_2c

http://aflow.org/prototype-encyclopedia/A3B_oP32_62_ab4c_2c



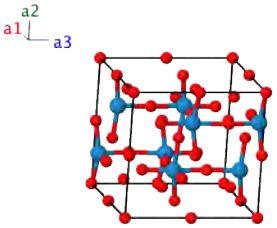
Prototype	:	O_3W
AFLOW prototype label	:	A3B_oP32_62_ab4c_2c
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP32
Space group number	:	62
Space group symbol	:	*Pnma*
AFLOW prototype command	:	`aflow --proto=A3B_oP32_62_ab4c_2c --params=`$a, b/a, c/a, x_3, z_3, x_4, z_4, x_5, z_5, x_6, z_6, x_7, z_7, x_8, z_8$

- All stable phases of WO_3 are distortions of the cubic α-ReO_3 ($D0_9$) phase. Based on (Woodward, 1997 and Vogt, 1999), the known stable phases and their approximate temperature ranges are:
  - α-WO_3 (1010-1170 K) (Vogt, 1999)
  - β-WO_3 (600-1170 K) (Vogt, 1999)
  - γ-WO_3 (290-600 K) (Vogt, 1999)
  - δ-WO_3 (230-290 K) (Diehl, 1978)
  - ϵ-WO_3 (below 23 K) (Woodward, 1997)

  Woodward's caption says "The transition temperatures display large hysteresis effects and universal agreement is not found in the literature."
- In addition, several other structures have been proposed and/or found:
  - The original $D0_{10}$ structure (Bräkken, 1931), (Hermann, 1937) superseded by δ-WO_3
  - Original β-WO_3 (Salje, 1977), this structure
  - Hexagonal WO_3 (Gerand, 1979) (metastable)

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}b\,\hat{\mathbf{y}}$	(4*a*)	O I
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	(4*b*)	O II
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	(4*b*)	O II
$\mathbf{B}_7$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*b*)	O II
$\mathbf{B}_8$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$	(4*b*)	O II
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	(4*c*)	O III
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_3\right)\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(4*c*)	O III
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	(4*c*)	O III
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(4*c*)	O III
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	(4*c*)	O IV
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - x_4\right)\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(4*c*)	O IV
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	(4*c*)	O IV
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_4\right)\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(4*c*)	O IV
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_5c\,\hat{\mathbf{z}}$	(4*c*)	O V
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} - x_5\right)\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	(4*c*)	O V
$\mathbf{B}_{19}$	=	$-x_5\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_5c\,\hat{\mathbf{z}}$	(4*c*)	O V
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} + x_5\right)\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	(4*c*)	O V
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_6c\,\hat{\mathbf{z}}$	(4*c*)	O VI
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} - x_6\right)\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\,\hat{\mathbf{z}}$	(4*c*)	O VI
$\mathbf{B}_{23}$	=	$-x_6\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$-x_6a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_6c\,\hat{\mathbf{z}}$	(4*c*)	O VI
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} + x_6\right)\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\,\hat{\mathbf{z}}$	(4*c*)	O VI
$\mathbf{B}_{25}$	=	$x_7\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$x_7a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_7c\,\hat{\mathbf{z}}$	(4*c*)	W I
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} - x_7\right)\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_7\right)a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right)c\,\hat{\mathbf{z}}$	(4*c*)	W I
$\mathbf{B}_{27}$	=	$-x_7\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$-x_7a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_7c\,\hat{\mathbf{z}}$	(4*c*)	W I

$$\mathbf{B}_{28} = \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2 + \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{x}} + \tfrac{1}{4}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_7\right)c\,\hat{\mathbf{z}} \quad (4c) \quad \text{W I}$$

$$\mathbf{B}_{29} = x_8\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2 + z_8\mathbf{a}_3 = x_8 a\,\hat{\mathbf{x}} + \tfrac{1}{4}b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}} \quad (4c) \quad \text{W II}$$

$$\mathbf{B}_{30} = \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_1 + \tfrac{3}{4}\mathbf{a}_2 + \left(\tfrac{1}{2}+z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{x}} + \tfrac{3}{4}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_8\right)c\,\hat{\mathbf{z}} \quad (4c) \quad \text{W II}$$

$$\mathbf{B}_{31} = -x_8\mathbf{a}_1 + \tfrac{3}{4}\mathbf{a}_2 - z_8\mathbf{a}_3 = -x_8 a\,\hat{\mathbf{x}} + \tfrac{3}{4}b\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}} \quad (4c) \quad \text{W II}$$

$$\mathbf{B}_{32} = \left(\tfrac{1}{2}+x_8\right)\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2 + \left(\tfrac{1}{2}-z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{x}} + \tfrac{1}{4}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_8\right)c\,\hat{\mathbf{z}} \quad (4c) \quad \text{W II}$$

References:

- P. M. Woodward, A. W. Sleight, and T. Vogt, *Ferroelectric Tungsten Trioxide*, J. Solid State Chem. **131**, 9–17 (1997), doi:10.1006/jssc.1997.7268.
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- R. Diehl, G. Brandt, and E. Salje, *The Crystal Structure of Triclinic WO_3*, Acta Crystallogr. Sect. B Struct. Sci. **34**, 1105–1111 (1978), doi:10.1107/S0567740878005014.
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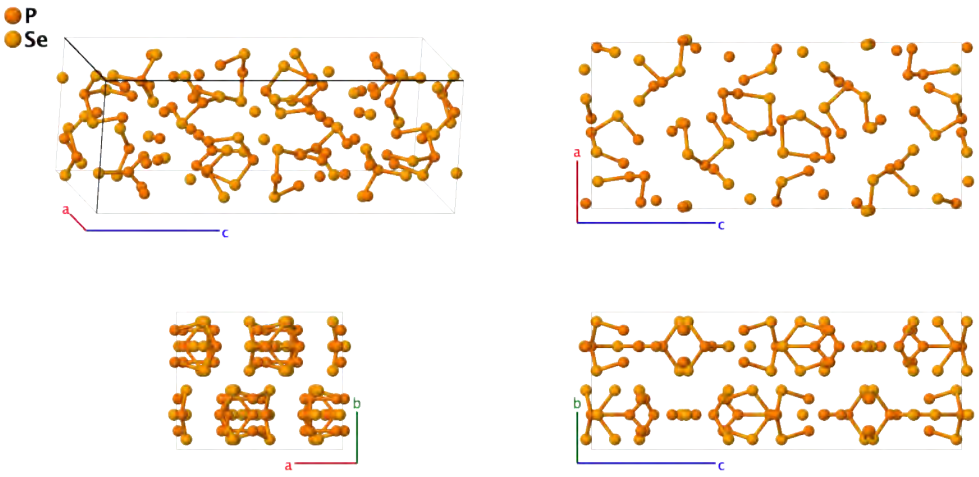
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).
- E. Salje, *The Orthorhombic Phase of WO_3*, Acta Crystallogr. Sect. B Struct. Sci. **33**, 574–577 (1977), doi:10.1107/S0567740877004130.
- B. Gerand, G. Nowogrocki, J. Guenot, and M. Figlarz, *Structural study of a new hexagonal form of tungsten trioxide*, J. Solid State Chem. **29**, 429–434 (1979), doi:10.1016/0022-4596(79)90199-3.

Geometry files:

- CIF: pp. 1636
- POSCAR: pp. 1636

P_4Se_3 Structure: A4B3_oP112_62_8c4d_4c4d

http://aflow.org/prototype-encyclopedia/A4B3_oP112_62_8c4d_4c4d

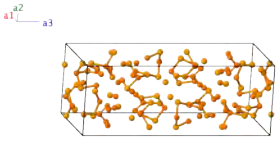


Prototype	:	P_4Se_3
AFLOW prototype label	:	A4B3_oP112_62_8c4d_4c4d
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP112
Space group number	:	62
Space group symbol	:	*Pnma*
AFLOW prototype command	:	`aflow --proto=A4B3_oP112_62_8c4d_4c4d --params=`$a, b/a, c/a, x_1, z_1, x_2, z_2, x_3, z_3, x_4, z_4, x_5, z_5, x_6, z_6, x_7, z_7, x_8, z_8, x_9, z_9, x_{10}, z_{10}, x_{11}, z_{11}, x_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}, x_{17}, y_{17}, z_{17}, x_{18}, y_{18}, z_{18}, x_{19}, y_{19}, z_{19}, x_{20}, y_{20}, z_{20}$

- (Keulen, 1959) give this structure in the *Pmnb* setting of space group #62. We used FINDSYM to translate this into the standard *Pnma* setting.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	($4c$)	P I
$\mathbf{B}_2$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	($4c$)	P I
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	($4c$)	P I
$\mathbf{B}_4$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right) c\,\hat{\mathbf{z}}$	($4c$)	P I

$$
\begin{array}{lcccccc}
\mathbf{B}_{5} & = & x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3 & = & x_2 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}} & (4c) & \text{P II} \\
\mathbf{B}_{6} & = & \left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}} & (4c) & \text{P II} \\
\mathbf{B}_{7} & = & -x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3 & = & -x_2 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}} & (4c) & \text{P II} \\
\mathbf{B}_{8} & = & \left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3 & = & \left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right) c\,\hat{\mathbf{z}} & (4c) & \text{P II} \\
\mathbf{B}_{9} & = & x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3 & = & x_3 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}} & (4c) & \text{P III} \\
\mathbf{B}_{10} & = & \left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_3\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}} & (4c) & \text{P III} \\
\mathbf{B}_{11} & = & -x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3 & = & -x_3 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}} & (4c) & \text{P III} \\
\mathbf{B}_{12} & = & \left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3 & = & \left(\frac{1}{2} + x_3\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right) c\,\hat{\mathbf{z}} & (4c) & \text{P III} \\
\mathbf{B}_{13} & = & x_4\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3 & = & x_4 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}} & (4c) & \text{P IV} \\
\mathbf{B}_{14} & = & \left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}} & (4c) & \text{P IV} \\
\mathbf{B}_{15} & = & -x_4\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3 & = & -x_4 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}} & (4c) & \text{P IV} \\
\mathbf{B}_{16} & = & \left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3 & = & \left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\,\hat{\mathbf{z}} & (4c) & \text{P IV} \\
\mathbf{B}_{17} & = & x_5\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_5\,\mathbf{a}_3 & = & x_5 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}} & (4c) & \text{P V} \\
\mathbf{B}_{18} & = & \left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_5\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}} & (4c) & \text{P V} \\
\mathbf{B}_{19} & = & -x_5\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_5\,\mathbf{a}_3 & = & -x_5 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}} & (4c) & \text{P V} \\
\mathbf{B}_{20} & = & \left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3 & = & \left(\frac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right) c\,\hat{\mathbf{z}} & (4c) & \text{P V} \\
\mathbf{B}_{21} & = & x_6\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 & = & x_6 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}} & (4c) & \text{P VI} \\
\mathbf{B}_{22} & = & \left(\frac{1}{2} - x_6\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}} & (4c) & \text{P VI} \\
\mathbf{B}_{23} & = & -x_6\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_6\,\mathbf{a}_3 & = & -x_6 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}} & (4c) & \text{P VI} \\
\mathbf{B}_{24} & = & \left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3 & = & \left(\frac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right) c\,\hat{\mathbf{z}} & (4c) & \text{P VI} \\
\mathbf{B}_{25} & = & x_7\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 & = & x_7 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}} & (4c) & \text{P VII} \\
\mathbf{B}_{26} & = & \left(\frac{1}{2} - x_7\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_7\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right) c\,\hat{\mathbf{z}} & (4c) & \text{P VII} \\
\mathbf{B}_{27} & = & -x_7\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_7\,\mathbf{a}_3 & = & -x_7 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}} & (4c) & \text{P VII} \\
\mathbf{B}_{28} & = & \left(\frac{1}{2} + x_7\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\mathbf{a}_3 & = & \left(\frac{1}{2} + x_7\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right) c\,\hat{\mathbf{z}} & (4c) & \text{P VII} \\
\mathbf{B}_{29} & = & x_8\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 & = & x_8 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}} & (4c) & \text{P VIII} \\
\mathbf{B}_{30} & = & \left(\frac{1}{2} - x_8\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_8\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right) c\,\hat{\mathbf{z}} & (4c) & \text{P VIII} \\
\mathbf{B}_{31} & = & -x_8\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_8\,\mathbf{a}_3 & = & -x_8 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}} & (4c) & \text{P VIII} \\
\mathbf{B}_{32} & = & \left(\frac{1}{2} + x_8\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\mathbf{a}_3 & = & \left(\frac{1}{2} + x_8\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right) c\,\hat{\mathbf{z}} & (4c) & \text{P VIII} \\
\mathbf{B}_{33} & = & x_9\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 & = & x_9 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}} & (4c) & \text{Se I} \\
\mathbf{B}_{34} & = & \left(\frac{1}{2} - x_9\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_9\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right) c\,\hat{\mathbf{z}} & (4c) & \text{Se I} \\
\mathbf{B}_{35} & = & -x_9\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_9\,\mathbf{a}_3 & = & -x_9 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}} & (4c) & \text{Se I} \\
\mathbf{B}_{36} & = & \left(\frac{1}{2} + x_9\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_9\right)\mathbf{a}_3 & = & \left(\frac{1}{2} + x_9\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_9\right) c\,\hat{\mathbf{z}} & (4c) & \text{Se I} \\
\mathbf{B}_{37} & = & x_{10}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 & = & x_{10} a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}} & (4c) & \text{Se II} \\
\mathbf{B}_{38} & = & \left(\frac{1}{2} - x_{10}\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{10}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_{10}\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{10}\right) c\,\hat{\mathbf{z}} & (4c) & \text{Se II} \\
\mathbf{B}_{39} & = & -x_{10}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 & = & -x_{10} a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_{10} c\,\hat{\mathbf{z}} & (4c) & \text{Se II} \\
\mathbf{B}_{40} & = & \left(\frac{1}{2} + x_{10}\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{10}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} + x_{10}\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{10}\right) c\,\hat{\mathbf{z}} & (4c) & \text{Se II}
\end{array}
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$\mathbf{B}_{41}$	=	$x_{11}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$x_{11}a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_{11}c\,\hat{\mathbf{z}}$	(4*c*)	Se III
$\mathbf{B}_{42}$	=	$\left(\frac{1}{2} - x_{11}\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{11}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{11}\right)a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{11}\right)c\,\hat{\mathbf{z}}$	(4*c*)	Se III
$\mathbf{B}_{43}$	=	$-x_{11}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3$	=	$-x_{11}a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_{11}c\,\hat{\mathbf{z}}$	(4*c*)	Se III
$\mathbf{B}_{44}$	=	$\left(\frac{1}{2} + x_{11}\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{11}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{11}\right)a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{11}\right)c\,\hat{\mathbf{z}}$	(4*c*)	Se III
$\mathbf{B}_{45}$	=	$x_{12}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3$	=	$x_{12}a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_{12}c\,\hat{\mathbf{z}}$	(4*c*)	Se IV
$\mathbf{B}_{46}$	=	$\left(\frac{1}{2} - x_{12}\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{12}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{12}\right)a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{12}\right)c\,\hat{\mathbf{z}}$	(4*c*)	Se IV
$\mathbf{B}_{47}$	=	$-x_{12}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3$	=	$-x_{12}a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_{12}c\,\hat{\mathbf{z}}$	(4*c*)	Se IV
$\mathbf{B}_{48}$	=	$\left(\frac{1}{2} + x_{12}\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{12}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{12}\right)a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{12}\right)c\,\hat{\mathbf{z}}$	(4*c*)	Se IV
$\mathbf{B}_{49}$	=	$x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3$	=	$x_{13}a\,\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} + z_{13}c\,\hat{\mathbf{z}}$	(8*d*)	P IX
$\mathbf{B}_{50}$	=	$\left(\frac{1}{2} - x_{13}\right)\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{13}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{13}\right)a\,\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{13}\right)c\,\hat{\mathbf{z}}$	(8*d*)	P IX
$\mathbf{B}_{51}$	=	$-x_{13}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{13}\right)\mathbf{a}_2 - z_{13}\,\mathbf{a}_3$	=	$-x_{13}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{13}\right)b\,\hat{\mathbf{y}} - z_{13}c\,\hat{\mathbf{z}}$	(8*d*)	P IX
$\mathbf{B}_{52}$	=	$\left(\frac{1}{2} + x_{13}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{13}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{13}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{13}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{13}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{13}\right)c\,\hat{\mathbf{z}}$	(8*d*)	P IX
$\mathbf{B}_{53}$	=	$-x_{13}\,\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 - z_{13}\,\mathbf{a}_3$	=	$-x_{13}a\,\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} - z_{13}c\,\hat{\mathbf{z}}$	(8*d*)	P IX
$\mathbf{B}_{54}$	=	$\left(\frac{1}{2} + x_{13}\right)\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{13}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{13}\right)a\,\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{13}\right)c\,\hat{\mathbf{z}}$	(8*d*)	P IX
$\mathbf{B}_{55}$	=	$x_{13}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{13}\right)\mathbf{a}_2 + z_{13}\,\mathbf{a}_3$	=	$x_{13}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{13}\right)b\,\hat{\mathbf{y}} + z_{13}c\,\hat{\mathbf{z}}$	(8*d*)	P IX
$\mathbf{B}_{56}$	=	$\left(\frac{1}{2} - x_{13}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{13}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{13}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{13}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{13}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{13}\right)c\,\hat{\mathbf{z}}$	(8*d*)	P IX
$\mathbf{B}_{57}$	=	$x_{14}\,\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3$	=	$x_{14}a\,\hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} + z_{14}c\,\hat{\mathbf{z}}$	(8*d*)	P X
$\mathbf{B}_{58}$	=	$\left(\frac{1}{2} - x_{14}\right)\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{14}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{14}\right)a\,\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{14}\right)c\,\hat{\mathbf{z}}$	(8*d*)	P X
$\mathbf{B}_{59}$	=	$-x_{14}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{14}\right)\mathbf{a}_2 - z_{14}\,\mathbf{a}_3$	=	$-x_{14}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{14}\right)b\,\hat{\mathbf{y}} - z_{14}c\,\hat{\mathbf{z}}$	(8*d*)	P X
$\mathbf{B}_{60}$	=	$\left(\frac{1}{2} + x_{14}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{14}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{14}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{14}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{14}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{14}\right)c\,\hat{\mathbf{z}}$	(8*d*)	P X
$\mathbf{B}_{61}$	=	$-x_{14}\,\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 - z_{14}\,\mathbf{a}_3$	=	$-x_{14}a\,\hat{\mathbf{x}} - y_{14}b\,\hat{\mathbf{y}} - z_{14}c\,\hat{\mathbf{z}}$	(8*d*)	P X
$\mathbf{B}_{62}$	=	$\left(\frac{1}{2} + x_{14}\right)\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{14}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{14}\right)a\,\hat{\mathbf{x}} + y_{14}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{14}\right)c\,\hat{\mathbf{z}}$	(8*d*)	P X
$\mathbf{B}_{63}$	=	$x_{14}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{14}\right)\mathbf{a}_2 + z_{14}\,\mathbf{a}_3$	=	$x_{14}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{14}\right)b\,\hat{\mathbf{y}} + z_{14}c\,\hat{\mathbf{z}}$	(8*d*)	P X
$\mathbf{B}_{64}$	=	$\left(\frac{1}{2} - x_{14}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{14}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{14}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{14}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{14}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{14}\right)c\,\hat{\mathbf{z}}$	(8*d*)	P X
$\mathbf{B}_{65}$	=	$x_{15}\,\mathbf{a}_1 + y_{15}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3$	=	$x_{15}a\,\hat{\mathbf{x}} + y_{15}b\,\hat{\mathbf{y}} + z_{15}c\,\hat{\mathbf{z}}$	(8*d*)	P XI
$\mathbf{B}_{66}$	=	$\left(\frac{1}{2} - x_{15}\right)\mathbf{a}_1 - y_{15}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{15}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{15}\right)a\,\hat{\mathbf{x}} - y_{15}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{15}\right)c\,\hat{\mathbf{z}}$	(8*d*)	P XI
$\mathbf{B}_{67}$	=	$-x_{15}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{15}\right)\mathbf{a}_2 - z_{15}\,\mathbf{a}_3$	=	$-x_{15}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{15}\right)b\,\hat{\mathbf{y}} - z_{15}c\,\hat{\mathbf{z}}$	(8*d*)	P XI
$\mathbf{B}_{68}$	=	$\left(\frac{1}{2} + x_{15}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{15}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{15}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{15}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{15}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{15}\right)c\,\hat{\mathbf{z}}$	(8*d*)	P XI
$\mathbf{B}_{69}$	=	$-x_{15}\,\mathbf{a}_1 - y_{15}\,\mathbf{a}_2 - z_{15}\,\mathbf{a}_3$	=	$-x_{15}a\,\hat{\mathbf{x}} - y_{15}b\,\hat{\mathbf{y}} - z_{15}c\,\hat{\mathbf{z}}$	(8*d*)	P XI
$\mathbf{B}_{70}$	=	$\left(\frac{1}{2} + x_{15}\right)\mathbf{a}_1 + y_{15}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{15}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{15}\right)a\,\hat{\mathbf{x}} + y_{15}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{15}\right)c\,\hat{\mathbf{z}}$	(8*d*)	P XI
$\mathbf{B}_{71}$	=	$x_{15}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{15}\right)\mathbf{a}_2 + z_{15}\,\mathbf{a}_3$	=	$x_{15}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{15}\right)b\,\hat{\mathbf{y}} + z_{15}c\,\hat{\mathbf{z}}$	(8*d*)	P XI
$\mathbf{B}_{72}$	=	$\left(\frac{1}{2} - x_{15}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{15}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{15}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{15}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{15}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{15}\right)c\,\hat{\mathbf{z}}$	(8*d*)	P XI

$\mathbf{B}_{73}$	$=$	$x_{16}\,\mathbf{a}_1 + y_{16}\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3$	$=$	$x_{16}a\,\hat{\mathbf{x}} + y_{16}b\,\hat{\mathbf{y}} + z_{16}c\,\hat{\mathbf{z}}$	(8*d*)	P XII
$\mathbf{B}_{74}$	$=$	$\left(\frac{1}{2} - x_{16}\right)\mathbf{a}_1 - y_{16}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{16}\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2} - x_{16}\right)a\,\hat{\mathbf{x}} - y_{16}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{16}\right)c\,\hat{\mathbf{z}}$	(8*d*)	P XII
$\mathbf{B}_{75}$	$=$	$-x_{16}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{16}\right)\mathbf{a}_2 - z_{16}\,\mathbf{a}_3$	$=$	$-x_{16}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{16}\right)b\,\hat{\mathbf{y}} - z_{16}c\,\hat{\mathbf{z}}$	(8*d*)	P XII
$\mathbf{B}_{76}$	$=$	$\left(\frac{1}{2} + x_{16}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{16}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{16}\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2} + x_{16}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{16}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{16}\right)c\,\hat{\mathbf{z}}$	(8*d*)	P XII
$\mathbf{B}_{77}$	$=$	$-x_{16}\,\mathbf{a}_1 - y_{16}\,\mathbf{a}_2 - z_{16}\,\mathbf{a}_3$	$=$	$-x_{16}a\,\hat{\mathbf{x}} - y_{16}b\,\hat{\mathbf{y}} - z_{16}c\,\hat{\mathbf{z}}$	(8*d*)	P XII
$\mathbf{B}_{78}$	$=$	$\left(\frac{1}{2} + x_{16}\right)\mathbf{a}_1 + y_{16}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{16}\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2} + x_{16}\right)a\,\hat{\mathbf{x}} + y_{16}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{16}\right)c\,\hat{\mathbf{z}}$	(8*d*)	P XII
$\mathbf{B}_{79}$	$=$	$x_{16}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{16}\right)\mathbf{a}_2 + z_{16}\,\mathbf{a}_3$	$=$	$x_{16}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{16}\right)b\,\hat{\mathbf{y}} + z_{16}c\,\hat{\mathbf{z}}$	(8*d*)	P XII
$\mathbf{B}_{80}$	$=$	$\left(\frac{1}{2} - x_{16}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{16}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{16}\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2} - x_{16}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{16}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{16}\right)c\,\hat{\mathbf{z}}$	(8*d*)	P XII
$\mathbf{B}_{81}$	$=$	$x_{17}\,\mathbf{a}_1 + y_{17}\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3$	$=$	$x_{17}a\,\hat{\mathbf{x}} + y_{17}b\,\hat{\mathbf{y}} + z_{17}c\,\hat{\mathbf{z}}$	(8*d*)	Se V
$\mathbf{B}_{82}$	$=$	$\left(\frac{1}{2} - x_{17}\right)\mathbf{a}_1 - y_{17}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{17}\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2} - x_{17}\right)a\,\hat{\mathbf{x}} - y_{17}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{17}\right)c\,\hat{\mathbf{z}}$	(8*d*)	Se V
$\mathbf{B}_{83}$	$=$	$-x_{17}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{17}\right)\mathbf{a}_2 - z_{17}\,\mathbf{a}_3$	$=$	$-x_{17}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{17}\right)b\,\hat{\mathbf{y}} - z_{17}c\,\hat{\mathbf{z}}$	(8*d*)	Se V
$\mathbf{B}_{84}$	$=$	$\left(\frac{1}{2} + x_{17}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{17}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{17}\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2} + x_{17}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{17}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{17}\right)c\,\hat{\mathbf{z}}$	(8*d*)	Se V
$\mathbf{B}_{85}$	$=$	$-x_{17}\,\mathbf{a}_1 - y_{17}\,\mathbf{a}_2 - z_{17}\,\mathbf{a}_3$	$=$	$-x_{17}a\,\hat{\mathbf{x}} - y_{17}b\,\hat{\mathbf{y}} - z_{17}c\,\hat{\mathbf{z}}$	(8*d*)	Se V
$\mathbf{B}_{86}$	$=$	$\left(\frac{1}{2} + x_{17}\right)\mathbf{a}_1 + y_{17}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{17}\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2} + x_{17}\right)a\,\hat{\mathbf{x}} + y_{17}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{17}\right)c\,\hat{\mathbf{z}}$	(8*d*)	Se V
$\mathbf{B}_{87}$	$=$	$x_{17}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{17}\right)\mathbf{a}_2 + z_{17}\,\mathbf{a}_3$	$=$	$x_{17}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{17}\right)b\,\hat{\mathbf{y}} + z_{17}c\,\hat{\mathbf{z}}$	(8*d*)	Se V
$\mathbf{B}_{88}$	$=$	$\left(\frac{1}{2} - x_{17}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{17}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{17}\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2} - x_{17}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{17}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{17}\right)c\,\hat{\mathbf{z}}$	(8*d*)	Se V
$\mathbf{B}_{89}$	$=$	$x_{18}\,\mathbf{a}_1 + y_{18}\,\mathbf{a}_2 + z_{18}\,\mathbf{a}_3$	$=$	$x_{18}a\,\hat{\mathbf{x}} + y_{18}b\,\hat{\mathbf{y}} + z_{18}c\,\hat{\mathbf{z}}$	(8*d*)	Se VI
$\mathbf{B}_{90}$	$=$	$\left(\frac{1}{2} - x_{18}\right)\mathbf{a}_1 - y_{18}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{18}\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2} - x_{18}\right)a\,\hat{\mathbf{x}} - y_{18}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{18}\right)c\,\hat{\mathbf{z}}$	(8*d*)	Se VI
$\mathbf{B}_{91}$	$=$	$-x_{18}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{18}\right)\mathbf{a}_2 - z_{18}\,\mathbf{a}_3$	$=$	$-x_{18}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{18}\right)b\,\hat{\mathbf{y}} - z_{18}c\,\hat{\mathbf{z}}$	(8*d*)	Se VI
$\mathbf{B}_{92}$	$=$	$\left(\frac{1}{2} + x_{18}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{18}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{18}\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2} + x_{18}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{18}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{18}\right)c\,\hat{\mathbf{z}}$	(8*d*)	Se VI
$\mathbf{B}_{93}$	$=$	$-x_{18}\,\mathbf{a}_1 - y_{18}\,\mathbf{a}_2 - z_{18}\,\mathbf{a}_3$	$=$	$-x_{18}a\,\hat{\mathbf{x}} - y_{18}b\,\hat{\mathbf{y}} - z_{18}c\,\hat{\mathbf{z}}$	(8*d*)	Se VI
$\mathbf{B}_{94}$	$=$	$\left(\frac{1}{2} + x_{18}\right)\mathbf{a}_1 + y_{18}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{18}\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2} + x_{18}\right)a\,\hat{\mathbf{x}} + y_{18}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{18}\right)c\,\hat{\mathbf{z}}$	(8*d*)	Se VI
$\mathbf{B}_{95}$	$=$	$x_{18}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{18}\right)\mathbf{a}_2 + z_{18}\,\mathbf{a}_3$	$=$	$x_{18}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{18}\right)b\,\hat{\mathbf{y}} + z_{18}c\,\hat{\mathbf{z}}$	(8*d*)	Se VI
$\mathbf{B}_{96}$	$=$	$\left(\frac{1}{2} - x_{18}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{18}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{18}\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2} - x_{18}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{18}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{18}\right)c\,\hat{\mathbf{z}}$	(8*d*)	Se VI
$\mathbf{B}_{97}$	$=$	$x_{19}\,\mathbf{a}_1 + y_{19}\,\mathbf{a}_2 + z_{19}\,\mathbf{a}_3$	$=$	$x_{19}a\,\hat{\mathbf{x}} + y_{19}b\,\hat{\mathbf{y}} + z_{19}c\,\hat{\mathbf{z}}$	(8*d*)	Se VII
$\mathbf{B}_{98}$	$=$	$\left(\frac{1}{2} - x_{19}\right)\mathbf{a}_1 - y_{19}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{19}\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2} - x_{19}\right)a\,\hat{\mathbf{x}} - y_{19}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{19}\right)c\,\hat{\mathbf{z}}$	(8*d*)	Se VII
$\mathbf{B}_{99}$	$=$	$-x_{19}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{19}\right)\mathbf{a}_2 - z_{19}\,\mathbf{a}_3$	$=$	$-x_{19}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{19}\right)b\,\hat{\mathbf{y}} - z_{19}c\,\hat{\mathbf{z}}$	(8*d*)	Se VII
$\mathbf{B}_{100}$	$=$	$\left(\frac{1}{2} + x_{19}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{19}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{19}\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2} + x_{19}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{19}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{19}\right)c\,\hat{\mathbf{z}}$	(8*d*)	Se VII
$\mathbf{B}_{101}$	$=$	$-x_{19}\,\mathbf{a}_1 - y_{19}\,\mathbf{a}_2 - z_{19}\,\mathbf{a}_3$	$=$	$-x_{19}a\,\hat{\mathbf{x}} - y_{19}b\,\hat{\mathbf{y}} - z_{19}c\,\hat{\mathbf{z}}$	(8*d*)	Se VII
$\mathbf{B}_{102}$	$=$	$\left(\frac{1}{2} + x_{19}\right)\mathbf{a}_1 + y_{19}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{19}\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2} + x_{19}\right)a\,\hat{\mathbf{x}} + y_{19}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{19}\right)c\,\hat{\mathbf{z}}$	(8*d*)	Se VII
$\mathbf{B}_{103}$	$=$	$x_{19}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{19}\right)\mathbf{a}_2 + z_{19}\,\mathbf{a}_3$	$=$	$x_{19}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{19}\right)b\,\hat{\mathbf{y}} + z_{19}c\,\hat{\mathbf{z}}$	(8*d*)	Se VII

$\mathbf{B}_{104}$	=	$\left(\frac{1}{2}-x_{19}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{19}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{19}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{19}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{19}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{19}\right)c\,\hat{\mathbf{z}}$	(8*d*)	Se VII
$\mathbf{B}_{105}$	=	$x_{20}\,\mathbf{a}_1+y_{20}\,\mathbf{a}_2+z_{20}\,\mathbf{a}_3$	=	$x_{20}a\,\hat{\mathbf{x}}+y_{20}b\,\hat{\mathbf{y}}+z_{20}c\,\hat{\mathbf{z}}$	(8*d*)	Se VIII
$\mathbf{B}_{106}$	=	$\left(\frac{1}{2}-x_{20}\right)\mathbf{a}_1-y_{20}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{20}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{20}\right)a\,\hat{\mathbf{x}}-y_{20}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{20}\right)c\,\hat{\mathbf{z}}$	(8*d*)	Se VIII
$\mathbf{B}_{107}$	=	$-x_{20}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{20}\right)\mathbf{a}_2-z_{20}\,\mathbf{a}_3$	=	$-x_{20}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{20}\right)b\,\hat{\mathbf{y}}-z_{20}c\,\hat{\mathbf{z}}$	(8*d*)	Se VIII
$\mathbf{B}_{108}$	=	$\left(\frac{1}{2}+x_{20}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{20}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{20}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{20}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{20}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{20}\right)c\,\hat{\mathbf{z}}$	(8*d*)	Se VIII
$\mathbf{B}_{109}$	=	$-x_{20}\,\mathbf{a}_1-y_{20}\,\mathbf{a}_2-z_{20}\,\mathbf{a}_3$	=	$-x_{20}a\,\hat{\mathbf{x}}-y_{20}b\,\hat{\mathbf{y}}-z_{20}c\,\hat{\mathbf{z}}$	(8*d*)	Se VIII
$\mathbf{B}_{110}$	=	$\left(\frac{1}{2}+x_{20}\right)\mathbf{a}_1+y_{20}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{20}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{20}\right)a\,\hat{\mathbf{x}}+y_{20}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{20}\right)c\,\hat{\mathbf{z}}$	(8*d*)	Se VIII
$\mathbf{B}_{111}$	=	$x_{20}\,\mathbf{a}_1+\left(\frac{1}{2}-y_{20}\right)\mathbf{a}_2+z_{20}\,\mathbf{a}_3$	=	$x_{20}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{20}\right)b\,\hat{\mathbf{y}}+z_{20}c\,\hat{\mathbf{z}}$	(8*d*)	Se VIII
$\mathbf{B}_{112}$	=	$\left(\frac{1}{2}-x_{20}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{20}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{20}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{20}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{20}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{20}\right)c\,\hat{\mathbf{z}}$	(8*d*)	Se VIII

References:

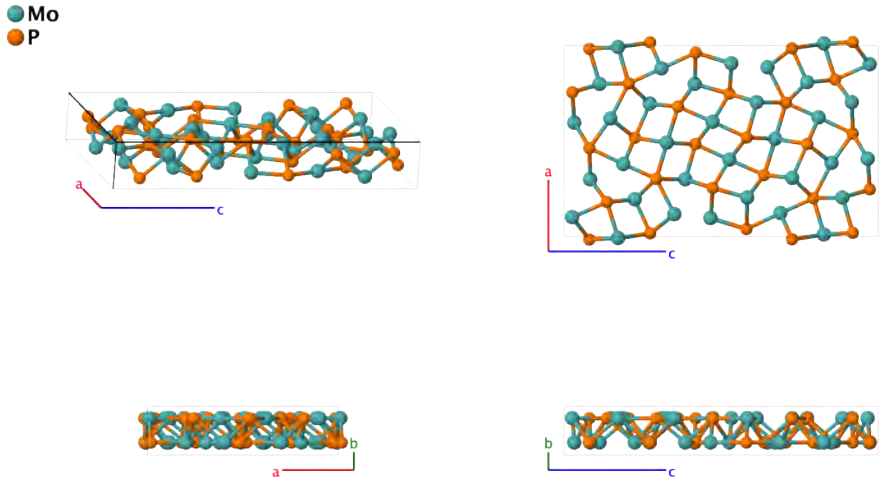
- E. Keulen and A. Vos, *The Crystal Structure of P_4Se_3*, Acta Cryst. **12**, 323–329 (1959), doi:10.1107/S0365110X59000950.

Geometry files:

- CIF: pp. 1636
- POSCAR: pp. 1637

Mo_4P_3 Structure: A4B3_oP56_62_8c_6c

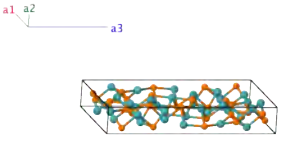
http://aflow.org/prototype-encyclopedia/A4B3_oP56_62_8c_6c



Prototype	:	Mo_4P_3
AFLOW prototype label	:	A4B3_oP56_62_8c_6c
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP56
Space group number	:	62
Space group symbol	:	*Pnma*
AFLOW prototype command	:	`aflow --proto=A4B3_oP56_62_8c_6c --params=`$a, b/a, c/a, x_1, z_1, x_2, z_2, x_3, z_3, x_4, z_4, x_5, z_5, x_6, z_6, x_7, z_7, x_8, z_8, x_9, z_9, x_{10}, z_{10}, x_{11}, z_{11}, x_{12}, z_{12}, x_{13}, z_{13}, x_{14}, z_{14}$

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(4*c*)	Mo I
$\mathbf{B}_2$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	Mo I
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	(4*c*)	Mo I
$\mathbf{B}_4$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	Mo I
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*c*)	Mo II

$$
\begin{aligned}
\mathbf{B}_{6} &= \left(\tfrac{1}{2} - x_2\right)\mathbf{a}_1 + \tfrac{3}{4}\mathbf{a}_2 + \left(\tfrac{1}{2} + z_2\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \tfrac{3}{4}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_2\right)c\,\hat{\mathbf{z}} && (4c) && \text{Mo II}\\
\mathbf{B}_{7} &= -x_2\,\mathbf{a}_1 + \tfrac{3}{4}\mathbf{a}_2 - z_2\,\mathbf{a}_3 &&= -x_2 a\,\hat{\mathbf{x}} + \tfrac{3}{4}b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}} && (4c) && \text{Mo II}\\
\mathbf{B}_{8} &= \left(\tfrac{1}{2} + x_2\right)\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2 + \left(\tfrac{1}{2} - z_2\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \tfrac{1}{4}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_2\right)c\,\hat{\mathbf{z}} && (4c) && \text{Mo II}\\
\mathbf{B}_{9} &= x_3\,\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2 + z_3\,\mathbf{a}_3 &&= x_3 a\,\hat{\mathbf{x}} + \tfrac{1}{4}b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}} && (4c) && \text{Mo III}\\
\mathbf{B}_{10} &= \left(\tfrac{1}{2} - x_3\right)\mathbf{a}_1 + \tfrac{3}{4}\mathbf{a}_2 + \left(\tfrac{1}{2} + z_3\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \tfrac{3}{4}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_3\right)c\,\hat{\mathbf{z}} && (4c) && \text{Mo III}\\
\mathbf{B}_{11} &= -x_3\,\mathbf{a}_1 + \tfrac{3}{4}\mathbf{a}_2 - z_3\,\mathbf{a}_3 &&= -x_3 a\,\hat{\mathbf{x}} + \tfrac{3}{4}b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}} && (4c) && \text{Mo III}\\
\mathbf{B}_{12} &= \left(\tfrac{1}{2} + x_3\right)\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2 + \left(\tfrac{1}{2} - z_3\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \tfrac{1}{4}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_3\right)c\,\hat{\mathbf{z}} && (4c) && \text{Mo III}\\
\mathbf{B}_{13} &= x_4\,\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2 + z_4\,\mathbf{a}_3 &&= x_4 a\,\hat{\mathbf{x}} + \tfrac{1}{4}b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}} && (4c) && \text{Mo IV}\\
\mathbf{B}_{14} &= \left(\tfrac{1}{2} - x_4\right)\mathbf{a}_1 + \tfrac{3}{4}\mathbf{a}_2 + \left(\tfrac{1}{2} + z_4\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \tfrac{3}{4}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_4\right)c\,\hat{\mathbf{z}} && (4c) && \text{Mo IV}\\
\mathbf{B}_{15} &= -x_4\,\mathbf{a}_1 + \tfrac{3}{4}\mathbf{a}_2 - z_4\,\mathbf{a}_3 &&= -x_4 a\,\hat{\mathbf{x}} + \tfrac{3}{4}b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}} && (4c) && \text{Mo IV}\\
\mathbf{B}_{16} &= \left(\tfrac{1}{2} + x_4\right)\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2 + \left(\tfrac{1}{2} - z_4\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \tfrac{1}{4}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_4\right)c\,\hat{\mathbf{z}} && (4c) && \text{Mo IV}\\
\mathbf{B}_{17} &= x_5\,\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2 + z_5\,\mathbf{a}_3 &&= x_5 a\,\hat{\mathbf{x}} + \tfrac{1}{4}b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}} && (4c) && \text{Mo V}\\
\mathbf{B}_{18} &= \left(\tfrac{1}{2} - x_5\right)\mathbf{a}_1 + \tfrac{3}{4}\mathbf{a}_2 + \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + \tfrac{3}{4}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_5\right)c\,\hat{\mathbf{z}} && (4c) && \text{Mo V}\\
\mathbf{B}_{19} &= -x_5\,\mathbf{a}_1 + \tfrac{3}{4}\mathbf{a}_2 - z_5\,\mathbf{a}_3 &&= -x_5 a\,\hat{\mathbf{x}} + \tfrac{3}{4}b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}} && (4c) && \text{Mo V}\\
\mathbf{B}_{20} &= \left(\tfrac{1}{2} + x_5\right)\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2 + \left(\tfrac{1}{2} - z_5\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + \tfrac{1}{4}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_5\right)c\,\hat{\mathbf{z}} && (4c) && \text{Mo V}\\
\mathbf{B}_{21} &= x_6\,\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2 + z_6\,\mathbf{a}_3 &&= x_6 a\,\hat{\mathbf{x}} + \tfrac{1}{4}b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}} && (4c) && \text{Mo VI}\\
\mathbf{B}_{22} &= \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_1 + \tfrac{3}{4}\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_6\right)a\,\hat{\mathbf{x}} + \tfrac{3}{4}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right)c\,\hat{\mathbf{z}} && (4c) && \text{Mo VI}\\
\mathbf{B}_{23} &= -x_6\,\mathbf{a}_1 + \tfrac{3}{4}\mathbf{a}_2 - z_6\,\mathbf{a}_3 &&= -x_6 a\,\hat{\mathbf{x}} + \tfrac{3}{4}b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}} && (4c) && \text{Mo VI}\\
\mathbf{B}_{24} &= \left(\tfrac{1}{2} + x_6\right)\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2 + \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_6\right)a\,\hat{\mathbf{x}} + \tfrac{1}{4}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_6\right)c\,\hat{\mathbf{z}} && (4c) && \text{Mo VI}\\
\mathbf{B}_{25} &= x_7\,\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2 + z_7\,\mathbf{a}_3 &&= x_7 a\,\hat{\mathbf{x}} + \tfrac{1}{4}b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}} && (4c) && \text{Mo VII}\\
\mathbf{B}_{26} &= \left(\tfrac{1}{2} - x_7\right)\mathbf{a}_1 + \tfrac{3}{4}\mathbf{a}_2 + \left(\tfrac{1}{2} + z_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_7\right)a\,\hat{\mathbf{x}} + \tfrac{3}{4}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_7\right)c\,\hat{\mathbf{z}} && (4c) && \text{Mo VII}\\
\mathbf{B}_{27} &= -x_7\,\mathbf{a}_1 + \tfrac{3}{4}\mathbf{a}_2 - z_7\,\mathbf{a}_3 &&= -x_7 a\,\hat{\mathbf{x}} + \tfrac{3}{4}b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}} && (4c) && \text{Mo VII}\\
\mathbf{B}_{28} &= \left(\tfrac{1}{2} + x_7\right)\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2 + \left(\tfrac{1}{2} - z_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_7\right)a\,\hat{\mathbf{x}} + \tfrac{1}{4}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_7\right)c\,\hat{\mathbf{z}} && (4c) && \text{Mo VII}\\
\mathbf{B}_{29} &= x_8\,\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2 + z_8\,\mathbf{a}_3 &&= x_8 a\,\hat{\mathbf{x}} + \tfrac{1}{4}b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}} && (4c) && \text{Mo VIII}\\
\mathbf{B}_{30} &= \left(\tfrac{1}{2} - x_8\right)\mathbf{a}_1 + \tfrac{3}{4}\mathbf{a}_2 + \left(\tfrac{1}{2} + z_8\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_8\right)a\,\hat{\mathbf{x}} + \tfrac{3}{4}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_8\right)c\,\hat{\mathbf{z}} && (4c) && \text{Mo VIII}\\
\mathbf{B}_{31} &= -x_8\,\mathbf{a}_1 + \tfrac{3}{4}\mathbf{a}_2 - z_8\,\mathbf{a}_3 &&= -x_8 a\,\hat{\mathbf{x}} + \tfrac{3}{4}b\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}} && (4c) && \text{Mo VIII}\\
\mathbf{B}_{32} &= \left(\tfrac{1}{2} + x_8\right)\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2 + \left(\tfrac{1}{2} - z_8\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_8\right)a\,\hat{\mathbf{x}} + \tfrac{1}{4}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_8\right)c\,\hat{\mathbf{z}} && (4c) && \text{Mo VIII}\\
\mathbf{B}_{33} &= x_9\,\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2 + z_9\,\mathbf{a}_3 &&= x_9 a\,\hat{\mathbf{x}} + \tfrac{1}{4}b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}} && (4c) && \text{P I}\\
\mathbf{B}_{34} &= \left(\tfrac{1}{2} - x_9\right)\mathbf{a}_1 + \tfrac{3}{4}\mathbf{a}_2 + \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_9\right)a\,\hat{\mathbf{x}} + \tfrac{3}{4}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_9\right)c\,\hat{\mathbf{z}} && (4c) && \text{P I}\\
\mathbf{B}_{35} &= -x_9\,\mathbf{a}_1 + \tfrac{3}{4}\mathbf{a}_2 - z_9\,\mathbf{a}_3 &&= -x_9 a\,\hat{\mathbf{x}} + \tfrac{3}{4}b\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}} && (4c) && \text{P I}\\
\mathbf{B}_{36} &= \left(\tfrac{1}{2} + x_9\right)\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2 + \left(\tfrac{1}{2} - z_9\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_9\right)a\,\hat{\mathbf{x}} + \tfrac{1}{4}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_9\right)c\,\hat{\mathbf{z}} && (4c) && \text{P I}\\
\mathbf{B}_{37} &= x_{10}\,\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 &&= x_{10} a\,\hat{\mathbf{x}} + \tfrac{1}{4}b\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}} && (4c) && \text{P II}\\
\mathbf{B}_{38} &= \left(\tfrac{1}{2} - x_{10}\right)\mathbf{a}_1 + \tfrac{3}{4}\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{10}\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_{10}\right)a\,\hat{\mathbf{x}} + \tfrac{3}{4}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{10}\right)c\,\hat{\mathbf{z}} && (4c) && \text{P II}\\
\mathbf{B}_{39} &= -x_{10}\,\mathbf{a}_1 + \tfrac{3}{4}\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 &&= -x_{10} a\,\hat{\mathbf{x}} + \tfrac{3}{4}b\,\hat{\mathbf{y}} - z_{10} c\,\hat{\mathbf{z}} && (4c) && \text{P II}\\
\mathbf{B}_{40} &= \left(\tfrac{1}{2} + x_{10}\right)\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{10}\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_{10}\right)a\,\hat{\mathbf{x}} + \tfrac{1}{4}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{10}\right)c\,\hat{\mathbf{z}} && (4c) && \text{P II}\\
\mathbf{B}_{41} &= x_{11}\,\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 &&= x_{11} a\,\hat{\mathbf{x}} + \tfrac{1}{4}b\,\hat{\mathbf{y}} + z_{11} c\,\hat{\mathbf{z}} && (4c) && \text{P III}
\end{aligned}
$$

$$\mathbf{B}_{42} = \left(\tfrac{1}{2} - x_{11}\right)\mathbf{a}_1 + \tfrac{3}{4}\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{11}\right)a\,\hat{\mathbf{x}} + \tfrac{3}{4}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{11}\right)c\,\hat{\mathbf{z}} \quad (4c) \quad \text{P III}$$

$$\mathbf{B}_{43} = -x_{11}\mathbf{a}_1 + \tfrac{3}{4}\mathbf{a}_2 - z_{11}\mathbf{a}_3 = -x_{11}a\,\hat{\mathbf{x}} + \tfrac{3}{4}b\,\hat{\mathbf{y}} - z_{11}c\,\hat{\mathbf{z}} \quad (4c) \quad \text{P III}$$

$$\mathbf{B}_{44} = \left(\tfrac{1}{2} + x_{11}\right)\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{11}\right)a\,\hat{\mathbf{x}} + \tfrac{1}{4}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{11}\right)c\,\hat{\mathbf{z}} \quad (4c) \quad \text{P III}$$

$$\mathbf{B}_{45} = x_{12}\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2 + z_{12}\mathbf{a}_3 = x_{12}a\,\hat{\mathbf{x}} + \tfrac{1}{4}b\,\hat{\mathbf{y}} + z_{12}c\,\hat{\mathbf{z}} \quad (4c) \quad \text{P IV}$$

$$\mathbf{B}_{46} = \left(\tfrac{1}{2} - x_{12}\right)\mathbf{a}_1 + \tfrac{3}{4}\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{12}\right)a\,\hat{\mathbf{x}} + \tfrac{3}{4}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{12}\right)c\,\hat{\mathbf{z}} \quad (4c) \quad \text{P IV}$$

$$\mathbf{B}_{47} = -x_{12}\mathbf{a}_1 + \tfrac{3}{4}\mathbf{a}_2 - z_{12}\mathbf{a}_3 = -x_{12}a\,\hat{\mathbf{x}} + \tfrac{3}{4}b\,\hat{\mathbf{y}} - z_{12}c\,\hat{\mathbf{z}} \quad (4c) \quad \text{P IV}$$

$$\mathbf{B}_{48} = \left(\tfrac{1}{2} + x_{12}\right)\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{12}\right)a\,\hat{\mathbf{x}} + \tfrac{1}{4}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{12}\right)c\,\hat{\mathbf{z}} \quad (4c) \quad \text{P IV}$$

$$\mathbf{B}_{49} = x_{13}\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2 + z_{13}\mathbf{a}_3 = x_{13}a\,\hat{\mathbf{x}} + \tfrac{1}{4}b\,\hat{\mathbf{y}} + z_{13}c\,\hat{\mathbf{z}} \quad (4c) \quad \text{P V}$$

$$\mathbf{B}_{50} = \left(\tfrac{1}{2} - x_{13}\right)\mathbf{a}_1 + \tfrac{3}{4}\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{13}\right)a\,\hat{\mathbf{x}} + \tfrac{3}{4}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{13}\right)c\,\hat{\mathbf{z}} \quad (4c) \quad \text{P V}$$

$$\mathbf{B}_{51} = -x_{13}\mathbf{a}_1 + \tfrac{3}{4}\mathbf{a}_2 - z_{13}\mathbf{a}_3 = -x_{13}a\,\hat{\mathbf{x}} + \tfrac{3}{4}b\,\hat{\mathbf{y}} - z_{13}c\,\hat{\mathbf{z}} \quad (4c) \quad \text{P V}$$

$$\mathbf{B}_{52} = \left(\tfrac{1}{2} + x_{13}\right)\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{13}\right)a\,\hat{\mathbf{x}} + \tfrac{1}{4}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{13}\right)c\,\hat{\mathbf{z}} \quad (4c) \quad \text{P V}$$

$$\mathbf{B}_{53} = x_{14}\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2 + z_{14}\mathbf{a}_3 = x_{14}a\,\hat{\mathbf{x}} + \tfrac{1}{4}b\,\hat{\mathbf{y}} + z_{14}c\,\hat{\mathbf{z}} \quad (4c) \quad \text{P VI}$$

$$\mathbf{B}_{54} = \left(\tfrac{1}{2} - x_{14}\right)\mathbf{a}_1 + \tfrac{3}{4}\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{14}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{14}\right)a\,\hat{\mathbf{x}} + \tfrac{3}{4}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{14}\right)c\,\hat{\mathbf{z}} \quad (4c) \quad \text{P VI}$$

$$\mathbf{B}_{55} = -x_{14}\mathbf{a}_1 + \tfrac{3}{4}\mathbf{a}_2 - z_{14}\mathbf{a}_3 = -x_{14}a\,\hat{\mathbf{x}} + \tfrac{3}{4}b\,\hat{\mathbf{y}} - z_{14}c\,\hat{\mathbf{z}} \quad (4c) \quad \text{P VI}$$

$$\mathbf{B}_{56} = \left(\tfrac{1}{2} + x_{14}\right)\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{14}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{14}\right)a\,\hat{\mathbf{x}} + \tfrac{1}{4}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{14}\right)c\,\hat{\mathbf{z}} \quad (4c) \quad \text{P VI}$$

References:

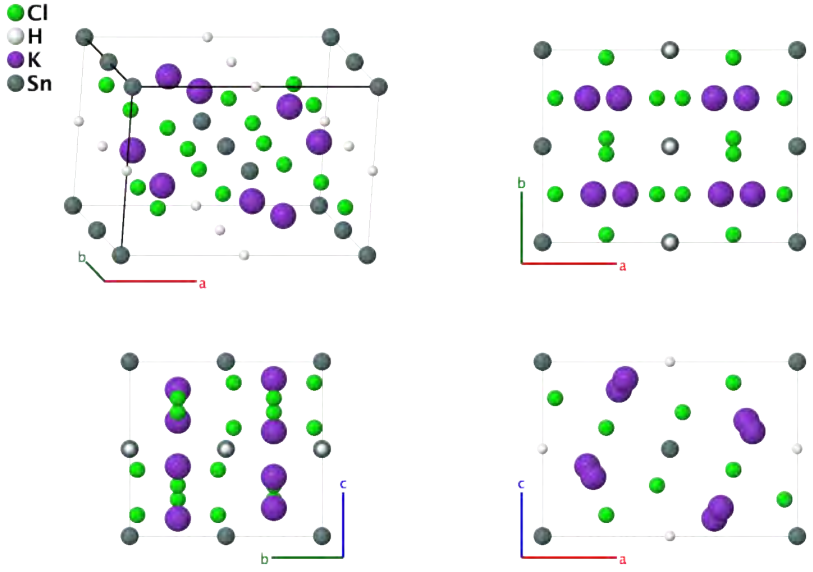
- S. Rundqvist, *The Crystal Structure of Mo_4P_3*, Acta Chem. Scand. **19**, 393–400 (1965), doi:10.3891/acta.chem.scand.19-0393.

Geometry files:

- CIF: pp. 1637
- POSCAR: pp. 1638

$K_2SnCl_4{\cdot}H_2O$ ($E3_5$) Structure: A4BC2D_oP32_62_2cd_b_2c_a

http://aflow.org/prototype-encyclopedia/A4BC2D_oP32_62_2cd_b_2c_a

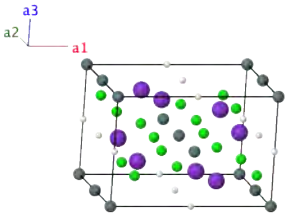


Prototype	:	$Cl_4(H_2O)K_2Sn$
AFLOW prototype label	:	A4BC2D_oP32_62_2cd_b_2c_a
***Strukturbericht* designation**	:	$E3_5$
Pearson symbol	:	oP32
Space group number	:	62
Space group symbol	:	*Pnma*
AFLOW prototype command	:	`aflow --proto=A4BC2D_oP32_62_2cd_b_2c_a --params=`$a, b/a, c/a, x_3, z_3, x_4, z_4, x_5, z_5, x_6, z_6, x_7, y_7, z_7$

- This is the structure listed in (Hermann, 1939) as the $E3_5$ structure.
- (Kamenar, 1962) proposed a somewhat different version of the structure for this crystal. We have no clear indication as to which structure is correct.
- The positions of the hydrogen atoms in the water molecules were not determined, so we only give the positions of the oxygen atoms (labeled as H_2O).
- (Brasseur, 1939) gives the crystal structure and atomic positions in the *Pbnm* setting of space group #62. We have used FINDSYM to show the structure in the standard *Pnma* setting.
- (Hermann, 1939) gives incorrect values for the Wyckoff positions of the Cl III atoms. The correct coordinates are ($x_{\mathrm{ClIII}} = 0.79$, $y_{\mathrm{ClIII}} = +0.05$).

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(4*a*)	Sn
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*a*)	Sn
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}b\,\hat{\mathbf{y}}$	(4*a*)	Sn
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*a*)	Sn
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	(4*b*)	H_2O
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	(4*b*)	H_2O
$\mathbf{B}_7$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*b*)	H_2O
$\mathbf{B}_8$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$	(4*b*)	H_2O
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	(4*c*)	Cl I
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(4*c*)	Cl I
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	(4*c*)	Cl I
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(4*c*)	Cl I
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	(4*c*)	Cl II
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(4*c*)	Cl II
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	(4*c*)	Cl II
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(4*c*)	Cl II
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_5c\,\hat{\mathbf{z}}$	(4*c*)	K I
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	(4*c*)	K I
$\mathbf{B}_{19}$	=	$-x_5\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_5c\,\hat{\mathbf{z}}$	(4*c*)	K I
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	(4*c*)	K I
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_6c\,\hat{\mathbf{z}}$	(4*c*)	K II
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\,\hat{\mathbf{z}}$	(4*c*)	K II
$\mathbf{B}_{23}$	=	$-x_6\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$-x_6a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_6c\,\hat{\mathbf{z}}$	(4*c*)	K II
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\,\hat{\mathbf{z}}$	(4*c*)	K II
$\mathbf{B}_{25}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$x_7a\,\hat{\mathbf{x}} + y_7b\,\hat{\mathbf{y}} + z_7c\,\hat{\mathbf{z}}$	(8*d*)	Cl III
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} - x_7\right)\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_7\right)a\,\hat{\mathbf{x}} - y_7b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right)c\,\hat{\mathbf{z}}$	(8*d*)	Cl III
$\mathbf{B}_{27}$	=	$-x_7\,\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$-x_7a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right)b\,\hat{\mathbf{y}} - z_7c\,\hat{\mathbf{z}}$	(8*d*)	Cl III

$\mathbf{B}_{28}$	=	$\left(\frac{1}{2}+x_7\right)\mathbf{a}_1+\left(\frac{1}{2}-y_7\right)\mathbf{a}_2+\left(\frac{1}{2}-z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_7\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_7\right)c\,\hat{\mathbf{z}}$	(8*d*)	Cl III
$\mathbf{B}_{29}$	=	$-x_7\,\mathbf{a}_1-y_7\,\mathbf{a}_2-z_7\,\mathbf{a}_3$	=	$-x_7a\,\hat{\mathbf{x}}-y_7b\,\hat{\mathbf{y}}-z_7c\,\hat{\mathbf{z}}$	(8*d*)	Cl III
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2}+x_7\right)\mathbf{a}_1+y_7\,\mathbf{a}_2+\left(\frac{1}{2}-z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{x}}+y_7b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_7\right)c\,\hat{\mathbf{z}}$	(8*d*)	Cl III
$\mathbf{B}_{31}$	=	$x_7\,\mathbf{a}_1+\left(\frac{1}{2}-y_7\right)\mathbf{a}_2+z_7\,\mathbf{a}_3$	=	$x_7a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_7\right)b\,\hat{\mathbf{y}}+z_7c\,\hat{\mathbf{z}}$	(8*d*)	Cl III
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2}-x_7\right)\mathbf{a}_1+\left(\frac{1}{2}+y_7\right)\mathbf{a}_2+\left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_7\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_7\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	(8*d*)	Cl III

References:

- H. Brasseur and A. de Rassenfosse, *The Crystal Structure of Hydrated Potassium Chlorostannite*, Zeitschrift für Kristallographie - Crystalline Materials **101**, 389–395 (1939), doi:10.1524/zkri.1939.101.1.389.
- B. Kamenar and D. Grdenić, *The crystal structure of potassium chloride trichlorostannite hydrate, KCl, $KSnCl_3$, H_2O*, J. Inorg. Nucl. Chem. **24**, 1039–1045 (1962), doi:10.1016/0022-1902(62)80247-4.

Found in:

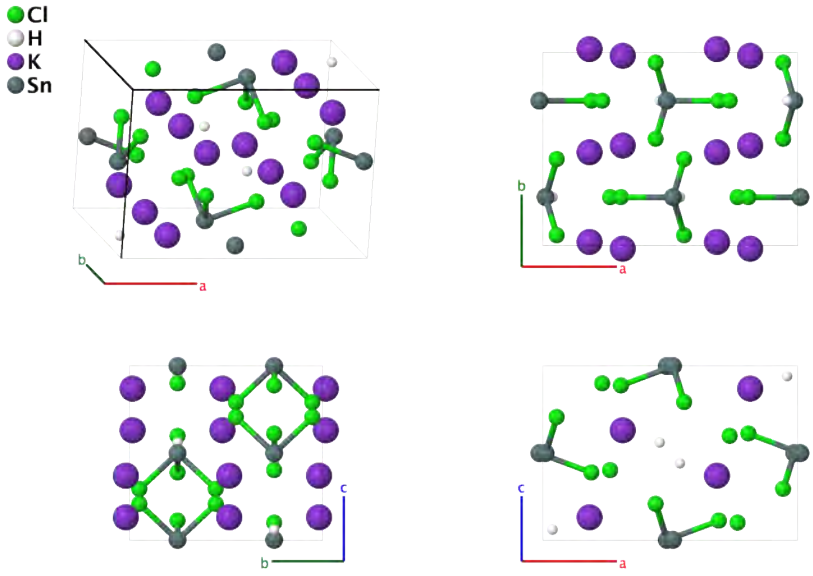
- K. Herrmann, ed., *Strukturbericht Band VII 1939* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1943).

Geometry files:

- CIF: pp. 1638
- POSCAR: pp. 1638

$K_2SnCl_4 \cdot H_2O$ Structure: A4BC2D_oP32_62_2cd_c_d_c

http://aflow.org/prototype-encyclopedia/A4BC2D_oP32_62_2cd_c_d_c

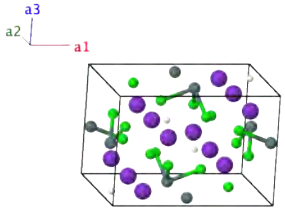


Prototype	:	$Cl_4(H_2O)K_2Sn$
AFLOW prototype label	:	A4BC2D_oP32_62_2cd_c_d_c
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP32
Space group number	:	62
Space group symbol	:	*Pnma*
AFLOW prototype command	:	`aflow --proto=A4BC2D_oP32_62_2cd_c_d_c --params=`$a, b/a, c/a, x_1, z_1, x_2, z_2, x_3, z_3, x_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6$

- This is *not* the version of the structure given in (Hermann, 1939) as *Strukturbericht* $E3_3$. That structure was determined by (Brasseur, 1939). We have no clear indication as to which structure is correct.
- The positions of the hydrogen atoms in the water molecules were not determined, so we only give the positions of the oxygen atoms (labeled as H_2O).
- (Kamenar, 1962) gives the crystal structure and atomic positions in the *Pbnm* setting of space group #62. We have used FINDSYM to show the structure in the standard *Pnma* setting.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(4c)	Cl I
$\mathbf{B}_2$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	(4c)	Cl I
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	(4c)	Cl I
$\mathbf{B}_4$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right) c\,\hat{\mathbf{z}}$	(4c)	Cl I
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4c)	Cl II
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	(4c)	Cl II
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(4c)	Cl II
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right) c\,\hat{\mathbf{z}}$	(4c)	Cl II
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4c)	H_2O
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	(4c)	H_2O
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(4c)	H_2O
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right) c\,\hat{\mathbf{z}}$	(4c)	H_2O
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(4c)	Sn
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	(4c)	Sn
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(4c)	Sn
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\,\hat{\mathbf{z}}$	(4c)	Sn
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(8d)	Cl III
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right) a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	(8d)	Cl III
$\mathbf{B}_{19}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right) b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(8d)	Cl III
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right) c\,\hat{\mathbf{z}}$	(8d)	Cl III
$\mathbf{B}_{21}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(8d)	Cl III
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + y_5\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right) c\,\hat{\mathbf{z}}$	(8d)	Cl III
$\mathbf{B}_{23}$	=	$x_5\,\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right) b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(8d)	Cl III
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	(8d)	Cl III
$\mathbf{B}_{25}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(8d)	K
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	(8d)	K
$\mathbf{B}_{27}$	=	$-x_6\,\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right) b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(8d)	K
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right) c\,\hat{\mathbf{z}}$	(8d)	K
$\mathbf{B}_{29}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(8d)	K
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + y_6\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right) c\,\hat{\mathbf{z}}$	(8d)	K
$\mathbf{B}_{31}$	=	$x_6\,\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right) b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(8d)	K

$$\mathbf{B}_{32} = \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_6\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_6\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right)c\,\hat{\mathbf{z}} \quad (8d) \quad \text{K}$$

References:

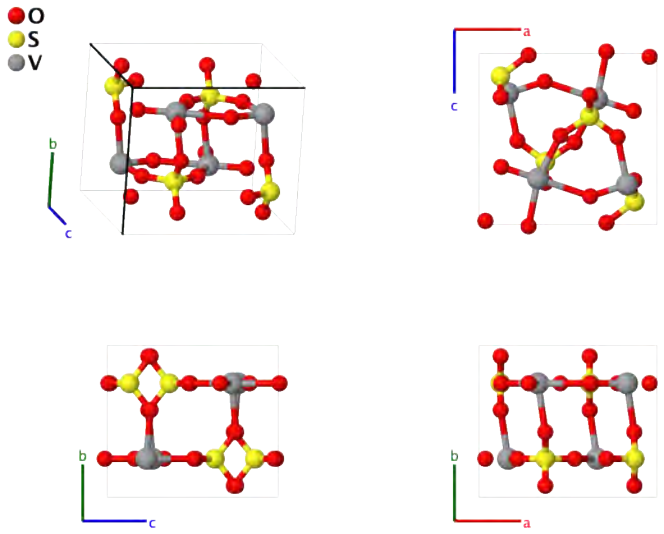
- B. Kamenar and D. Grdenić, *The crystal structure of potassium chloride trichlorostannite hydrate, KCl, $KSnCl_3$, H_2O*, J. Inorg. Nucl. Chem. **24**, 1039–1045 (1962), doi:10.1016/0022-1902(62)80247-4.
- K. Herrmann, ed., *Strukturbericht Band VII 1939* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1943).
- H. Brasseur and A. de Rassenfosse, *The Crystal Structure of Hydrated Potassium Chlorostannite*, Zeitschrift für Kristallographie - Crystalline Materials **101**, 389–395 (1939), doi:10.1524/zkri.1939.101.1.389.

Geometry files:

- CIF: pp. 1639
- POSCAR: pp. 1639

$VOSO_4$ Structure: A5BC_oP28_62_3cd_c_c

http://aflow.org/prototype-encyclopedia/A5BC_oP28_62_3cd_c_c

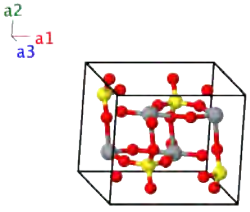


Prototype	:	O_5SV
AFLOW prototype label	:	A5BC_oP28_62_3cd_c_c
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP28
Space group number	:	62
Space group symbol	:	*Pnma*
AFLOW prototype command	:	`aflow --proto=A5BC_oP28_62_3cd_c_c --params=`$a, b/a, c/a, x_1, z_1, x_2, z_2, x_3, z_3, x_4, z_4, x_5, z_5, x_6, y_6, z_6$

- This structure consists of distorted SO_4 tetrahedra linked to distorted VO_6 tetrahedra.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_2$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	(4*c*)	O I

$\mathbf{B}_4$	=	$\left(\frac{1}{2}+x_1\right)\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+\left(\frac{1}{2}-z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_1\right)a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_1\right)c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_5$	=	$x_2\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+z_2\mathbf{a}_3$	=	$x_2a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+z_2c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_6$	=	$\left(\frac{1}{2}-x_2\right)\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2+\left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_2\right)a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_2\right)c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_7$	=	$-x_2\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2-z_2\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}-z_2c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_8$	=	$\left(\frac{1}{2}+x_2\right)\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+\left(\frac{1}{2}-z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_2\right)a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_2\right)c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_9$	=	$x_3\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+z_3\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+z_3c\,\hat{\mathbf{z}}$	(4*c*)	O III
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2}-x_3\right)\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2+\left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_3\right)c\,\hat{\mathbf{z}}$	(4*c*)	O III
$\mathbf{B}_{11}$	=	$-x_3\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2-z_3\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}-z_3c\,\hat{\mathbf{z}}$	(4*c*)	O III
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2}+x_3\right)\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+\left(\frac{1}{2}-z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_3\right)c\,\hat{\mathbf{z}}$	(4*c*)	O III
$\mathbf{B}_{13}$	=	$x_4\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+z_4\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+z_4c\,\hat{\mathbf{z}}$	(4*c*)	S
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2}-x_4\right)\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2+\left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_4\right)c\,\hat{\mathbf{z}}$	(4*c*)	S
$\mathbf{B}_{15}$	=	$-x_4\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2-z_4\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}-z_4c\,\hat{\mathbf{z}}$	(4*c*)	S
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2}+x_4\right)\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+\left(\frac{1}{2}-z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_4\right)c\,\hat{\mathbf{z}}$	(4*c*)	S
$\mathbf{B}_{17}$	=	$x_5\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+z_5\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+z_5c\,\hat{\mathbf{z}}$	(4*c*)	V
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2}-x_5\right)\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	(4*c*)	V
$\mathbf{B}_{19}$	=	$-x_5\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2-z_5\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}-z_5c\,\hat{\mathbf{z}}$	(4*c*)	V
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2}+x_5\right)\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+\left(\frac{1}{2}-z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_5\right)c\,\hat{\mathbf{z}}$	(4*c*)	V
$\mathbf{B}_{21}$	=	$x_6\mathbf{a}_1+y_6\mathbf{a}_2+z_6\mathbf{a}_3$	=	$x_6a\,\hat{\mathbf{x}}+y_6b\,\hat{\mathbf{y}}+z_6c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2}-x_6\right)\mathbf{a}_1-y_6\mathbf{a}_2+\left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{x}}-y_6b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{23}$	=	$-x_6\mathbf{a}_1+\left(\frac{1}{2}+y_6\right)\mathbf{a}_2-z_6\mathbf{a}_3$	=	$-x_6a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_6\right)b\,\hat{\mathbf{y}}-z_6c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2}+x_6\right)\mathbf{a}_1+\left(\frac{1}{2}-y_6\right)\mathbf{a}_2+\left(\frac{1}{2}-z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_6\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_6\right)c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{25}$	=	$-x_6\mathbf{a}_1-y_6\mathbf{a}_2-z_6\mathbf{a}_3$	=	$-x_6a\,\hat{\mathbf{x}}-y_6b\,\hat{\mathbf{y}}-z_6c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2}+x_6\right)\mathbf{a}_1+y_6\mathbf{a}_2+\left(\frac{1}{2}-z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{x}}+y_6b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_6\right)c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{27}$	=	$x_6\mathbf{a}_1+\left(\frac{1}{2}-y_6\right)\mathbf{a}_2+z_6\mathbf{a}_3$	=	$x_6a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_6\right)b\,\hat{\mathbf{y}}+z_6c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2}-x_6\right)\mathbf{a}_1+\left(\frac{1}{2}+y_6\right)\mathbf{a}_2+\left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_6\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	(8*d*)	O IV

References:

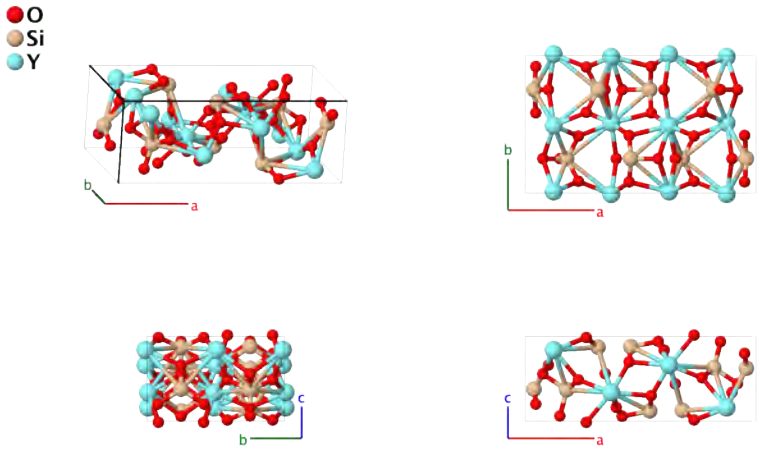
- P. Kierkegaard, J. M. Longo, and B.-O. Marinder, *Note on the Crystal Structure of $VOSO_4$*, Acta Chem. Scand. **19**, 763–764 (1965), doi:10.3891/acta.chem.scand.19-0763.

Geometry files:

- CIF: pp. 1639
- POSCAR: pp. 1640

Possible δ-$Y_2Si_2O_7$ Structure: A7B2C2_oP44_62_3c2d_2c_d

http://aflow.org/prototype-encyclopedia/A7B2C2_oP44_62_3c2d_2c_d



Prototype	:	$O_7Si_2Y_2$
AFLOW prototype label	:	A7B2C2_oP44_62_3c2d_2c_d
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP44
Space group number	:	62
Space group symbol	:	*Pnma*
AFLOW prototype command	:	`aflow --proto=A7B2C2_oP44_62_3c2d_2c_d --params=`$a, b/a, c/a, x_1, z_1, x_2, z_2, x_3, z_3, x_4, z_4, x_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8$

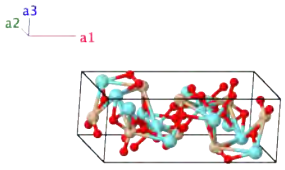
Other compounds with this structure

- δ-$Ho_2O_7Si_2$, δ-$Dy_2O_7Si_2$, δ-$Y_2O_7Si_2$, and $Li_2S_2O_7$

- (Dias, 1990) found that some structures of $RE_2Si_2O_7$ (RE = Ho, Dy, Gd, Y) were in the centro-symmetric orthorhombic *Pnma* #62 space group, in which case this would be the prototype of δ-$RE_2Si_2O_7$ (Becerro, 2004). However, (Smolin, 1970) found δ-$Gd_2Si_2O_7$ to be in the non-centro-symmetric $Pna2_1$ #33 space group. (Becerro, 2004) found only one yttrium site in the δ-structure, supporting (Dias, 1990). In addition, if we allow a small amount of uncertainty (0.2 Å) in positions of the $Pna2_1$ structure, AFLOW-SYM and FINDSYM place this structure in the *Pnma* group. Nevertheless we have found no work explicitly stating that the structure of (Smolin, 1970) is in error, and indeed (Christensen, 1994) found δ-$Y_2Si_2O_7$ in space group $Pna2_1$. Given this ambiguity, we list $Y_2Si_2O_7$ only as a possible prototype for the δ-phase pyrosilicates.

Simple Orthorhombic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= a\,\hat{\mathbf{x}}\\ \mathbf{a}_2 &= b\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	$x_1 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	$(4c)$	O I
$\mathbf{B}_2$	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	$(4c)$	O I
$\mathbf{B}_3$	$-x_1\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	$-x_1 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	$(4c)$	O I
$\mathbf{B}_4$	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_1\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right) c\,\hat{\mathbf{z}}$	$(4c)$	O I
$\mathbf{B}_5$	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	$x_2 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(4c)$	O II
$\mathbf{B}_6$	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	$(4c)$	O II
$\mathbf{B}_7$	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	$-x_2 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	$(4c)$	O II
$\mathbf{B}_8$	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right) c\,\hat{\mathbf{z}}$	$(4c)$	O II
$\mathbf{B}_9$	$x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$x_3 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(4c)$	O III
$\mathbf{B}_{10}$	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_3\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	$(4c)$	O III
$\mathbf{B}_{11}$	$-x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	$-x_3 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(4c)$	O III
$\mathbf{B}_{12}$	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_3\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right) c\,\hat{\mathbf{z}}$	$(4c)$	O III
$\mathbf{B}_{13}$	$x_4\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$x_4 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(4c)$	Si I
$\mathbf{B}_{14}$	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	$(4c)$	Si I
$\mathbf{B}_{15}$	$-x_4\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	$-x_4 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	$(4c)$	Si I
$\mathbf{B}_{16}$	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\,\hat{\mathbf{z}}$	$(4c)$	Si I
$\mathbf{B}_{17}$	$x_5\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	$x_5 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(4c)$	Si II
$\mathbf{B}_{18}$	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_5\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	$(4c)$	Si II
$\mathbf{B}_{19}$	$-x_5\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	$-x_5 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	$(4c)$	Si II
$\mathbf{B}_{20}$	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right) c\,\hat{\mathbf{z}}$	$(4c)$	Si II
$\mathbf{B}_{21}$	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(8d)$	O IV
$\mathbf{B}_{22}$	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	$(8d)$	O IV
$\mathbf{B}_{23}$	$-x_6\,\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3$	$-x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right) b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	$(8d)$	O IV
$\mathbf{B}_{24}$	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right) c\,\hat{\mathbf{z}}$	$(8d)$	O IV
$\mathbf{B}_{25}$	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	$-x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	$(8d)$	O IV
$\mathbf{B}_{26}$	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + y_6\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right) c\,\hat{\mathbf{z}}$	$(8d)$	O IV
$\mathbf{B}_{27}$	$x_6\,\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + z_6\,\mathbf{a}_3$	$x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right) b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(8d)$	O IV

$\mathbf{B}_{28}$	=	$\left(\frac{1}{2}-x_6\right)\mathbf{a}_1+\left(\frac{1}{2}+y_6\right)\mathbf{a}_2+\left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_6\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{29}$	=	$x_7\,\mathbf{a}_1+y_7\,\mathbf{a}_2+z_7\,\mathbf{a}_3$	=	$x_7a\,\hat{\mathbf{x}}+y_7b\,\hat{\mathbf{y}}+z_7c\,\hat{\mathbf{z}}$	(8*d*)	O V
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2}-x_7\right)\mathbf{a}_1-y_7\,\mathbf{a}_2+\left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_7\right)a\,\hat{\mathbf{x}}-y_7b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	(8*d*)	O V
$\mathbf{B}_{31}$	=	$-x_7\,\mathbf{a}_1+\left(\frac{1}{2}+y_7\right)\mathbf{a}_2-z_7\,\mathbf{a}_3$	=	$-x_7a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_7\right)b\,\hat{\mathbf{y}}-z_7c\,\hat{\mathbf{z}}$	(8*d*)	O V
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2}+x_7\right)\mathbf{a}_1+\left(\frac{1}{2}-y_7\right)\mathbf{a}_2+\left(\frac{1}{2}-z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_7\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_7\right)c\,\hat{\mathbf{z}}$	(8*d*)	O V
$\mathbf{B}_{33}$	=	$-x_7\,\mathbf{a}_1-y_7\,\mathbf{a}_2-z_7\,\mathbf{a}_3$	=	$-x_7a\,\hat{\mathbf{x}}-y_7b\,\hat{\mathbf{y}}-z_7c\,\hat{\mathbf{z}}$	(8*d*)	O V
$\mathbf{B}_{34}$	=	$\left(\frac{1}{2}+x_7\right)\mathbf{a}_1+y_7\,\mathbf{a}_2+\left(\frac{1}{2}-z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{x}}+y_7b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_7\right)c\,\hat{\mathbf{z}}$	(8*d*)	O V
$\mathbf{B}_{35}$	=	$x_7\,\mathbf{a}_1+\left(\frac{1}{2}-y_7\right)\mathbf{a}_2+z_7\,\mathbf{a}_3$	=	$x_7a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_7\right)b\,\hat{\mathbf{y}}+z_7c\,\hat{\mathbf{z}}$	(8*d*)	O V
$\mathbf{B}_{36}$	=	$\left(\frac{1}{2}-x_7\right)\mathbf{a}_1+\left(\frac{1}{2}+y_7\right)\mathbf{a}_2+\left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_7\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_7\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	(8*d*)	O V
$\mathbf{B}_{37}$	=	$x_8\,\mathbf{a}_1+y_8\,\mathbf{a}_2+z_8\,\mathbf{a}_3$	=	$x_8a\,\hat{\mathbf{x}}+y_8b\,\hat{\mathbf{y}}+z_8c\,\hat{\mathbf{z}}$	(8*d*)	Y
$\mathbf{B}_{38}$	=	$\left(\frac{1}{2}-x_8\right)\mathbf{a}_1-y_8\,\mathbf{a}_2+\left(\frac{1}{2}+z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_8\right)a\,\hat{\mathbf{x}}-y_8b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_8\right)c\,\hat{\mathbf{z}}$	(8*d*)	Y
$\mathbf{B}_{39}$	=	$-x_8\,\mathbf{a}_1+\left(\frac{1}{2}+y_8\right)\mathbf{a}_2-z_8\,\mathbf{a}_3$	=	$-x_8a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_8\right)b\,\hat{\mathbf{y}}-z_8c\,\hat{\mathbf{z}}$	(8*d*)	Y
$\mathbf{B}_{40}$	=	$\left(\frac{1}{2}+x_8\right)\mathbf{a}_1+\left(\frac{1}{2}-y_8\right)\mathbf{a}_2+\left(\frac{1}{2}-z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_8\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_8\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_8\right)c\,\hat{\mathbf{z}}$	(8*d*)	Y
$\mathbf{B}_{41}$	=	$-x_8\,\mathbf{a}_1-y_8\,\mathbf{a}_2-z_8\,\mathbf{a}_3$	=	$-x_8a\,\hat{\mathbf{x}}-y_8b\,\hat{\mathbf{y}}-z_8c\,\hat{\mathbf{z}}$	(8*d*)	Y
$\mathbf{B}_{42}$	=	$\left(\frac{1}{2}+x_8\right)\mathbf{a}_1+y_8\,\mathbf{a}_2+\left(\frac{1}{2}-z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_8\right)a\,\hat{\mathbf{x}}+y_8b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_8\right)c\,\hat{\mathbf{z}}$	(8*d*)	Y
$\mathbf{B}_{43}$	=	$x_8\,\mathbf{a}_1+\left(\frac{1}{2}-y_8\right)\mathbf{a}_2+z_8\,\mathbf{a}_3$	=	$x_8a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_8\right)b\,\hat{\mathbf{y}}+z_8c\,\hat{\mathbf{z}}$	(8*d*)	Y
$\mathbf{B}_{44}$	=	$\left(\frac{1}{2}-x_8\right)\mathbf{a}_1+\left(\frac{1}{2}+y_8\right)\mathbf{a}_2+\left(\frac{1}{2}+z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_8\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_8\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_8\right)c\,\hat{\mathbf{z}}$	(8*d*)	Y

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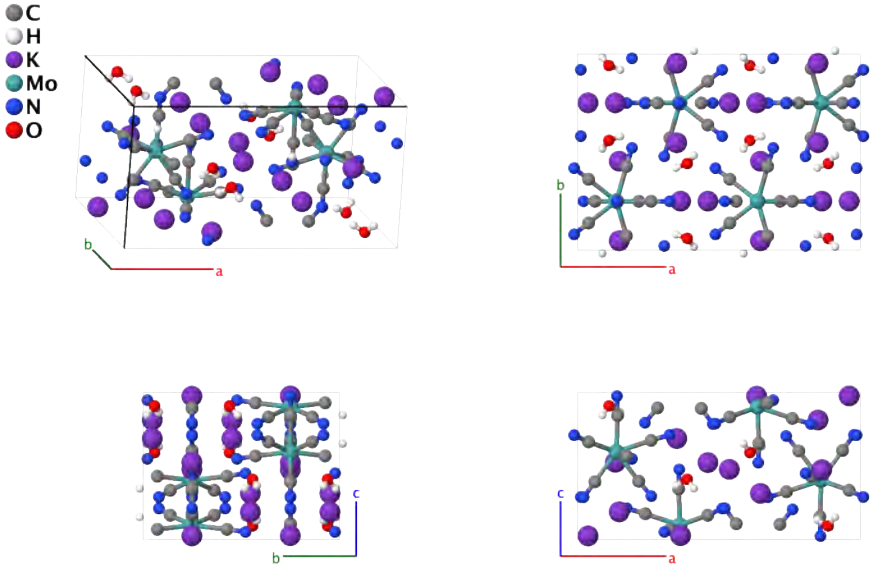
Geometry files:

- CIF: pp. 1640
- POSCAR: pp. 1640

$K_4[Mo(CN)_8]\cdot 2H_2O$ ($F2_1$) Structure: A8B4C4DE8F2_oP108_62_4c2d_2d_2cd_c_4c2d_d

http://aflow.org/prototype-encyclopedia/A8B4C4DE8F2_oP108_62_4c2d_2d_2cd_c_4c2d_d

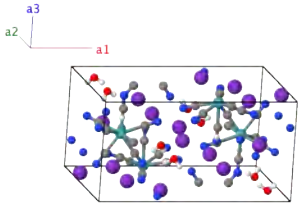


Prototype	:	$C_8H_4K_4MoN_8O_2$
AFLOW prototype label	:	A8B4C4DE8F2_oP108_62_4c2d_2d_2cd_c_4c2d_d
***Strukturbericht* designation**	:	$F2_1$
Pearson symbol	:	oP108
Space group number	:	62
Space group symbol	:	*Pnma*
AFLOW prototype command	:	`aflow --proto=A8B4C4DE8F2_oP108_62_4c2d_2d_2cd_c_4c2d_d --params=`$a, b/a, c/a, x_1, z_1, x_2, z_2, x_3, z_3, x_4, z_4, x_5, z_5, x_6, z_6, x_7, z_7, x_8, z_8, x_9, z_9, x_{10}, z_{10}, x_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}, x_{17}, y_{17}, z_{17}, x_{18}, y_{18}, z_{18}, x_{19}, y_{19}, z_{19}$

- (Hoard, 1939) originally determined this structure, but were unable to locate the hydrogen atoms. (Herrmann, 1943) gave this the *Strukturbericht* designation $F2_1$. (Typilo, 2010) were able to refine the structure at 173 K, including the hydrogen positions. Since the space group and Wyckoff positions are otherwise unchanged we use the newer structure as our prototype.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	($4c$)	C I
$\mathbf{B}_2$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	($4c$)	C I
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	($4c$)	C I
$\mathbf{B}_4$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right) c\,\hat{\mathbf{z}}$	($4c$)	C I
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	($4c$)	C II
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	($4c$)	C II
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	($4c$)	C II
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right) c\,\hat{\mathbf{z}}$	($4c$)	C II
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	($4c$)	C III
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	($4c$)	C III
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	($4c$)	C III
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right) c\,\hat{\mathbf{z}}$	($4c$)	C III
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	($4c$)	C IV
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	($4c$)	C IV
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	($4c$)	C IV
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\,\hat{\mathbf{z}}$	($4c$)	C IV
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	($4c$)	K I
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	($4c$)	K I
$\mathbf{B}_{19}$	=	$-x_5\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	($4c$)	K I
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right) c\,\hat{\mathbf{z}}$	($4c$)	K I
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	($4c$)	K II
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	($4c$)	K II
$\mathbf{B}_{23}$	=	$-x_6\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	($4c$)	K II
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right) c\,\hat{\mathbf{z}}$	($4c$)	K II
$\mathbf{B}_{25}$	=	$x_7\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	($4c$)	Mo
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} - x_7\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_7\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right) c\,\hat{\mathbf{z}}$	($4c$)	Mo
$\mathbf{B}_{27}$	=	$-x_7\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	($4c$)	Mo
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} + x_7\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_7\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right) c\,\hat{\mathbf{z}}$	($4c$)	Mo
$\mathbf{B}_{29}$	=	$x_8\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	($4c$)	N I
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2} - x_8\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_8\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right) c\,\hat{\mathbf{z}}$	($4c$)	N I
$\mathbf{B}_{31}$	=	$-x_8\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	($4c$)	N I
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2} + x_8\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_8\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right) c\,\hat{\mathbf{z}}$	($4c$)	N I
$\mathbf{B}_{33}$	=	$x_9\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	($4c$)	N II
$\mathbf{B}_{34}$	=	$\left(\frac{1}{2} - x_9\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_9\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right) c\,\hat{\mathbf{z}}$	($4c$)	N II

$\mathbf{B}_{35}$	=	$-x_9\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}}$	(4*c*)	N II
$\mathbf{B}_{36}$	=	$\left(\frac{1}{2} + x_9\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_9\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_9\right) c\,\hat{\mathbf{z}}$	(4*c*)	N II
$\mathbf{B}_{37}$	=	$x_{10}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$x_{10} a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}}$	(4*c*)	N III
$\mathbf{B}_{38}$	=	$\left(\frac{1}{2} - x_{10}\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{10}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{10}\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{10}\right) c\,\hat{\mathbf{z}}$	(4*c*)	N III
$\mathbf{B}_{39}$	=	$-x_{10}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3$	=	$-x_{10} a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_{10} c\,\hat{\mathbf{z}}$	(4*c*)	N III
$\mathbf{B}_{40}$	=	$\left(\frac{1}{2} + x_{10}\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{10}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{10}\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{10}\right) c\,\hat{\mathbf{z}}$	(4*c*)	N III
$\mathbf{B}_{41}$	=	$x_{11}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$x_{11} a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_{11} c\,\hat{\mathbf{z}}$	(4*c*)	N IV
$\mathbf{B}_{42}$	=	$\left(\frac{1}{2} - x_{11}\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{11}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{11}\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{11}\right) c\,\hat{\mathbf{z}}$	(4*c*)	N IV
$\mathbf{B}_{43}$	=	$-x_{11}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3$	=	$-x_{11} a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_{11} c\,\hat{\mathbf{z}}$	(4*c*)	N IV
$\mathbf{B}_{44}$	=	$\left(\frac{1}{2} + x_{11}\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{11}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{11}\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{11}\right) c\,\hat{\mathbf{z}}$	(4*c*)	N IV
$\mathbf{B}_{45}$	=	$x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3$	=	$x_{12} a\,\hat{\mathbf{x}} + y_{12} b\,\hat{\mathbf{y}} + z_{12} c\,\hat{\mathbf{z}}$	(8*d*)	C V
$\mathbf{B}_{46}$	=	$\left(\frac{1}{2} - x_{12}\right)\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{12}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{12}\right) a\,\hat{\mathbf{x}} - y_{12} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{12}\right) c\,\hat{\mathbf{z}}$	(8*d*)	C V
$\mathbf{B}_{47}$	=	$-x_{12}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{12}\right)\mathbf{a}_2 - z_{12}\,\mathbf{a}_3$	=	$-x_{12} a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{12}\right) b\,\hat{\mathbf{y}} - z_{12} c\,\hat{\mathbf{z}}$	(8*d*)	C V
$\mathbf{B}_{48}$	=	$\left(\frac{1}{2} + x_{12}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{12}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{12}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{12}\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{12}\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{12}\right) c\,\hat{\mathbf{z}}$	(8*d*)	C V
$\mathbf{B}_{49}$	=	$-x_{12}\,\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3$	=	$-x_{12} a\,\hat{\mathbf{x}} - y_{12} b\,\hat{\mathbf{y}} - z_{12} c\,\hat{\mathbf{z}}$	(8*d*)	C V
$\mathbf{B}_{50}$	=	$\left(\frac{1}{2} + x_{12}\right)\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{12}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{12}\right) a\,\hat{\mathbf{x}} + y_{12} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{12}\right) c\,\hat{\mathbf{z}}$	(8*d*)	C V
$\mathbf{B}_{51}$	=	$x_{12}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{12}\right)\mathbf{a}_2 + z_{12}\,\mathbf{a}_3$	=	$x_{12} a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{12}\right) b\,\hat{\mathbf{y}} + z_{12} c\,\hat{\mathbf{z}}$	(8*d*)	C V
$\mathbf{B}_{52}$	=	$\left(\frac{1}{2} - x_{12}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{12}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{12}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{12}\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{12}\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{12}\right) c\,\hat{\mathbf{z}}$	(8*d*)	C V
$\mathbf{B}_{53}$	=	$x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3$	=	$x_{13} a\,\hat{\mathbf{x}} + y_{13} b\,\hat{\mathbf{y}} + z_{13} c\,\hat{\mathbf{z}}$	(8*d*)	C VI
$\mathbf{B}_{54}$	=	$\left(\frac{1}{2} - x_{13}\right)\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{13}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{13}\right) a\,\hat{\mathbf{x}} - y_{13} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{13}\right) c\,\hat{\mathbf{z}}$	(8*d*)	C VI
$\mathbf{B}_{55}$	=	$-x_{13}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{13}\right)\mathbf{a}_2 - z_{13}\,\mathbf{a}_3$	=	$-x_{13} a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{13}\right) b\,\hat{\mathbf{y}} - z_{13} c\,\hat{\mathbf{z}}$	(8*d*)	C VI
$\mathbf{B}_{56}$	=	$\left(\frac{1}{2} + x_{13}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{13}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{13}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{13}\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{13}\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{13}\right) c\,\hat{\mathbf{z}}$	(8*d*)	C VI
$\mathbf{B}_{57}$	=	$-x_{13}\,\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 - z_{13}\,\mathbf{a}_3$	=	$-x_{13} a\,\hat{\mathbf{x}} - y_{13} b\,\hat{\mathbf{y}} - z_{13} c\,\hat{\mathbf{z}}$	(8*d*)	C VI
$\mathbf{B}_{58}$	=	$\left(\frac{1}{2} + x_{13}\right)\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{13}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{13}\right) a\,\hat{\mathbf{x}} + y_{13} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{13}\right) c\,\hat{\mathbf{z}}$	(8*d*)	C VI
$\mathbf{B}_{59}$	=	$x_{13}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{13}\right)\mathbf{a}_2 + z_{13}\,\mathbf{a}_3$	=	$x_{13} a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{13}\right) b\,\hat{\mathbf{y}} + z_{13} c\,\hat{\mathbf{z}}$	(8*d*)	C VI
$\mathbf{B}_{60}$	=	$\left(\frac{1}{2} - x_{13}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{13}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{13}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{13}\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{13}\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{13}\right) c\,\hat{\mathbf{z}}$	(8*d*)	C VI
$\mathbf{B}_{61}$	=	$x_{14}\,\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3$	=	$x_{14} a\,\hat{\mathbf{x}} + y_{14} b\,\hat{\mathbf{y}} + z_{14} c\,\hat{\mathbf{z}}$	(8*d*)	H I
$\mathbf{B}_{62}$	=	$\left(\frac{1}{2} - x_{14}\right)\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{14}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{14}\right) a\,\hat{\mathbf{x}} - y_{14} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{14}\right) c\,\hat{\mathbf{z}}$	(8*d*)	H I
$\mathbf{B}_{63}$	=	$-x_{14}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{14}\right)\mathbf{a}_2 - z_{14}\,\mathbf{a}_3$	=	$-x_{14} a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{14}\right) b\,\hat{\mathbf{y}} - z_{14} c\,\hat{\mathbf{z}}$	(8*d*)	H I
$\mathbf{B}_{64}$	=	$\left(\frac{1}{2} + x_{14}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{14}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{14}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{14}\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{14}\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{14}\right) c\,\hat{\mathbf{z}}$	(8*d*)	H I
$\mathbf{B}_{65}$	=	$-x_{14}\,\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 - z_{14}\,\mathbf{a}_3$	=	$-x_{14} a\,\hat{\mathbf{x}} - y_{14} b\,\hat{\mathbf{y}} - z_{14} c\,\hat{\mathbf{z}}$	(8*d*)	H I
$\mathbf{B}_{66}$	=	$\left(\frac{1}{2} + x_{14}\right)\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{14}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{14}\right) a\,\hat{\mathbf{x}} + y_{14} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{14}\right) c\,\hat{\mathbf{z}}$	(8*d*)	H I

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\begin{array}{lcccccc}
\mathbf{B}_{67} & = & x_{14}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{14}\right)\mathbf{a}_2 + z_{14}\,\mathbf{a}_3 & = & x_{14}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{14}\right)b\,\hat{\mathbf{y}} + z_{14}c\,\hat{\mathbf{z}} & (8d) & \text{H I} \\
\mathbf{B}_{68} & = & \left(\frac{1}{2} - x_{14}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{14}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{14}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_{14}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{14}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{14}\right)c\,\hat{\mathbf{z}} & (8d) & \text{H I} \\
\mathbf{B}_{69} & = & x_{15}\,\mathbf{a}_1 + y_{15}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 & = & x_{15}a\,\hat{\mathbf{x}} + y_{15}b\,\hat{\mathbf{y}} + z_{15}c\,\hat{\mathbf{z}} & (8d) & \text{H II} \\
\mathbf{B}_{70} & = & \left(\frac{1}{2} - x_{15}\right)\mathbf{a}_1 - y_{15}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{15}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_{15}\right)a\,\hat{\mathbf{x}} - y_{15}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{15}\right)c\,\hat{\mathbf{z}} & (8d) & \text{H II} \\
\mathbf{B}_{71} & = & -x_{15}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{15}\right)\mathbf{a}_2 - z_{15}\,\mathbf{a}_3 & = & -x_{15}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{15}\right)b\,\hat{\mathbf{y}} - z_{15}c\,\hat{\mathbf{z}} & (8d) & \text{H II} \\
\mathbf{B}_{72} & = & \left(\frac{1}{2} + x_{15}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{15}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{15}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} + x_{15}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{15}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{15}\right)c\,\hat{\mathbf{z}} & (8d) & \text{H II} \\
\mathbf{B}_{73} & = & -x_{15}\,\mathbf{a}_1 - y_{15}\,\mathbf{a}_2 - z_{15}\,\mathbf{a}_3 & = & -x_{15}a\,\hat{\mathbf{x}} - y_{15}b\,\hat{\mathbf{y}} - z_{15}c\,\hat{\mathbf{z}} & (8d) & \text{H II} \\
\mathbf{B}_{74} & = & \left(\frac{1}{2} + x_{15}\right)\mathbf{a}_1 + y_{15}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{15}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} + x_{15}\right)a\,\hat{\mathbf{x}} + y_{15}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{15}\right)c\,\hat{\mathbf{z}} & (8d) & \text{H II} \\
\mathbf{B}_{75} & = & x_{15}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{15}\right)\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 & = & x_{15}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{15}\right)b\,\hat{\mathbf{y}} + z_{15}c\,\hat{\mathbf{z}} & (8d) & \text{H II} \\
\mathbf{B}_{76} & = & \left(\frac{1}{2} - x_{15}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{15}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{15}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_{15}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{15}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{15}\right)c\,\hat{\mathbf{z}} & (8d) & \text{H II} \\
\mathbf{B}_{77} & = & x_{16}\,\mathbf{a}_1 + y_{16}\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 & = & x_{16}a\,\hat{\mathbf{x}} + y_{16}b\,\hat{\mathbf{y}} + z_{16}c\,\hat{\mathbf{z}} & (8d) & \text{K III} \\
\mathbf{B}_{78} & = & \left(\frac{1}{2} - x_{16}\right)\mathbf{a}_1 - y_{16}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{16}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_{16}\right)a\,\hat{\mathbf{x}} - y_{16}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{16}\right)c\,\hat{\mathbf{z}} & (8d) & \text{K III} \\
\mathbf{B}_{79} & = & -x_{16}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{16}\right)\mathbf{a}_2 - z_{16}\,\mathbf{a}_3 & = & -x_{16}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{16}\right)b\,\hat{\mathbf{y}} - z_{16}c\,\hat{\mathbf{z}} & (8d) & \text{K III} \\
\mathbf{B}_{80} & = & \left(\frac{1}{2} + x_{16}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{16}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{16}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} + x_{16}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{16}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{16}\right)c\,\hat{\mathbf{z}} & (8d) & \text{K III} \\
\mathbf{B}_{81} & = & -x_{16}\,\mathbf{a}_1 - y_{16}\,\mathbf{a}_2 - z_{16}\,\mathbf{a}_3 & = & -x_{16}a\,\hat{\mathbf{x}} - y_{16}b\,\hat{\mathbf{y}} - z_{16}c\,\hat{\mathbf{z}} & (8d) & \text{K III} \\
\mathbf{B}_{82} & = & \left(\frac{1}{2} + x_{16}\right)\mathbf{a}_1 + y_{16}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{16}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} + x_{16}\right)a\,\hat{\mathbf{x}} + y_{16}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{16}\right)c\,\hat{\mathbf{z}} & (8d) & \text{K III} \\
\mathbf{B}_{83} & = & x_{16}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{16}\right)\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 & = & x_{16}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{16}\right)b\,\hat{\mathbf{y}} + z_{16}c\,\hat{\mathbf{z}} & (8d) & \text{K III} \\
\mathbf{B}_{84} & = & \left(\frac{1}{2} - x_{16}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{16}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{16}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_{16}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{16}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{16}\right)c\,\hat{\mathbf{z}} & (8d) & \text{K III} \\
\mathbf{B}_{85} & = & x_{17}\,\mathbf{a}_1 + y_{17}\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 & = & x_{17}a\,\hat{\mathbf{x}} + y_{17}b\,\hat{\mathbf{y}} + z_{17}c\,\hat{\mathbf{z}} & (8d) & \text{N V} \\
\mathbf{B}_{86} & = & \left(\frac{1}{2} - x_{17}\right)\mathbf{a}_1 - y_{17}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{17}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_{17}\right)a\,\hat{\mathbf{x}} - y_{17}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{17}\right)c\,\hat{\mathbf{z}} & (8d) & \text{N V} \\
\mathbf{B}_{87} & = & -x_{17}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{17}\right)\mathbf{a}_2 - z_{17}\,\mathbf{a}_3 & = & -x_{17}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{17}\right)b\,\hat{\mathbf{y}} - z_{17}c\,\hat{\mathbf{z}} & (8d) & \text{N V} \\
\mathbf{B}_{88} & = & \left(\frac{1}{2} + x_{17}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{17}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{17}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} + x_{17}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{17}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{17}\right)c\,\hat{\mathbf{z}} & (8d) & \text{N V} \\
\mathbf{B}_{89} & = & -x_{17}\,\mathbf{a}_1 - y_{17}\,\mathbf{a}_2 - z_{17}\,\mathbf{a}_3 & = & -x_{17}a\,\hat{\mathbf{x}} - y_{17}b\,\hat{\mathbf{y}} - z_{17}c\,\hat{\mathbf{z}} & (8d) & \text{N V} \\
\mathbf{B}_{90} & = & \left(\frac{1}{2} + x_{17}\right)\mathbf{a}_1 + y_{17}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{17}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} + x_{17}\right)a\,\hat{\mathbf{x}} + y_{17}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{17}\right)c\,\hat{\mathbf{z}} & (8d) & \text{N V} \\
\mathbf{B}_{91} & = & x_{17}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{17}\right)\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 & = & x_{17}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{17}\right)b\,\hat{\mathbf{y}} + z_{17}c\,\hat{\mathbf{z}} & (8d) & \text{N V} \\
\mathbf{B}_{92} & = & \left(\frac{1}{2} - x_{17}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{17}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{17}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_{17}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{17}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{17}\right)c\,\hat{\mathbf{z}} & (8d) & \text{N V} \\
\mathbf{B}_{93} & = & x_{18}\,\mathbf{a}_1 + y_{18}\,\mathbf{a}_2 + z_{18}\,\mathbf{a}_3 & = & x_{18}a\,\hat{\mathbf{x}} + y_{18}b\,\hat{\mathbf{y}} + z_{18}c\,\hat{\mathbf{z}} & (8d) & \text{N VI} \\
\mathbf{B}_{94} & = & \left(\frac{1}{2} - x_{18}\right)\mathbf{a}_1 - y_{18}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{18}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_{18}\right)a\,\hat{\mathbf{x}} - y_{18}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{18}\right)c\,\hat{\mathbf{z}} & (8d) & \text{N VI} \\
\mathbf{B}_{95} & = & -x_{18}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{18}\right)\mathbf{a}_2 - z_{18}\,\mathbf{a}_3 & = & -x_{18}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{18}\right)b\,\hat{\mathbf{y}} - z_{18}c\,\hat{\mathbf{z}} & (8d) & \text{N VI} \\
\mathbf{B}_{96} & = & \left(\frac{1}{2} + x_{18}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{18}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{18}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} + x_{18}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{18}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{18}\right)c\,\hat{\mathbf{z}} & (8d) & \text{N VI}
\end{array}
$$

$\mathbf{B}_{97}$	=	$-x_{18}\,\mathbf{a}_1 - y_{18}\,\mathbf{a}_2 - z_{18}\,\mathbf{a}_3$	=	$-x_{18}a\,\hat{\mathbf{x}} - y_{18}b\,\hat{\mathbf{y}} - z_{18}c\,\hat{\mathbf{z}}$	(8*d*)	N VI
$\mathbf{B}_{98}$	=	$\left(\frac{1}{2}+x_{18}\right)\mathbf{a}_1 + y_{18}\,\mathbf{a}_2 + \left(\frac{1}{2}-z_{18}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{18}\right)a\,\hat{\mathbf{x}} + y_{18}b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_{18}\right)c\,\hat{\mathbf{z}}$	(8*d*)	N VI
$\mathbf{B}_{99}$	=	$x_{18}\,\mathbf{a}_1 + \left(\frac{1}{2}-y_{18}\right)\mathbf{a}_2 + z_{18}\,\mathbf{a}_3$	=	$x_{18}a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_{18}\right)b\,\hat{\mathbf{y}} + z_{18}c\,\hat{\mathbf{z}}$	(8*d*)	N VI
$\mathbf{B}_{100}$	=	$\left(\frac{1}{2}-x_{18}\right)\mathbf{a}_1 + \left(\frac{1}{2}+y_{18}\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_{18}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{18}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_{18}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_{18}\right)c\,\hat{\mathbf{z}}$	(8*d*)	N VI
$\mathbf{B}_{101}$	=	$x_{19}\,\mathbf{a}_1 + y_{19}\,\mathbf{a}_2 + z_{19}\,\mathbf{a}_3$	=	$x_{19}a\,\hat{\mathbf{x}} + y_{19}b\,\hat{\mathbf{y}} + z_{19}c\,\hat{\mathbf{z}}$	(8*d*)	O
$\mathbf{B}_{102}$	=	$\left(\frac{1}{2}-x_{19}\right)\mathbf{a}_1 - y_{19}\,\mathbf{a}_2 + \left(\frac{1}{2}+z_{19}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{19}\right)a\,\hat{\mathbf{x}} - y_{19}b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_{19}\right)c\,\hat{\mathbf{z}}$	(8*d*)	O
$\mathbf{B}_{103}$	=	$-x_{19}\,\mathbf{a}_1 + \left(\frac{1}{2}+y_{19}\right)\mathbf{a}_2 - z_{19}\,\mathbf{a}_3$	=	$-x_{19}a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_{19}\right)b\,\hat{\mathbf{y}} - z_{19}c\,\hat{\mathbf{z}}$	(8*d*)	O
$\mathbf{B}_{104}$	=	$\left(\frac{1}{2}+x_{19}\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_{19}\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_{19}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{19}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_{19}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_{19}\right)c\,\hat{\mathbf{z}}$	(8*d*)	O
$\mathbf{B}_{105}$	=	$-x_{19}\,\mathbf{a}_1 - y_{19}\,\mathbf{a}_2 - z_{19}\,\mathbf{a}_3$	=	$-x_{19}a\,\hat{\mathbf{x}} - y_{19}b\,\hat{\mathbf{y}} - z_{19}c\,\hat{\mathbf{z}}$	(8*d*)	O
$\mathbf{B}_{106}$	=	$\left(\frac{1}{2}+x_{19}\right)\mathbf{a}_1 + y_{19}\,\mathbf{a}_2 + \left(\frac{1}{2}-z_{19}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{19}\right)a\,\hat{\mathbf{x}} + y_{19}b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_{19}\right)c\,\hat{\mathbf{z}}$	(8*d*)	O
$\mathbf{B}_{107}$	=	$x_{19}\,\mathbf{a}_1 + \left(\frac{1}{2}-y_{19}\right)\mathbf{a}_2 + z_{19}\,\mathbf{a}_3$	=	$x_{19}a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_{19}\right)b\,\hat{\mathbf{y}} + z_{19}c\,\hat{\mathbf{z}}$	(8*d*)	O
$\mathbf{B}_{108}$	=	$\left(\frac{1}{2}-x_{19}\right)\mathbf{a}_1 + \left(\frac{1}{2}+y_{19}\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_{19}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{19}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_{19}\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_{19}\right)c\,\hat{\mathbf{z}}$	(8*d*)	O

References:

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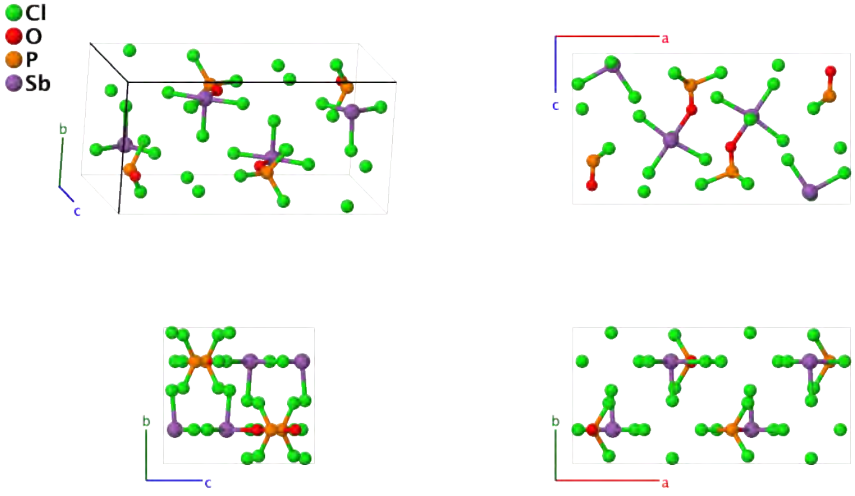
- J. L. Hoard and H. H. Nordsieck, *The Structure of Potassium Molybdocyanide Dihydrate. The Configuration of the Molybdenum Octocyanide Group*, J. Am. Chem. Soc. **61**, 2853–2863 (1939), doi:10.1021/ja01265a083.
- K. Herrmann, ed., *Strukturbericht Band VII 1939* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1943).

Geometry files:

- CIF: pp. 1641
- POSCAR: pp. 1641

$SbCl_5 \cdot POCl_3$ Structure: A8BCD_oP44_62_4c2d_c_c_c

http://aflow.org/prototype-encyclopedia/A8BCD_oP44_62_4c2d_c_c_c



Prototype	:	Cl_8OPSb
AFLOW prototype label	:	A8BCD_oP44_62_4c2d_c_c_c
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP44
Space group number	:	62
Space group symbol	:	*Pnma*
AFLOW prototype command	:	`aflow --proto=A8BCD_oP44_62_4c2d_c_c_c --params=`$a, b/a, c/a, x_1, z_1, x_2, z_2, x_3, z_3, x_4, z_4, x_5, z_5, x_6, z_6, x_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9$

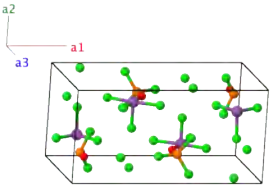
Other compounds with this structure

- $SbCl_5 \cdot POCl_3(CH_3)_3$

- (Lindqvist, 1959) give the data for this structure in setting *Pmnb* of space group #62. We used FINDSYM to exchange the *a*- and *b*-axes so that the structure is now in the standard *Pnma* orientation.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(4*c*)	Cl I
$\mathbf{B}_2$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	Cl I
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	(4*c*)	Cl I
$\mathbf{B}_4$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	Cl I
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*c*)	Cl II
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	(4*c*)	Cl II
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(4*c*)	Cl II
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right) c\,\hat{\mathbf{z}}$	(4*c*)	Cl II
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*c*)	Cl III
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	(4*c*)	Cl III
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(4*c*)	Cl III
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right) c\,\hat{\mathbf{z}}$	(4*c*)	Cl III
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(4*c*)	Cl IV
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	(4*c*)	Cl IV
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(4*c*)	Cl IV
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\,\hat{\mathbf{z}}$	(4*c*)	Cl IV
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(4*c*)	O
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	(4*c*)	O
$\mathbf{B}_{19}$	=	$-x_5\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(4*c*)	O
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right) c\,\hat{\mathbf{z}}$	(4*c*)	O
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(4*c*)	P
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	(4*c*)	P
$\mathbf{B}_{23}$	=	$-x_6\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(4*c*)	P
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right) c\,\hat{\mathbf{z}}$	(4*c*)	P
$\mathbf{B}_{25}$	=	$x_7\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(4*c*)	Sb
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} - x_7\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_7\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right) c\,\hat{\mathbf{z}}$	(4*c*)	Sb
$\mathbf{B}_{27}$	=	$-x_7\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	(4*c*)	Sb
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} + x_7\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_7\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right) c\,\hat{\mathbf{z}}$	(4*c*)	Sb
$\mathbf{B}_{29}$	=	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(8*d*)	Cl V
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2} - x_8\right)\mathbf{a}_1 - y_8\,\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_8\right) a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right) c\,\hat{\mathbf{z}}$	(8*d*)	Cl V
$\mathbf{B}_{31}$	=	$-x_8\,\mathbf{a}_1 + \left(\frac{1}{2} + y_8\right)\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_8\right) b\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(8*d*)	Cl V
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2} + x_8\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_8\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_8\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_8\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right) c\,\hat{\mathbf{z}}$	(8*d*)	Cl V
$\mathbf{B}_{33}$	=	$-x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(8*d*)	Cl V
$\mathbf{B}_{34}$	=	$\left(\frac{1}{2} + x_8\right)\mathbf{a}_1 + y_8\,\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_8\right) a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right) c\,\hat{\mathbf{z}}$	(8*d*)	Cl V
$\mathbf{B}_{35}$	=	$x_8\,\mathbf{a}_1 + \left(\frac{1}{2} - y_8\right)\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_8\right) b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(8*d*)	Cl V

$\mathbf{B}_{36}$	=	$\left(\frac{1}{2} - x_8\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_8\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_8\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right)c\,\hat{\mathbf{z}}$	(8*d*)	Cl V
$\mathbf{B}_{37}$	=	$x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$x_9a\,\hat{\mathbf{x}} + y_9b\,\hat{\mathbf{y}} + z_9c\,\hat{\mathbf{z}}$	(8*d*)	Cl VI
$\mathbf{B}_{38}$	=	$\left(\frac{1}{2} - x_9\right)\mathbf{a}_1 - y_9\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_9\right)a\,\hat{\mathbf{x}} - y_9b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right)c\,\hat{\mathbf{z}}$	(8*d*)	Cl VI
$\mathbf{B}_{39}$	=	$-x_9\,\mathbf{a}_1 + \left(\frac{1}{2} + y_9\right)\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$-x_9a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_9\right)b\,\hat{\mathbf{y}} - z_9c\,\hat{\mathbf{z}}$	(8*d*)	Cl VI
$\mathbf{B}_{40}$	=	$\left(\frac{1}{2} + x_9\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_9\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_9\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_9\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_9\right)c\,\hat{\mathbf{z}}$	(8*d*)	Cl VI
$\mathbf{B}_{41}$	=	$-x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$-x_9a\,\hat{\mathbf{x}} - y_9b\,\hat{\mathbf{y}} - z_9c\,\hat{\mathbf{z}}$	(8*d*)	Cl VI
$\mathbf{B}_{42}$	=	$\left(\frac{1}{2} + x_9\right)\mathbf{a}_1 + y_9\,\mathbf{a}_2 + \left(\frac{1}{2} - z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_9\right)a\,\hat{\mathbf{x}} + y_9b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_9\right)c\,\hat{\mathbf{z}}$	(8*d*)	Cl VI
$\mathbf{B}_{43}$	=	$x_9\,\mathbf{a}_1 + \left(\frac{1}{2} - y_9\right)\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$x_9a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_9\right)b\,\hat{\mathbf{y}} + z_9c\,\hat{\mathbf{z}}$	(8*d*)	Cl VI
$\mathbf{B}_{44}$	=	$\left(\frac{1}{2} - x_9\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_9\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_9\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_9\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right)c\,\hat{\mathbf{z}}$	(8*d*)	Cl VI

References:

- I. Lindqvist and C.-I. Brändén, *The Crystal Structure of $SbCl_5{\cdot}POCl_3$*, Acta Cryst. **12**, 642–645 (1959), doi:10.1107/S0365110X5900189X.

Found in:

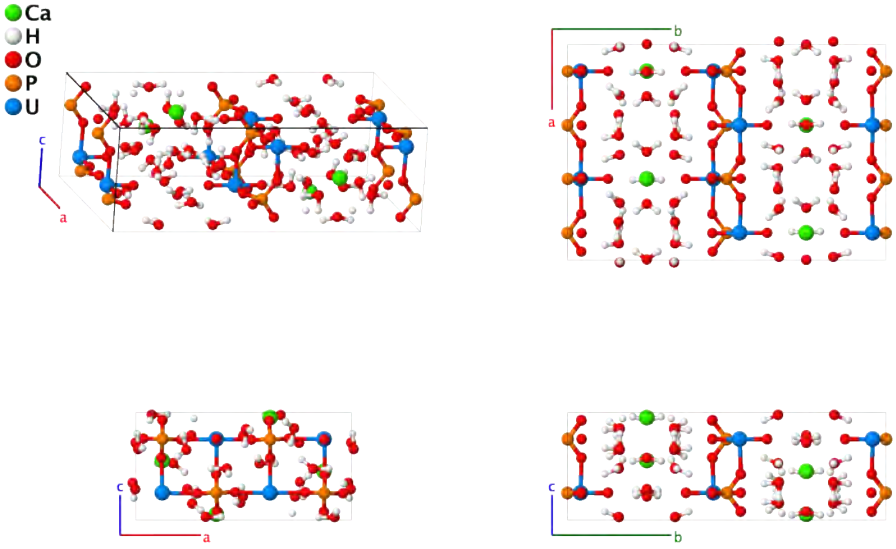
- C.-I. Brändén and I. Lindqvist, *The Crystal Structure of $SbCl_5.PO(CH_3)_3$*, Acta Chem. Scand. **15**, 167–174 (1961), doi:10.3891/acta.chem.scand.15-0167.

Geometry files:

- CIF: pp. 1642
- POSCAR: pp. 1642

Autunite $\{Ca[(UO_2)(PO_4)]_2(H_2O)_{11}\}$ Structure: AB22C23D2E2_oP200_62_c_11d_3c10d_d_d

http://aflow.org/prototype-encyclopedia/AB22C23D2E2_oP200_62_c_11d_3c10d_d_d

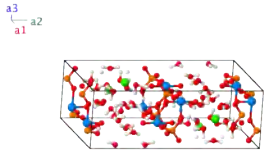


Prototype	:	$CaH_{22}O_{23}P_2U_2$
AFLOW prototype label	:	AB22C23D2E2_oP200_62_c_11d_3c10d_d_d
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP200
Space group number	:	62
Space group symbol	:	*Pnma*
AFLOW prototype command	:	`aflow --proto=AB22C23D2E2_oP200_62_c_11d_3c10d_d_d --params=`$a, b/a, c/a, x_1, z_1, x_2, z_2, x_3, z_3, x_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}, x_{17}, y_{17}, z_{17}, x_{18}, y_{18}, z_{18}, x_{19}, y_{19}, z_{19}, x_{20}, y_{20}, z_{20}, x_{21}, y_{21}, z_{21}, x_{22}, y_{22}, z_{22}, x_{23}, y_{23}, z_{23}, x_{24}, y_{24}, z_{24}, x_{25}, y_{25}, z_{25}, x_{26}, y_{26}, z_{26}, x_{27}, y_{27}, z_{27}$

- Autunite, $Ca(UO_2)_2(PO_4)_2{\cdot}nH_2O$, is found in three varieties: naturally occurring Autunite, with $n \gtrsim 10$, and meta-autunite (I), which is partially dehydrated, $6 \gtrsim n \gtrsim 10$. Further dehydration in the laboratory produces meta-autunite (II).
- The original determination of the autunite structure designated $H5_9$ by (Herrmann, 1941) was a tetragonal structure, and none of the positions of the oxygen atom or water molecules were determined. (Locock, 2003) find a pseudo-tetragonal ($a \approx 2c$) unit cell which doubles the size of the original cell. They were able to locate all of the atoms in the structure.
- (Locock, 2003) found the ($4c$) calcium site to be occupied only 86% of the time.

Simple Orthorhombic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= a\,\hat{\mathbf{x}}\\ \mathbf{a}_2 &= b\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	$x_1 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(4*c*)	Ca
$\mathbf{B}_2$	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	Ca
$\mathbf{B}_3$	$-x_1\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	$-x_1 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	(4*c*)	Ca
$\mathbf{B}_4$	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_1\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	Ca
$\mathbf{B}_5$	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	$x_2 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_6$	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_7$	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	$-x_2 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_8$	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right) c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_9$	$x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$x_3 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_{10}$	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_3\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_{11}$	$-x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	$-x_3 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_{12}$	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_3\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right) c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_{13}$	$x_4\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$x_4 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(4*c*)	O III
$\mathbf{B}_{14}$	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	(4*c*)	O III
$\mathbf{B}_{15}$	$-x_4\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	$-x_4 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(4*c*)	O III
$\mathbf{B}_{16}$	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\,\hat{\mathbf{z}}$	(4*c*)	O III
$\mathbf{B}_{17}$	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(8*d*)	H I
$\mathbf{B}_{18}$	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_5\right) a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	(8*d*)	H I
$\mathbf{B}_{19}$	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3$	$-x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right) b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(8*d*)	H I
$\mathbf{B}_{20}$	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right) c\,\hat{\mathbf{z}}$	(8*d*)	H I
$\mathbf{B}_{21}$	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(8*d*)	H I
$\mathbf{B}_{22}$	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + y_5\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right) c\,\hat{\mathbf{z}}$	(8*d*)	H I
$\mathbf{B}_{23}$	$x_5\,\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + z_5\,\mathbf{a}_3$	$x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right) b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(8*d*)	H I
$\mathbf{B}_{24}$	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_5\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	(8*d*)	H I
$\mathbf{B}_{25}$	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(8*d*)	H II
$\mathbf{B}_{26}$	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	(8*d*)	H II

$$\mathbf{B}_{27} = -x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3 = -x_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_6\right) b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H II}$$

$$\mathbf{B}_{28} = \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_6\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_6\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_6\right) c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H II}$$

$$\mathbf{B}_{29} = -x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3 = -x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H II}$$

$$\mathbf{B}_{30} = \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_1 + y_6\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_6\right) a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_6\right) c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H II}$$

$$\mathbf{B}_{31} = x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_6\right)\mathbf{a}_2 + z_6\,\mathbf{a}_3 = x_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_6\right) b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H II}$$

$$\mathbf{B}_{32} = \left(\tfrac{1}{2}-x_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_6\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_6\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_6\right) c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H II}$$

$$\mathbf{B}_{33} = x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 = x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H III}$$

$$\mathbf{B}_{34} = \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_7\right) a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_7\right) c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H III}$$

$$\mathbf{B}_{35} = -x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_7\right)\mathbf{a}_2 - z_7\,\mathbf{a}_3 = -x_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_7\right) b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H III}$$

$$\mathbf{B}_{36} = \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_7\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_7\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_7\right) c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H III}$$

$$\mathbf{B}_{37} = -x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3 = -x_7 a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H III}$$

$$\mathbf{B}_{38} = \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1 + y_7\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_7\right) a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_7\right) c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H III}$$

$$\mathbf{B}_{39} = x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_7\right)\mathbf{a}_2 + z_7\,\mathbf{a}_3 = x_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_7\right) b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H III}$$

$$\mathbf{B}_{40} = \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_7\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_7\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_7\right) c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H III}$$

$$\mathbf{B}_{41} = x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 = x_8 a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H IV}$$

$$\mathbf{B}_{42} = \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_1 - y_8\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_8\right) a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_8\right) c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H IV}$$

$$\mathbf{B}_{43} = -x_8\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_8\right)\mathbf{a}_2 - z_8\,\mathbf{a}_3 = -x_8 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_8\right) b\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H IV}$$

$$\mathbf{B}_{44} = \left(\tfrac{1}{2}+x_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_8\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_8\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_8\right) c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H IV}$$

$$\mathbf{B}_{45} = -x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3 = -x_8 a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H IV}$$

$$\mathbf{B}_{46} = \left(\tfrac{1}{2}+x_8\right)\mathbf{a}_1 + y_8\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_8\right) a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_8\right) c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H IV}$$

$$\mathbf{B}_{47} = x_8\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_8\right)\mathbf{a}_2 + z_8\,\mathbf{a}_3 = x_8 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_8\right) b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H IV}$$

$$\mathbf{B}_{48} = \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_8\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_8\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_8\right) c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H IV}$$

$$\mathbf{B}_{49} = x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 = x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H V}$$

$$\mathbf{B}_{50} = \left(\tfrac{1}{2}-x_9\right)\mathbf{a}_1 - y_9\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_9\right) a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_9\right) c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H V}$$

$$\mathbf{B}_{51} = -x_9\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_9\right)\mathbf{a}_2 - z_9\,\mathbf{a}_3 = -x_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_9\right) b\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H V}$$

$$\mathbf{B}_{52} = \left(\tfrac{1}{2}+x_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_9\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_9\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_9\right) c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H V}$$

$$\mathbf{B}_{53} = -x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3 = -x_9 a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H V}$$

$$\mathbf{B}_{54} = \left(\tfrac{1}{2}+x_9\right)\mathbf{a}_1 + y_9\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_9\right) a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_9\right) c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H V}$$

$$\mathbf{B}_{55} = x_9\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_9\right)\mathbf{a}_2 + z_9\,\mathbf{a}_3 = x_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_9\right) b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H V}$$

$$\mathbf{B}_{56} = \left(\tfrac{1}{2}-x_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_9\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_9\right) b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_9\right) c\,\hat{\mathbf{z}} \quad (8d) \quad \text{H V}$$

$$
\begin{array}{lcccccc}
\mathbf{B}_{57} & = & x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 & = & x_{10}a\,\hat{\mathbf{x}} + y_{10}b\,\hat{\mathbf{y}} + z_{10}c\,\hat{\mathbf{z}} & (8d) & \text{H VI} \\
\mathbf{B}_{58} & = & \left(\tfrac{1}{2} - x_{10}\right)\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{10}\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} - x_{10}\right)a\,\hat{\mathbf{x}} - y_{10}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{10}\right)c\,\hat{\mathbf{z}} & (8d) & \text{H VI} \\
\mathbf{B}_{59} & = & -x_{10}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{10}\right)\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 & = & -x_{10}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{10}\right)b\,\hat{\mathbf{y}} - z_{10}c\,\hat{\mathbf{z}} & (8d) & \text{H VI} \\
\mathbf{B}_{60} & = & \left(\tfrac{1}{2} + x_{10}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{10}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{10}\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} + x_{10}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{10}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{10}\right)c\,\hat{\mathbf{z}} & (8d) & \text{H VI} \\
\mathbf{B}_{61} & = & -x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 & = & -x_{10}a\,\hat{\mathbf{x}} - y_{10}b\,\hat{\mathbf{y}} - z_{10}c\,\hat{\mathbf{z}} & (8d) & \text{H VI} \\
\mathbf{B}_{62} & = & \left(\tfrac{1}{2} + x_{10}\right)\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{10}\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} + x_{10}\right)a\,\hat{\mathbf{x}} + y_{10}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{10}\right)c\,\hat{\mathbf{z}} & (8d) & \text{H VI} \\
\mathbf{B}_{63} & = & x_{10}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{10}\right)\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 & = & x_{10}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{10}\right)b\,\hat{\mathbf{y}} + z_{10}c\,\hat{\mathbf{z}} & (8d) & \text{H VI} \\
\mathbf{B}_{64} & = & \left(\tfrac{1}{2} - x_{10}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{10}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{10}\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} - x_{10}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{10}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{10}\right)c\,\hat{\mathbf{z}} & (8d) & \text{H VI} \\
\mathbf{B}_{65} & = & x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 & = & x_{11}a\,\hat{\mathbf{x}} + y_{11}b\,\hat{\mathbf{y}} + z_{11}c\,\hat{\mathbf{z}} & (8d) & \text{H VII} \\
\mathbf{B}_{66} & = & \left(\tfrac{1}{2} - x_{11}\right)\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{11}\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} - x_{11}\right)a\,\hat{\mathbf{x}} - y_{11}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{11}\right)c\,\hat{\mathbf{z}} & (8d) & \text{H VII} \\
\mathbf{B}_{67} & = & -x_{11}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{11}\right)\mathbf{a}_2 - z_{11}\,\mathbf{a}_3 & = & -x_{11}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{11}\right)b\,\hat{\mathbf{y}} - z_{11}c\,\hat{\mathbf{z}} & (8d) & \text{H VII} \\
\mathbf{B}_{68} & = & \left(\tfrac{1}{2} + x_{11}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{11}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{11}\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} + x_{11}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{11}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{11}\right)c\,\hat{\mathbf{z}} & (8d) & \text{H VII} \\
\mathbf{B}_{69} & = & -x_{11}\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3 & = & -x_{11}a\,\hat{\mathbf{x}} - y_{11}b\,\hat{\mathbf{y}} - z_{11}c\,\hat{\mathbf{z}} & (8d) & \text{H VII} \\
\mathbf{B}_{70} & = & \left(\tfrac{1}{2} + x_{11}\right)\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{11}\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} + x_{11}\right)a\,\hat{\mathbf{x}} + y_{11}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{11}\right)c\,\hat{\mathbf{z}} & (8d) & \text{H VII} \\
\mathbf{B}_{71} & = & x_{11}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{11}\right)\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 & = & x_{11}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{11}\right)b\,\hat{\mathbf{y}} + z_{11}c\,\hat{\mathbf{z}} & (8d) & \text{H VII} \\
\mathbf{B}_{72} & = & \left(\tfrac{1}{2} - x_{11}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{11}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{11}\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} - x_{11}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{11}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{11}\right)c\,\hat{\mathbf{z}} & (8d) & \text{H VII} \\
\mathbf{B}_{73} & = & x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 & = & x_{12}a\,\hat{\mathbf{x}} + y_{12}b\,\hat{\mathbf{y}} + z_{12}c\,\hat{\mathbf{z}} & (8d) & \text{H VIII} \\
\mathbf{B}_{74} & = & \left(\tfrac{1}{2} - x_{12}\right)\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{12}\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} - x_{12}\right)a\,\hat{\mathbf{x}} - y_{12}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{12}\right)c\,\hat{\mathbf{z}} & (8d) & \text{H VIII} \\
\mathbf{B}_{75} & = & -x_{12}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{12}\right)\mathbf{a}_2 - z_{12}\,\mathbf{a}_3 & = & -x_{12}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{12}\right)b\,\hat{\mathbf{y}} - z_{12}c\,\hat{\mathbf{z}} & (8d) & \text{H VIII} \\
\mathbf{B}_{76} & = & \left(\tfrac{1}{2} + x_{12}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{12}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{12}\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} + x_{12}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{12}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{12}\right)c\,\hat{\mathbf{z}} & (8d) & \text{H VIII} \\
\mathbf{B}_{77} & = & -x_{12}\,\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 - z_{12}\,\mathbf{a}_3 & = & -x_{12}a\,\hat{\mathbf{x}} - y_{12}b\,\hat{\mathbf{y}} - z_{12}c\,\hat{\mathbf{z}} & (8d) & \text{H VIII} \\
\mathbf{B}_{78} & = & \left(\tfrac{1}{2} + x_{12}\right)\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{12}\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} + x_{12}\right)a\,\hat{\mathbf{x}} + y_{12}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{12}\right)c\,\hat{\mathbf{z}} & (8d) & \text{H VIII} \\
\mathbf{B}_{79} & = & x_{12}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{12}\right)\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 & = & x_{12}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{12}\right)b\,\hat{\mathbf{y}} + z_{12}c\,\hat{\mathbf{z}} & (8d) & \text{H VIII} \\
\mathbf{B}_{80} & = & \left(\tfrac{1}{2} - x_{12}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{12}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{12}\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} - x_{12}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{12}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{12}\right)c\,\hat{\mathbf{z}} & (8d) & \text{H VIII} \\
\mathbf{B}_{81} & = & x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3 & = & x_{13}a\,\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} + z_{13}c\,\hat{\mathbf{z}} & (8d) & \text{H IX} \\
\mathbf{B}_{82} & = & \left(\tfrac{1}{2} - x_{13}\right)\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{13}\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} - x_{13}\right)a\,\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{13}\right)c\,\hat{\mathbf{z}} & (8d) & \text{H IX} \\
\mathbf{B}_{83} & = & -x_{13}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{13}\right)\mathbf{a}_2 - z_{13}\,\mathbf{a}_3 & = & -x_{13}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{13}\right)b\,\hat{\mathbf{y}} - z_{13}c\,\hat{\mathbf{z}} & (8d) & \text{H IX} \\
\mathbf{B}_{84} & = & \left(\tfrac{1}{2} + x_{13}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{13}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{13}\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} + x_{13}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{13}\right)b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{13}\right)c\,\hat{\mathbf{z}} & (8d) & \text{H IX} \\
\mathbf{B}_{85} & = & -x_{13}\,\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 - z_{13}\,\mathbf{a}_3 & = & -x_{13}a\,\hat{\mathbf{x}} - y_{13}b\,\hat{\mathbf{y}} - z_{13}c\,\hat{\mathbf{z}} & (8d) & \text{H IX} \\
\mathbf{B}_{86} & = & \left(\tfrac{1}{2} + x_{13}\right)\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{13}\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} + x_{13}\right)a\,\hat{\mathbf{x}} + y_{13}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{13}\right)c\,\hat{\mathbf{z}} & (8d) & \text{H IX} \\
\mathbf{B}_{87} & = & x_{13}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{13}\right)\mathbf{a}_2 + z_{13}\,\mathbf{a}_3 & = & x_{13}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{13}\right)b\,\hat{\mathbf{y}} + z_{13}c\,\hat{\mathbf{z}} & (8d) & \text{H IX}
\end{array}
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\begin{array}{lcccccc}
\mathbf{B}_{88} & = & \left(\frac{1}{2}-x_{13}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{13}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{13}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{13}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{13}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{13}\right)c\,\hat{\mathbf{z}} & (8d) & \text{H IX}\\
\mathbf{B}_{89} & = & x_{14}\,\mathbf{a}_1+y_{14}\,\mathbf{a}_2+z_{14}\,\mathbf{a}_3 & = & x_{14}a\,\hat{\mathbf{x}}+y_{14}b\,\hat{\mathbf{y}}+z_{14}c\,\hat{\mathbf{z}} & (8d) & \text{H X}\\
\mathbf{B}_{90} & = & \left(\frac{1}{2}-x_{14}\right)\mathbf{a}_1-y_{14}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{14}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{14}\right)a\,\hat{\mathbf{x}}-y_{14}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{14}\right)c\,\hat{\mathbf{z}} & (8d) & \text{H X}\\
\mathbf{B}_{91} & = & -x_{14}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{14}\right)\mathbf{a}_2-z_{14}\,\mathbf{a}_3 & = & -x_{14}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{14}\right)b\,\hat{\mathbf{y}}-z_{14}c\,\hat{\mathbf{z}} & (8d) & \text{H X}\\
\mathbf{B}_{92} & = & \left(\frac{1}{2}+x_{14}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{14}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{14}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{14}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{14}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{14}\right)c\,\hat{\mathbf{z}} & (8d) & \text{H X}\\
\mathbf{B}_{93} & = & -x_{14}\,\mathbf{a}_1-y_{14}\,\mathbf{a}_2-z_{14}\,\mathbf{a}_3 & = & -x_{14}a\,\hat{\mathbf{x}}-y_{14}b\,\hat{\mathbf{y}}-z_{14}c\,\hat{\mathbf{z}} & (8d) & \text{H X}\\
\mathbf{B}_{94} & = & \left(\frac{1}{2}+x_{14}\right)\mathbf{a}_1+y_{14}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{14}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{14}\right)a\,\hat{\mathbf{x}}+y_{14}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{14}\right)c\,\hat{\mathbf{z}} & (8d) & \text{H X}\\
\mathbf{B}_{95} & = & x_{14}\,\mathbf{a}_1+\left(\frac{1}{2}-y_{14}\right)\mathbf{a}_2+z_{14}\,\mathbf{a}_3 & = & x_{14}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{14}\right)b\,\hat{\mathbf{y}}+z_{14}c\,\hat{\mathbf{z}} & (8d) & \text{H X}\\
\mathbf{B}_{96} & = & \left(\frac{1}{2}-x_{14}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{14}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{14}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{14}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{14}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{14}\right)c\,\hat{\mathbf{z}} & (8d) & \text{H X}\\
\mathbf{B}_{97} & = & x_{15}\,\mathbf{a}_1+y_{15}\,\mathbf{a}_2+z_{15}\,\mathbf{a}_3 & = & x_{15}a\,\hat{\mathbf{x}}+y_{15}b\,\hat{\mathbf{y}}+z_{15}c\,\hat{\mathbf{z}} & (8d) & \text{H XI}\\
\mathbf{B}_{98} & = & \left(\frac{1}{2}-x_{15}\right)\mathbf{a}_1-y_{15}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{15}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{15}\right)a\,\hat{\mathbf{x}}-y_{15}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{15}\right)c\,\hat{\mathbf{z}} & (8d) & \text{H XI}\\
\mathbf{B}_{99} & = & -x_{15}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{15}\right)\mathbf{a}_2-z_{15}\,\mathbf{a}_3 & = & -x_{15}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{15}\right)b\,\hat{\mathbf{y}}-z_{15}c\,\hat{\mathbf{z}} & (8d) & \text{H XI}\\
\mathbf{B}_{100} & = & \left(\frac{1}{2}+x_{15}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{15}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{15}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{15}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{15}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{15}\right)c\,\hat{\mathbf{z}} & (8d) & \text{H XI}\\
\mathbf{B}_{101} & = & -x_{15}\,\mathbf{a}_1-y_{15}\,\mathbf{a}_2-z_{15}\,\mathbf{a}_3 & = & -x_{15}a\,\hat{\mathbf{x}}-y_{15}b\,\hat{\mathbf{y}}-z_{15}c\,\hat{\mathbf{z}} & (8d) & \text{H XI}\\
\mathbf{B}_{102} & = & \left(\frac{1}{2}+x_{15}\right)\mathbf{a}_1+y_{15}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{15}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{15}\right)a\,\hat{\mathbf{x}}+y_{15}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{15}\right)c\,\hat{\mathbf{z}} & (8d) & \text{H XI}\\
\mathbf{B}_{103} & = & x_{15}\,\mathbf{a}_1+\left(\frac{1}{2}-y_{15}\right)\mathbf{a}_2+z_{15}\,\mathbf{a}_3 & = & x_{15}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{15}\right)b\,\hat{\mathbf{y}}+z_{15}c\,\hat{\mathbf{z}} & (8d) & \text{H XI}\\
\mathbf{B}_{104} & = & \left(\frac{1}{2}-x_{15}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{15}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{15}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{15}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{15}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{15}\right)c\,\hat{\mathbf{z}} & (8d) & \text{H XI}\\
\mathbf{B}_{105} & = & x_{16}\,\mathbf{a}_1+y_{16}\,\mathbf{a}_2+z_{16}\,\mathbf{a}_3 & = & x_{16}a\,\hat{\mathbf{x}}+y_{16}b\,\hat{\mathbf{y}}+z_{16}c\,\hat{\mathbf{z}} & (8d) & \text{O IV}\\
\mathbf{B}_{106} & = & \left(\frac{1}{2}-x_{16}\right)\mathbf{a}_1-y_{16}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{16}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{16}\right)a\,\hat{\mathbf{x}}-y_{16}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{16}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O IV}\\
\mathbf{B}_{107} & = & -x_{16}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{16}\right)\mathbf{a}_2-z_{16}\,\mathbf{a}_3 & = & -x_{16}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{16}\right)b\,\hat{\mathbf{y}}-z_{16}c\,\hat{\mathbf{z}} & (8d) & \text{O IV}\\
\mathbf{B}_{108} & = & \left(\frac{1}{2}+x_{16}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{16}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{16}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{16}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{16}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{16}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O IV}\\
\mathbf{B}_{109} & = & -x_{16}\,\mathbf{a}_1-y_{16}\,\mathbf{a}_2-z_{16}\,\mathbf{a}_3 & = & -x_{16}a\,\hat{\mathbf{x}}-y_{16}b\,\hat{\mathbf{y}}-z_{16}c\,\hat{\mathbf{z}} & (8d) & \text{O IV}\\
\mathbf{B}_{110} & = & \left(\frac{1}{2}+x_{16}\right)\mathbf{a}_1+y_{16}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{16}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{16}\right)a\,\hat{\mathbf{x}}+y_{16}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{16}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O IV}\\
\mathbf{B}_{111} & = & x_{16}\,\mathbf{a}_1+\left(\frac{1}{2}-y_{16}\right)\mathbf{a}_2+z_{16}\,\mathbf{a}_3 & = & x_{16}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{16}\right)b\,\hat{\mathbf{y}}+z_{16}c\,\hat{\mathbf{z}} & (8d) & \text{O IV}\\
\mathbf{B}_{112} & = & \left(\frac{1}{2}-x_{16}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{16}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{16}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{16}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{16}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{16}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O IV}\\
\mathbf{B}_{113} & = & x_{17}\,\mathbf{a}_1+y_{17}\,\mathbf{a}_2+z_{17}\,\mathbf{a}_3 & = & x_{17}a\,\hat{\mathbf{x}}+y_{17}b\,\hat{\mathbf{y}}+z_{17}c\,\hat{\mathbf{z}} & (8d) & \text{O V}\\
\mathbf{B}_{114} & = & \left(\frac{1}{2}-x_{17}\right)\mathbf{a}_1-y_{17}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{17}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{17}\right)a\,\hat{\mathbf{x}}-y_{17}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{17}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O V}\\
\mathbf{B}_{115} & = & -x_{17}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{17}\right)\mathbf{a}_2-z_{17}\,\mathbf{a}_3 & = & -x_{17}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{17}\right)b\,\hat{\mathbf{y}}-z_{17}c\,\hat{\mathbf{z}} & (8d) & \text{O V}\\
\mathbf{B}_{116} & = & \left(\frac{1}{2}+x_{17}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{17}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{17}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{17}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{17}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{17}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O V}\\
\mathbf{B}_{117} & = & -x_{17}\,\mathbf{a}_1-y_{17}\,\mathbf{a}_2-z_{17}\,\mathbf{a}_3 & = & -x_{17}a\,\hat{\mathbf{x}}-y_{17}b\,\hat{\mathbf{y}}-z_{17}c\,\hat{\mathbf{z}} & (8d) & \text{O V}
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\begin{array}{lcccccc}
\mathbf{B}_{118} & = & \left(\frac{1}{2}+x_{17}\right)\mathbf{a}_1+y_{17}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{17}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{17}\right)a\,\hat{\mathbf{x}}+y_{17}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{17}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O V}\\
\mathbf{B}_{119} & = & x_{17}\,\mathbf{a}_1+\left(\frac{1}{2}-y_{17}\right)\mathbf{a}_2+z_{17}\,\mathbf{a}_3 & = & x_{17}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{17}\right)b\,\hat{\mathbf{y}}+z_{17}c\,\hat{\mathbf{z}} & (8d) & \text{O V}\\
\mathbf{B}_{120} & = & \left(\frac{1}{2}-x_{17}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{17}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{17}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{17}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{17}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{17}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O V}\\
\mathbf{B}_{121} & = & x_{18}\,\mathbf{a}_1+y_{18}\,\mathbf{a}_2+z_{18}\,\mathbf{a}_3 & = & x_{18}a\,\hat{\mathbf{x}}+y_{18}b\,\hat{\mathbf{y}}+z_{18}c\,\hat{\mathbf{z}} & (8d) & \text{O VI}\\
\mathbf{B}_{122} & = & \left(\frac{1}{2}-x_{18}\right)\mathbf{a}_1-y_{18}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{18}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{18}\right)a\,\hat{\mathbf{x}}-y_{18}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{18}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O VI}\\
\mathbf{B}_{123} & = & -x_{18}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{18}\right)\mathbf{a}_2-z_{18}\,\mathbf{a}_3 & = & -x_{18}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{18}\right)b\,\hat{\mathbf{y}}-z_{18}c\,\hat{\mathbf{z}} & (8d) & \text{O VI}\\
\mathbf{B}_{124} & = & \left(\frac{1}{2}+x_{18}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{18}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{18}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{18}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{18}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{18}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O VI}\\
\mathbf{B}_{125} & = & -x_{18}\,\mathbf{a}_1-y_{18}\,\mathbf{a}_2-z_{18}\,\mathbf{a}_3 & = & -x_{18}a\,\hat{\mathbf{x}}-y_{18}b\,\hat{\mathbf{y}}-z_{18}c\,\hat{\mathbf{z}} & (8d) & \text{O VI}\\
\mathbf{B}_{126} & = & \left(\frac{1}{2}+x_{18}\right)\mathbf{a}_1+y_{18}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{18}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{18}\right)a\,\hat{\mathbf{x}}+y_{18}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{18}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O VI}\\
\mathbf{B}_{127} & = & x_{18}\,\mathbf{a}_1+\left(\frac{1}{2}-y_{18}\right)\mathbf{a}_2+z_{18}\,\mathbf{a}_3 & = & x_{18}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{18}\right)b\,\hat{\mathbf{y}}+z_{18}c\,\hat{\mathbf{z}} & (8d) & \text{O VI}\\
\mathbf{B}_{128} & = & \left(\frac{1}{2}-x_{18}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{18}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{18}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{18}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{18}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{18}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O VI}\\
\mathbf{B}_{129} & = & x_{19}\,\mathbf{a}_1+y_{19}\,\mathbf{a}_2+z_{19}\,\mathbf{a}_3 & = & x_{19}a\,\hat{\mathbf{x}}+y_{19}b\,\hat{\mathbf{y}}+z_{19}c\,\hat{\mathbf{z}} & (8d) & \text{O VII}\\
\mathbf{B}_{130} & = & \left(\frac{1}{2}-x_{19}\right)\mathbf{a}_1-y_{19}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{19}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{19}\right)a\,\hat{\mathbf{x}}-y_{19}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{19}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O VII}\\
\mathbf{B}_{131} & = & -x_{19}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{19}\right)\mathbf{a}_2-z_{19}\,\mathbf{a}_3 & = & -x_{19}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{19}\right)b\,\hat{\mathbf{y}}-z_{19}c\,\hat{\mathbf{z}} & (8d) & \text{O VII}\\
\mathbf{B}_{132} & = & \left(\frac{1}{2}+x_{19}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{19}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{19}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{19}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{19}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{19}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O VII}\\
\mathbf{B}_{133} & = & -x_{19}\,\mathbf{a}_1-y_{19}\,\mathbf{a}_2-z_{19}\,\mathbf{a}_3 & = & -x_{19}a\,\hat{\mathbf{x}}-y_{19}b\,\hat{\mathbf{y}}-z_{19}c\,\hat{\mathbf{z}} & (8d) & \text{O VII}\\
\mathbf{B}_{134} & = & \left(\frac{1}{2}+x_{19}\right)\mathbf{a}_1+y_{19}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{19}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{19}\right)a\,\hat{\mathbf{x}}+y_{19}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{19}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O VII}\\
\mathbf{B}_{135} & = & x_{19}\,\mathbf{a}_1+\left(\frac{1}{2}-y_{19}\right)\mathbf{a}_2+z_{19}\,\mathbf{a}_3 & = & x_{19}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{19}\right)b\,\hat{\mathbf{y}}+z_{19}c\,\hat{\mathbf{z}} & (8d) & \text{O VII}\\
\mathbf{B}_{136} & = & \left(\frac{1}{2}-x_{19}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{19}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{19}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{19}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{19}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{19}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O VII}\\
\mathbf{B}_{137} & = & x_{20}\,\mathbf{a}_1+y_{20}\,\mathbf{a}_2+z_{20}\,\mathbf{a}_3 & = & x_{20}a\,\hat{\mathbf{x}}+y_{20}b\,\hat{\mathbf{y}}+z_{20}c\,\hat{\mathbf{z}} & (8d) & \text{O VIII}\\
\mathbf{B}_{138} & = & \left(\frac{1}{2}-x_{20}\right)\mathbf{a}_1-y_{20}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{20}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{20}\right)a\,\hat{\mathbf{x}}-y_{20}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{20}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O VIII}\\
\mathbf{B}_{139} & = & -x_{20}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{20}\right)\mathbf{a}_2-z_{20}\,\mathbf{a}_3 & = & -x_{20}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{20}\right)b\,\hat{\mathbf{y}}-z_{20}c\,\hat{\mathbf{z}} & (8d) & \text{O VIII}\\
\mathbf{B}_{140} & = & \left(\frac{1}{2}+x_{20}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{20}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{20}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{20}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{20}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{20}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O VIII}\\
\mathbf{B}_{141} & = & -x_{20}\,\mathbf{a}_1-y_{20}\,\mathbf{a}_2-z_{20}\,\mathbf{a}_3 & = & -x_{20}a\,\hat{\mathbf{x}}-y_{20}b\,\hat{\mathbf{y}}-z_{20}c\,\hat{\mathbf{z}} & (8d) & \text{O VIII}\\
\mathbf{B}_{142} & = & \left(\frac{1}{2}+x_{20}\right)\mathbf{a}_1+y_{20}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{20}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{20}\right)a\,\hat{\mathbf{x}}+y_{20}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{20}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O VIII}\\
\mathbf{B}_{143} & = & x_{20}\,\mathbf{a}_1+\left(\frac{1}{2}-y_{20}\right)\mathbf{a}_2+z_{20}\,\mathbf{a}_3 & = & x_{20}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{20}\right)b\,\hat{\mathbf{y}}+z_{20}c\,\hat{\mathbf{z}} & (8d) & \text{O VIII}\\
\mathbf{B}_{144} & = & \left(\frac{1}{2}-x_{20}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{20}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{20}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{20}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{20}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{20}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O VIII}\\
\mathbf{B}_{145} & = & x_{21}\,\mathbf{a}_1+y_{21}\,\mathbf{a}_2+z_{21}\,\mathbf{a}_3 & = & x_{21}a\,\hat{\mathbf{x}}+y_{21}b\,\hat{\mathbf{y}}+z_{21}c\,\hat{\mathbf{z}} & (8d) & \text{O IX}\\
\mathbf{B}_{146} & = & \left(\frac{1}{2}-x_{21}\right)\mathbf{a}_1-y_{21}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{21}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{21}\right)a\,\hat{\mathbf{x}}-y_{21}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{21}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O IX}\\
\mathbf{B}_{147} & = & -x_{21}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{21}\right)\mathbf{a}_2-z_{21}\,\mathbf{a}_3 & = & -x_{21}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{21}\right)b\,\hat{\mathbf{y}}-z_{21}c\,\hat{\mathbf{z}} & (8d) & \text{O IX}
\end{array}
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\begin{array}{lcccll}
\mathbf{B}_{148} & = & \left(\frac{1}{2}+x_{21}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{21}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{21}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{21}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{21}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{21}\right)c\,\hat{\mathbf{z}} & (8d)\ \text{O IX}\\
\mathbf{B}_{149} & = & -x_{21}\,\mathbf{a}_1-y_{21}\,\mathbf{a}_2-z_{21}\,\mathbf{a}_3 & = & -x_{21}a\,\hat{\mathbf{x}}-y_{21}b\,\hat{\mathbf{y}}-z_{21}c\,\hat{\mathbf{z}} & (8d)\ \text{O IX}\\
\mathbf{B}_{150} & = & \left(\frac{1}{2}+x_{21}\right)\mathbf{a}_1+y_{21}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{21}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{21}\right)a\,\hat{\mathbf{x}}+y_{21}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{21}\right)c\,\hat{\mathbf{z}} & (8d)\ \text{O IX}\\
\mathbf{B}_{151} & = & x_{21}\,\mathbf{a}_1+\left(\frac{1}{2}-y_{21}\right)\mathbf{a}_2+z_{21}\,\mathbf{a}_3 & = & x_{21}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{21}\right)b\,\hat{\mathbf{y}}+z_{21}c\,\hat{\mathbf{z}} & (8d)\ \text{O IX}\\
\mathbf{B}_{152} & = & \left(\frac{1}{2}-x_{21}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{21}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{21}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{21}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{21}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{21}\right)c\,\hat{\mathbf{z}} & (8d)\ \text{O IX}\\
\mathbf{B}_{153} & = & x_{22}\,\mathbf{a}_1+y_{22}\,\mathbf{a}_2+z_{22}\,\mathbf{a}_3 & = & x_{22}a\,\hat{\mathbf{x}}+y_{22}b\,\hat{\mathbf{y}}+z_{22}c\,\hat{\mathbf{z}} & (8d)\ \text{O X}\\
\mathbf{B}_{154} & = & \left(\frac{1}{2}-x_{22}\right)\mathbf{a}_1-y_{22}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{22}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{22}\right)a\,\hat{\mathbf{x}}-y_{22}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{22}\right)c\,\hat{\mathbf{z}} & (8d)\ \text{O X}\\
\mathbf{B}_{155} & = & -x_{22}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{22}\right)\mathbf{a}_2-z_{22}\,\mathbf{a}_3 & = & -x_{22}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{22}\right)b\,\hat{\mathbf{y}}-z_{22}c\,\hat{\mathbf{z}} & (8d)\ \text{O X}\\
\mathbf{B}_{156} & = & \left(\frac{1}{2}+x_{22}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{22}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{22}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{22}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{22}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{22}\right)c\,\hat{\mathbf{z}} & (8d)\ \text{O X}\\
\mathbf{B}_{157} & = & -x_{22}\,\mathbf{a}_1-y_{22}\,\mathbf{a}_2-z_{22}\,\mathbf{a}_3 & = & -x_{22}a\,\hat{\mathbf{x}}-y_{22}b\,\hat{\mathbf{y}}-z_{22}c\,\hat{\mathbf{z}} & (8d)\ \text{O X}\\
\mathbf{B}_{158} & = & \left(\frac{1}{2}+x_{22}\right)\mathbf{a}_1+y_{22}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{22}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{22}\right)a\,\hat{\mathbf{x}}+y_{22}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{22}\right)c\,\hat{\mathbf{z}} & (8d)\ \text{O X}\\
\mathbf{B}_{159} & = & x_{22}\,\mathbf{a}_1+\left(\frac{1}{2}-y_{22}\right)\mathbf{a}_2+z_{22}\,\mathbf{a}_3 & = & x_{22}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{22}\right)b\,\hat{\mathbf{y}}+z_{22}c\,\hat{\mathbf{z}} & (8d)\ \text{O X}\\
\mathbf{B}_{160} & = & \left(\frac{1}{2}-x_{22}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{22}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{22}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{22}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{22}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{22}\right)c\,\hat{\mathbf{z}} & (8d)\ \text{O X}\\
\mathbf{B}_{161} & = & x_{23}\,\mathbf{a}_1+y_{23}\,\mathbf{a}_2+z_{23}\,\mathbf{a}_3 & = & x_{23}a\,\hat{\mathbf{x}}+y_{23}b\,\hat{\mathbf{y}}+z_{23}c\,\hat{\mathbf{z}} & (8d)\ \text{O XI}\\
\mathbf{B}_{162} & = & \left(\frac{1}{2}-x_{23}\right)\mathbf{a}_1-y_{23}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{23}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{23}\right)a\,\hat{\mathbf{x}}-y_{23}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{23}\right)c\,\hat{\mathbf{z}} & (8d)\ \text{O XI}\\
\mathbf{B}_{163} & = & -x_{23}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{23}\right)\mathbf{a}_2-z_{23}\,\mathbf{a}_3 & = & -x_{23}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{23}\right)b\,\hat{\mathbf{y}}-z_{23}c\,\hat{\mathbf{z}} & (8d)\ \text{O XI}\\
\mathbf{B}_{164} & = & \left(\frac{1}{2}+x_{23}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{23}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{23}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{23}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{23}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{23}\right)c\,\hat{\mathbf{z}} & (8d)\ \text{O XI}\\
\mathbf{B}_{165} & = & -x_{23}\,\mathbf{a}_1-y_{23}\,\mathbf{a}_2-z_{23}\,\mathbf{a}_3 & = & -x_{23}a\,\hat{\mathbf{x}}-y_{23}b\,\hat{\mathbf{y}}-z_{23}c\,\hat{\mathbf{z}} & (8d)\ \text{O XI}\\
\mathbf{B}_{166} & = & \left(\frac{1}{2}+x_{23}\right)\mathbf{a}_1+y_{23}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{23}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{23}\right)a\,\hat{\mathbf{x}}+y_{23}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{23}\right)c\,\hat{\mathbf{z}} & (8d)\ \text{O XI}\\
\mathbf{B}_{167} & = & x_{23}\,\mathbf{a}_1+\left(\frac{1}{2}-y_{23}\right)\mathbf{a}_2+z_{23}\,\mathbf{a}_3 & = & x_{23}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{23}\right)b\,\hat{\mathbf{y}}+z_{23}c\,\hat{\mathbf{z}} & (8d)\ \text{O XI}\\
\mathbf{B}_{168} & = & \left(\frac{1}{2}-x_{23}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{23}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{23}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{23}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{23}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{23}\right)c\,\hat{\mathbf{z}} & (8d)\ \text{O XI}\\
\mathbf{B}_{169} & = & x_{24}\,\mathbf{a}_1+y_{24}\,\mathbf{a}_2+z_{24}\,\mathbf{a}_3 & = & x_{24}a\,\hat{\mathbf{x}}+y_{24}b\,\hat{\mathbf{y}}+z_{24}c\,\hat{\mathbf{z}} & (8d)\ \text{O XII}\\
\mathbf{B}_{170} & = & \left(\frac{1}{2}-x_{24}\right)\mathbf{a}_1-y_{24}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{24}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{24}\right)a\,\hat{\mathbf{x}}-y_{24}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{24}\right)c\,\hat{\mathbf{z}} & (8d)\ \text{O XII}\\
\mathbf{B}_{171} & = & -x_{24}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{24}\right)\mathbf{a}_2-z_{24}\,\mathbf{a}_3 & = & -x_{24}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{24}\right)b\,\hat{\mathbf{y}}-z_{24}c\,\hat{\mathbf{z}} & (8d)\ \text{O XII}\\
\mathbf{B}_{172} & = & \left(\frac{1}{2}+x_{24}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{24}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{24}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{24}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{24}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{24}\right)c\,\hat{\mathbf{z}} & (8d)\ \text{O XII}\\
\mathbf{B}_{173} & = & -x_{24}\,\mathbf{a}_1-y_{24}\,\mathbf{a}_2-z_{24}\,\mathbf{a}_3 & = & -x_{24}a\,\hat{\mathbf{x}}-y_{24}b\,\hat{\mathbf{y}}-z_{24}c\,\hat{\mathbf{z}} & (8d)\ \text{O XII}\\
\mathbf{B}_{174} & = & \left(\frac{1}{2}+x_{24}\right)\mathbf{a}_1+y_{24}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{24}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{24}\right)a\,\hat{\mathbf{x}}+y_{24}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{24}\right)c\,\hat{\mathbf{z}} & (8d)\ \text{O XII}\\
\mathbf{B}_{175} & = & x_{24}\,\mathbf{a}_1+\left(\frac{1}{2}-y_{24}\right)\mathbf{a}_2+z_{24}\,\mathbf{a}_3 & = & x_{24}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{24}\right)b\,\hat{\mathbf{y}}+z_{24}c\,\hat{\mathbf{z}} & (8d)\ \text{O XII}\\
\mathbf{B}_{176} & = & \left(\frac{1}{2}-x_{24}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{24}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{24}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{24}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{24}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{24}\right)c\,\hat{\mathbf{z}} & (8d)\ \text{O XII}\\
\mathbf{B}_{177} & = & x_{25}\,\mathbf{a}_1+y_{25}\,\mathbf{a}_2+z_{25}\,\mathbf{a}_3 & = & x_{25}a\,\hat{\mathbf{x}}+y_{25}b\,\hat{\mathbf{y}}+z_{25}c\,\hat{\mathbf{z}} & (8d)\ \text{O XIII}
\end{array}
$$

$\mathbf{B}_{178}$	=	$\left(\frac{1}{2}-x_{25}\right)\mathbf{a}_1-y_{25}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{25}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{25}\right)a\,\hat{\mathbf{x}}-y_{25}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{25}\right)c\,\hat{\mathbf{z}}$	(8*d*)	O XIII
$\mathbf{B}_{179}$	=	$-x_{25}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{25}\right)\mathbf{a}_2-z_{25}\,\mathbf{a}_3$	=	$-x_{25}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{25}\right)b\,\hat{\mathbf{y}}-z_{25}c\,\hat{\mathbf{z}}$	(8*d*)	O XIII
$\mathbf{B}_{180}$	=	$\left(\frac{1}{2}+x_{25}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{25}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{25}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{25}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{25}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{25}\right)c\,\hat{\mathbf{z}}$	(8*d*)	O XIII
$\mathbf{B}_{181}$	=	$-x_{25}\,\mathbf{a}_1-y_{25}\,\mathbf{a}_2-z_{25}\,\mathbf{a}_3$	=	$-x_{25}a\,\hat{\mathbf{x}}-y_{25}b\,\hat{\mathbf{y}}-z_{25}c\,\hat{\mathbf{z}}$	(8*d*)	O XIII
$\mathbf{B}_{182}$	=	$\left(\frac{1}{2}+x_{25}\right)\mathbf{a}_1+y_{25}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{25}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{25}\right)a\,\hat{\mathbf{x}}+y_{25}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{25}\right)c\,\hat{\mathbf{z}}$	(8*d*)	O XIII
$\mathbf{B}_{183}$	=	$x_{25}\,\mathbf{a}_1+\left(\frac{1}{2}-y_{25}\right)\mathbf{a}_2+z_{25}\,\mathbf{a}_3$	=	$x_{25}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{25}\right)b\,\hat{\mathbf{y}}+z_{25}c\,\hat{\mathbf{z}}$	(8*d*)	O XIII
$\mathbf{B}_{184}$	=	$\left(\frac{1}{2}-x_{25}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{25}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{25}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{25}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{25}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{25}\right)c\,\hat{\mathbf{z}}$	(8*d*)	O XIII
$\mathbf{B}_{185}$	=	$x_{26}\,\mathbf{a}_1+y_{26}\,\mathbf{a}_2+z_{26}\,\mathbf{a}_3$	=	$x_{26}a\,\hat{\mathbf{x}}+y_{26}b\,\hat{\mathbf{y}}+z_{26}c\,\hat{\mathbf{z}}$	(8*d*)	P
$\mathbf{B}_{186}$	=	$\left(\frac{1}{2}-x_{26}\right)\mathbf{a}_1-y_{26}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{26}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{26}\right)a\,\hat{\mathbf{x}}-y_{26}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{26}\right)c\,\hat{\mathbf{z}}$	(8*d*)	P
$\mathbf{B}_{187}$	=	$-x_{26}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{26}\right)\mathbf{a}_2-z_{26}\,\mathbf{a}_3$	=	$-x_{26}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{26}\right)b\,\hat{\mathbf{y}}-z_{26}c\,\hat{\mathbf{z}}$	(8*d*)	P
$\mathbf{B}_{188}$	=	$\left(\frac{1}{2}+x_{26}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{26}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{26}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{26}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{26}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{26}\right)c\,\hat{\mathbf{z}}$	(8*d*)	P
$\mathbf{B}_{189}$	=	$-x_{26}\,\mathbf{a}_1-y_{26}\,\mathbf{a}_2-z_{26}\,\mathbf{a}_3$	=	$-x_{26}a\,\hat{\mathbf{x}}-y_{26}b\,\hat{\mathbf{y}}-z_{26}c\,\hat{\mathbf{z}}$	(8*d*)	P
$\mathbf{B}_{190}$	=	$\left(\frac{1}{2}+x_{26}\right)\mathbf{a}_1+y_{26}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{26}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{26}\right)a\,\hat{\mathbf{x}}+y_{26}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{26}\right)c\,\hat{\mathbf{z}}$	(8*d*)	P
$\mathbf{B}_{191}$	=	$x_{26}\,\mathbf{a}_1+\left(\frac{1}{2}-y_{26}\right)\mathbf{a}_2+z_{26}\,\mathbf{a}_3$	=	$x_{26}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{26}\right)b\,\hat{\mathbf{y}}+z_{26}c\,\hat{\mathbf{z}}$	(8*d*)	P
$\mathbf{B}_{192}$	=	$\left(\frac{1}{2}-x_{26}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{26}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{26}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{26}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{26}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{26}\right)c\,\hat{\mathbf{z}}$	(8*d*)	P
$\mathbf{B}_{193}$	=	$x_{27}\,\mathbf{a}_1+y_{27}\,\mathbf{a}_2+z_{27}\,\mathbf{a}_3$	=	$x_{27}a\,\hat{\mathbf{x}}+y_{27}b\,\hat{\mathbf{y}}+z_{27}c\,\hat{\mathbf{z}}$	(8*d*)	U
$\mathbf{B}_{194}$	=	$\left(\frac{1}{2}-x_{27}\right)\mathbf{a}_1-y_{27}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{27}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{27}\right)a\,\hat{\mathbf{x}}-y_{27}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{27}\right)c\,\hat{\mathbf{z}}$	(8*d*)	U
$\mathbf{B}_{195}$	=	$-x_{27}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{27}\right)\mathbf{a}_2-z_{27}\,\mathbf{a}_3$	=	$-x_{27}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{27}\right)b\,\hat{\mathbf{y}}-z_{27}c\,\hat{\mathbf{z}}$	(8*d*)	U
$\mathbf{B}_{196}$	=	$\left(\frac{1}{2}+x_{27}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{27}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{27}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{27}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{27}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{27}\right)c\,\hat{\mathbf{z}}$	(8*d*)	U
$\mathbf{B}_{197}$	=	$-x_{27}\,\mathbf{a}_1-y_{27}\,\mathbf{a}_2-z_{27}\,\mathbf{a}_3$	=	$-x_{27}a\,\hat{\mathbf{x}}-y_{27}b\,\hat{\mathbf{y}}-z_{27}c\,\hat{\mathbf{z}}$	(8*d*)	U
$\mathbf{B}_{198}$	=	$\left(\frac{1}{2}+x_{27}\right)\mathbf{a}_1+y_{27}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{27}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{27}\right)a\,\hat{\mathbf{x}}+y_{27}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{27}\right)c\,\hat{\mathbf{z}}$	(8*d*)	U
$\mathbf{B}_{199}$	=	$x_{27}\,\mathbf{a}_1+\left(\frac{1}{2}-y_{27}\right)\mathbf{a}_2+z_{27}\,\mathbf{a}_3$	=	$x_{27}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{27}\right)b\,\hat{\mathbf{y}}+z_{27}c\,\hat{\mathbf{z}}$	(8*d*)	U
$\mathbf{B}_{200}$	=	$\left(\frac{1}{2}-x_{27}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{27}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{27}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{27}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{27}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{27}\right)c\,\hat{\mathbf{z}}$	(8*d*)	U

References:

- A. J. Locock and P. C. Burns, *The crystal structure of synthetic autunite, Ca[(UO$_2$)(PO$_4$)]$_2$(H$_2$O)$_{11}$*, Am. Mineral. **88**, 240–244 (2003), doi:10.2138/am-2003-0128.
- K. Herrmann, ed., *Strukturbericht Band VI 1938* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1941).

Found in:

- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

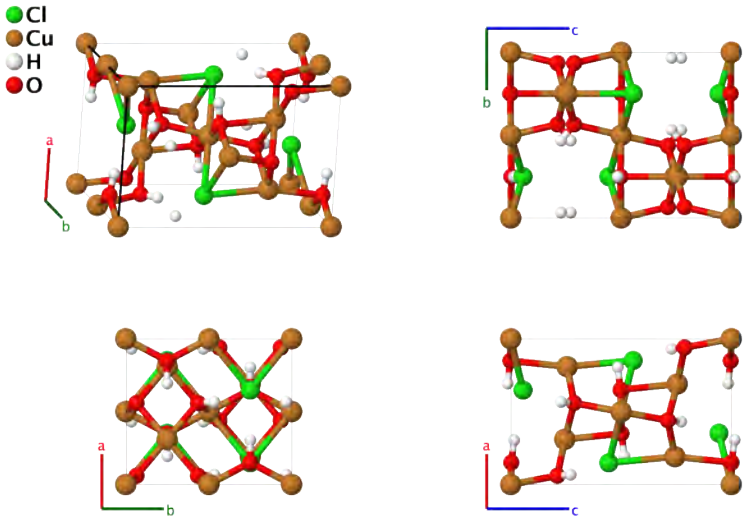
Geometry files:

- CIF: pp. 1642

- POSCAR: pp. 1643

Atacamite ($Cu_2(OH)_3Cl$) Structure: AB2C3D3_oP36_62_c_ac_cd_cd

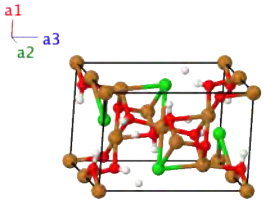
http://aflow.org/prototype-encyclopedia/AB2C3D3_oP36_62_c_ac_cd_cd



Prototype	:	$ClCu_2H_3O_3$
AFLOW prototype label	:	AB2C3D3_oP36_62_c_ac_cd_cd
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP36
Space group number	:	62
Space group symbol	:	*Pnma*
AFLOW prototype command	:	`aflow --proto=AB2C3D3_oP36_62_c_ac_cd_cd --params=`$a, b/a, c/a, x_2, z_2, x_3, z_3, x_4, z_4, x_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7$

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(4*a*)	Cu I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*a*)	Cu I
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}b\,\hat{\mathbf{y}}$	(4*a*)	Cu I
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*a*)	Cu I

$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*c*)	Cl
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	(4*c*)	Cl
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(4*c*)	Cl
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right) c\,\hat{\mathbf{z}}$	(4*c*)	Cl
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*c*)	Cu II
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	(4*c*)	Cu II
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(4*c*)	Cu II
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right) c\,\hat{\mathbf{z}}$	(4*c*)	Cu II
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(4*c*)	H I
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	(4*c*)	H I
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(4*c*)	H I
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\,\hat{\mathbf{z}}$	(4*c*)	H I
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_{19}$	=	$-x_5\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right) c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(8*d*)	H II
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	(8*d*)	H II
$\mathbf{B}_{23}$	=	$-x_6\,\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right) b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(8*d*)	H II
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right) c\,\hat{\mathbf{z}}$	(8*d*)	H II
$\mathbf{B}_{25}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(8*d*)	H II
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + y_6\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right) c\,\hat{\mathbf{z}}$	(8*d*)	H II
$\mathbf{B}_{27}$	=	$x_6\,\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right) b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(8*d*)	H II
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	(8*d*)	H II
$\mathbf{B}_{29}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2} - x_7\right)\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_7\right) a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right) c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{31}$	=	$-x_7\,\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right) b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2} + x_7\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_7\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_7\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_7\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right) c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{33}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{34}$	=	$\left(\frac{1}{2} + x_7\right)\mathbf{a}_1 + y_7\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_7\right) a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right) c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{35}$	=	$x_7\,\mathbf{a}_1 + \left(\frac{1}{2} - y_7\right)\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_7\right) b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{36}$	=	$\left(\frac{1}{2} - x_7\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_7\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right) c\,\hat{\mathbf{z}}$	(8*d*)	O II

References:

- J. B. Parise and B. G. Hyde, *The structure of atacamite and its relationship to spinel*, Acta Crystallogr. C **42**, 1277–1280 (1986), doi:10.1107/S0108270186092570.

Found in:

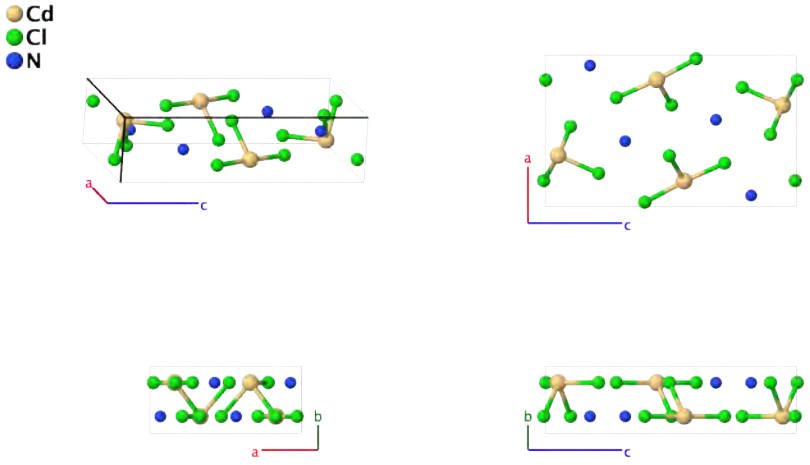
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1644
- POSCAR: pp. 1644

NH_4CdCl_3 ($E2_4$) Structure: AB3C_oP20_62_c_3c_c

http://aflow.org/prototype-encyclopedia/AB3C_oP20_62_c_3c_c



Prototype	:	$CdCl_3(NH_4)$
AFLOW prototype label	:	AB3C_oP20_62_c_3c_c
***Strukturbericht* designation**	:	$E2_4$
Pearson symbol	:	oP20
Space group number	:	62
Space group symbol	:	*Pnma*
AFLOW prototype command	:	`aflow --proto=AB3C_oP20_62_c_3c_c --params=`$a, b/a, c/a, x_1, z_1, x_2, z_2, x_3, z_3, x_4, z_4, x_5, z_5$

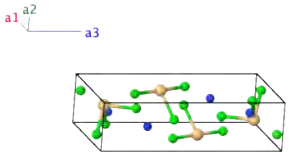
Other compounds with this structure

- $LaCrS_3$, $CeCrS_3$, $SmCrS_3$, $NdCrS_3$, $CeCrSe_3$, $PbSnS_3$, $RbCdBr_3$, $RbCdCl_3$, and $(Ti,Sn)_2S_3$

- (Brasseur, 1938) gave the Wyckoff positions in the *Pnam* setting of space group #62. We used FINDSYM to rotate the structure and put it in the standard *Pnma* setting.
- The hydrogen atoms for the NH_4 ions were not determined, so we only give the positions of the nitrogen atoms (labeled as NH_4).

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(4*c*)	Cd
$\mathbf{B}_2$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	Cd
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	(4*c*)	Cd
$\mathbf{B}_4$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	Cd
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*c*)	Cl I
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	(4*c*)	Cl I
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(4*c*)	Cl I
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right) c\,\hat{\mathbf{z}}$	(4*c*)	Cl I
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*c*)	Cl II
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	(4*c*)	Cl II
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(4*c*)	Cl II
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right) c\,\hat{\mathbf{z}}$	(4*c*)	Cl II
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(4*c*)	Cl III
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	(4*c*)	Cl III
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(4*c*)	Cl III
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\,\hat{\mathbf{z}}$	(4*c*)	Cl III
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(4*c*)	NH_4
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	(4*c*)	NH_4
$\mathbf{B}_{19}$	=	$-x_5\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(4*c*)	NH_4
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right) c\,\hat{\mathbf{z}}$	(4*c*)	NH_4

References:

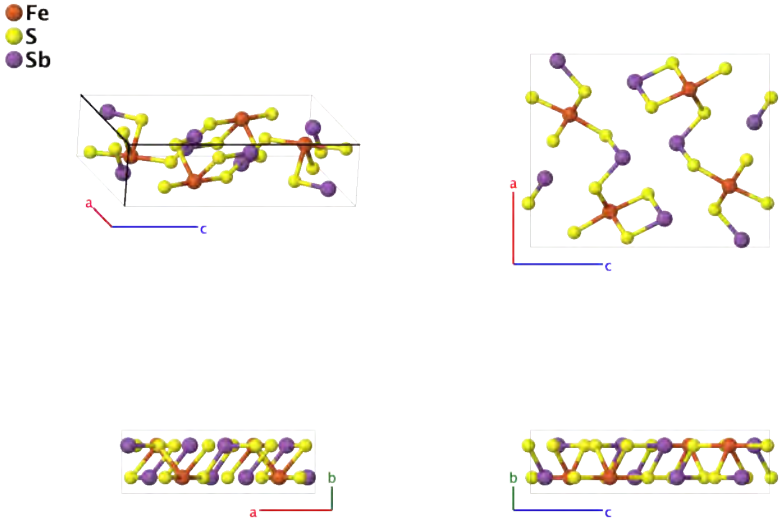
- H. Brasseur and L. Pauling, *The Crystal Structure of Ammonium Cadmium Chloride, NH_4CdCl_3*, J. Am. Chem. Soc. **60**, 2886–2890 (1938), doi:10.1021/ja01279a016.

Geometry files:

- CIF: pp. 1645
- POSCAR: pp. 1645

Berthierite ($FeSb_2S_4$, $E3_3$) Structure: AB4C2_oP28_62_c_4c_2c

http://aflow.org/prototype-encyclopedia/AB4C2_oP28_62_c_4c_2c

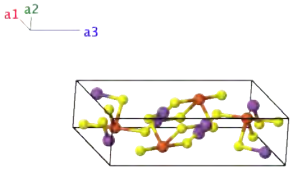


Prototype	:	FeS_4Sb_2
AFLOW prototype label	:	AB4C2_oP28_62_c_4c_2c
***Strukturbericht* designation**	:	$E3_3$
Pearson symbol	:	oP28
Space group number	:	62
Space group symbol	:	*Pnma*
AFLOW prototype command	:	`aflow --proto=AB4C2_oP28_62_c_4c_2c --params=`$a, b/a, c/a, x_1, z_1, x_2, z_2, x_3, z_3, x_4, z_4, x_5, z_5, x_6, z_6, x_7, z_7$

- The data for this structure was given in the *Pnam* orientation of space group #62. We used FINDSYM to place it in the standard *Pnma* orientation.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$\mathbf{B}_{1}$	=	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(4*c*)	Fe
$\mathbf{B}_{2}$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	Fe
$\mathbf{B}_{3}$	=	$-x_1\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	(4*c*)	Fe
$\mathbf{B}_{4}$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	Fe
$\mathbf{B}_{5}$	=	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*c*)	S I
$\mathbf{B}_{6}$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	(4*c*)	S I
$\mathbf{B}_{7}$	=	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(4*c*)	S I
$\mathbf{B}_{8}$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right) c\,\hat{\mathbf{z}}$	(4*c*)	S I
$\mathbf{B}_{9}$	=	$x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*c*)	S II
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	(4*c*)	S II
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(4*c*)	S II
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right) c\,\hat{\mathbf{z}}$	(4*c*)	S II
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(4*c*)	S III
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	(4*c*)	S III
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(4*c*)	S III
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\,\hat{\mathbf{z}}$	(4*c*)	S III
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(4*c*)	S IV
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	(4*c*)	S IV
$\mathbf{B}_{19}$	=	$-x_5\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(4*c*)	S IV
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right) c\,\hat{\mathbf{z}}$	(4*c*)	S IV
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(4*c*)	Sb I
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	(4*c*)	Sb I
$\mathbf{B}_{23}$	=	$-x_6\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(4*c*)	Sb I
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right) c\,\hat{\mathbf{z}}$	(4*c*)	Sb I
$\mathbf{B}_{25}$	=	$x_7\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(4*c*)	Sb II
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} - x_7\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_7\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right) c\,\hat{\mathbf{z}}$	(4*c*)	Sb II
$\mathbf{B}_{27}$	=	$-x_7\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	(4*c*)	Sb II
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} + x_7\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_7\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right) c\,\hat{\mathbf{z}}$	(4*c*)	Sb II

References:

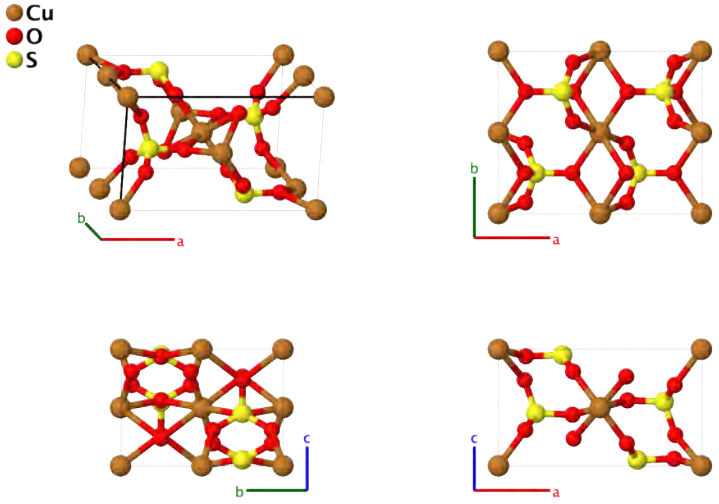
- M. J. Buerger and T. Hahn, *The Crystal Structure of Berthierite, $FeSb_2S_4$*, http://www.dtic.mil/dtic/tr/fulltext/u2/022245.pdf (1953). ONR Technical Report 2, Project NR 032 346.

Geometry files:

- CIF: pp. 1645
- POSCAR: pp. 1645

Chalcocyanite ($CuSO_4$) Structure: AB4C_oP24_62_a_2cd_c

http://aflow.org/prototype-encyclopedia/AB4C_oP24_62_a_2cd_c



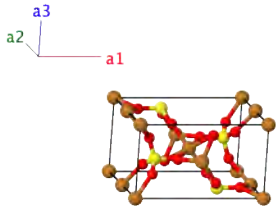
Prototype	:	CuO_4S
AFLOW prototype label	:	AB4C_oP24_62_a_2cd_c
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP24
Space group number	:	62
Space group symbol	:	*Pnma*
AFLOW prototype command	:	`aflow --proto=AB4C_oP24_62_a_2cd_c --params=`$a, b/a, c/a, x_2, z_2, x_3, z_3, x_4, z_4, x_5, y_5, z_5$

Other compounds with this structure

- $ZnSO_4$ (zincosite)

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(4a)$	Cu
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(4a)$	Cu

$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}b\,\hat{\mathbf{y}}$	($4a$)	Cu
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4a$)	Cu
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_2c\,\hat{\mathbf{z}}$	($4c$)	O I
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	($4c$)	O I
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_2c\,\hat{\mathbf{z}}$	($4c$)	O I
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	($4c$)	O I
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	($4c$)	O II
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	($4c$)	O II
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	($4c$)	O II
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	($4c$)	O II
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	($4c$)	S
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	($4c$)	S
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	($4c$)	S
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	($4c$)	S
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}} + y_5b\,\hat{\mathbf{y}} + z_5c\,\hat{\mathbf{z}}$	($8d$)	O III
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} - y_5b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	($8d$)	O III
$\mathbf{B}_{19}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right)b\,\hat{\mathbf{y}} - z_5c\,\hat{\mathbf{z}}$	($8d$)	O III
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	($8d$)	O III
$\mathbf{B}_{21}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}} - y_5b\,\hat{\mathbf{y}} - z_5c\,\hat{\mathbf{z}}$	($8d$)	O III
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + y_5\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + y_5b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	($8d$)	O III
$\mathbf{B}_{23}$	=	$x_5\,\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right)b\,\hat{\mathbf{y}} + z_5c\,\hat{\mathbf{z}}$	($8d$)	O III
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	($8d$)	O III

References:

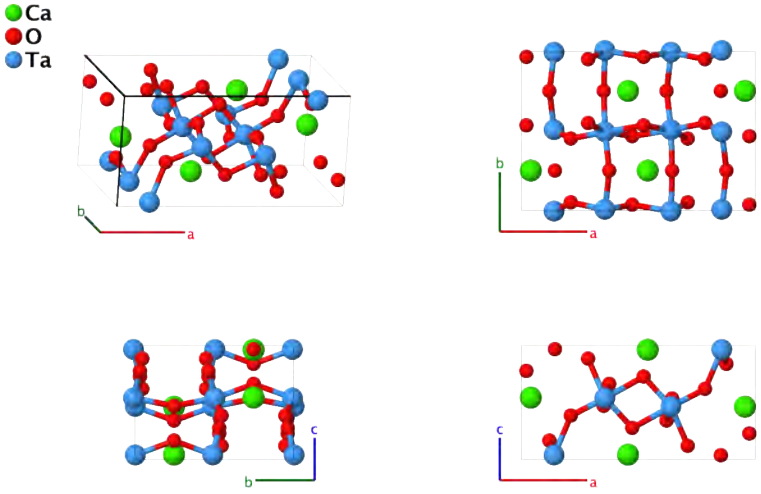
- M. Wildner and G. Giester, *Crystal structure refinements of synthetic chalcocyanite ($CuSO_4$) and zincosite ($ZnSO_4$)*, Mineral. Petrol. **39**, 201–209 (1988), doi:10.1007/BF01163035.

Geometry files:

- CIF: pp. 1646
- POSCAR: pp. 1646

Rynersonite (Orthorhombic $CaTa_2O_6$) Structure: AB6C2_oP36_62_c_2c2d_d

http://aflow.org/prototype-encyclopedia/AB6C2_oP36_62_c_2c2d_d



Prototype	:	CaO_6Ta_2
AFLOW prototype label	:	AB6C2_oP36_62_c_2c2d_d
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP36
Space group number	:	62
Space group symbol	:	*Pnma*
AFLOW prototype command	:	`aflow --proto=AB6C2_oP36_62_c_2c2d_d --params=`$a, b/a, c/a, x_1, z_1, x_2, z_2, x_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6$

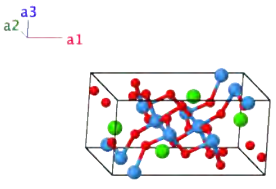
Other compounds with this structure

- $Ca(Ta,Ti)O_6$, $CaTa_2O_6$, $EuTa_2O_6$, $GdTa_2O_6$, $GdTi_2O_6$, $La(Nb,Ti)O_6$, $LaTa_2O_6$, $LaTi_2O_6$, $Nd(Ta,Ti)O_6$, $Pr(Ta,Ti)O_6$, $SrNb_2O_6$, $Tb(Ta,Ti)O_6$, YTa_2O_6, and $(Sr_{0.7}La_{0.3})Nb_2O_6$

- This is the room-temperature structure of $CaTa_2O_6$. At 500 °C, this takes on the cubic perovskite structure with a half-filled calcium site.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(4*c*)	Ca
$\mathbf{B}_2$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	Ca
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	(4*c*)	Ca
$\mathbf{B}_4$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	Ca
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right) c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right) c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right) b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{17}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{19}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right) b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{21}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right) a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{23}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right) b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right) c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{25}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + y_5\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right) c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{27}$	=	$x_5\,\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right) b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{29}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(8*d*)	Ta
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	(8*d*)	Ta
$\mathbf{B}_{31}$	=	$-x_6\,\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right) b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(8*d*)	Ta

$\mathbf{B}_{32}$	=	$\left(\frac{1}{2}+x_6\right)\mathbf{a}_1+\left(\frac{1}{2}-y_6\right)\mathbf{a}_2+\left(\frac{1}{2}-z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_6\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_6\right)c\,\hat{\mathbf{z}}$	(8*d*)	Ta
$\mathbf{B}_{33}$	=	$-x_6\,\mathbf{a}_1-y_6\,\mathbf{a}_2-z_6\,\mathbf{a}_3$	=	$-x_6a\,\hat{\mathbf{x}}-y_6b\,\hat{\mathbf{y}}-z_6c\,\hat{\mathbf{z}}$	(8*d*)	Ta
$\mathbf{B}_{34}$	=	$\left(\frac{1}{2}+x_6\right)\mathbf{a}_1+y_6\,\mathbf{a}_2+\left(\frac{1}{2}-z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{x}}+y_6b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_6\right)c\,\hat{\mathbf{z}}$	(8*d*)	Ta
$\mathbf{B}_{35}$	=	$x_6\,\mathbf{a}_1+\left(\frac{1}{2}-y_6\right)\mathbf{a}_2+z_6\,\mathbf{a}_3$	=	$x_6a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_6\right)b\,\hat{\mathbf{y}}+z_6c\,\hat{\mathbf{z}}$	(8*d*)	Ta
$\mathbf{B}_{36}$	=	$\left(\frac{1}{2}-x_6\right)\mathbf{a}_1+\left(\frac{1}{2}+y_6\right)\mathbf{a}_2+\left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_6\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	(8*d*)	Ta

References:

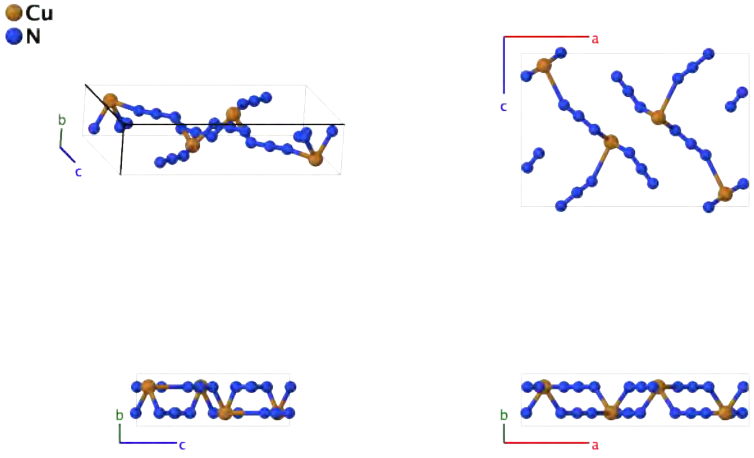
- L. Jahnberg, *Crystal Structure of Orthorhombic $CaTa_2O_6$*, Acta Chem. Scand. **71**, 2548–2559 (1963), doi:10.3891/acta.chem.scand.17-2548.

Geometry files:

- CIF: pp. 1646
- POSCAR: pp. 1646

Copper (II) Azide [$Cu(N_3)_2$] Structure: AB6_oP28_62_c_6c

http://aflow.org/prototype-encyclopedia/AB6_oP28_62_c_6c

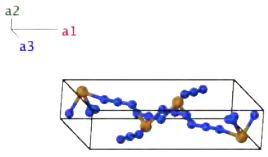


Prototype	:	CuN_6
AFLOW prototype label	:	AB6_oP28_62_c_6c
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP28
Space group number	:	62
Space group symbol	:	*Pnma*
AFLOW prototype command	:	`aflow --proto=AB6_oP28_62_c_6c --params=`$a, b/a, c/a, x_1, z_1, x_2, z_2, x_3, z_3, x_4, z_4, x_5, z_5, x_6, z_6, x_7, z_7$

- Not to be confused with Copper (I) Azide, CuN_3.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(4*c*)	Cu
$\mathbf{B}_2$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	Cu
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	(4*c*)	Cu

$\mathbf{B}_4$	$=$	$\left(\frac{1}{2}+x_1\right)\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+\left(\frac{1}{2}-z_1\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}+x_1\right)a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_1\right)c\,\hat{\mathbf{z}}$	(4*c*)	Cu
$\mathbf{B}_5$	$=$	$x_2\,\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+z_2\,\mathbf{a}_3$	$=$	$x_2a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+z_2c\,\hat{\mathbf{z}}$	(4*c*)	N I
$\mathbf{B}_6$	$=$	$\left(\frac{1}{2}-x_2\right)\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2+\left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}-x_2\right)a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_2\right)c\,\hat{\mathbf{z}}$	(4*c*)	N I
$\mathbf{B}_7$	$=$	$-x_2\,\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2-z_2\,\mathbf{a}_3$	$=$	$-x_2a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}-z_2c\,\hat{\mathbf{z}}$	(4*c*)	N I
$\mathbf{B}_8$	$=$	$\left(\frac{1}{2}+x_2\right)\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+\left(\frac{1}{2}-z_2\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}+x_2\right)a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_2\right)c\,\hat{\mathbf{z}}$	(4*c*)	N I
$\mathbf{B}_9$	$=$	$x_3\,\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+z_3\,\mathbf{a}_3$	$=$	$x_3a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+z_3c\,\hat{\mathbf{z}}$	(4*c*)	N II
$\mathbf{B}_{10}$	$=$	$\left(\frac{1}{2}-x_3\right)\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2+\left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_3\right)c\,\hat{\mathbf{z}}$	(4*c*)	N II
$\mathbf{B}_{11}$	$=$	$-x_3\,\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2-z_3\,\mathbf{a}_3$	$=$	$-x_3a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}-z_3c\,\hat{\mathbf{z}}$	(4*c*)	N II
$\mathbf{B}_{12}$	$=$	$\left(\frac{1}{2}+x_3\right)\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+\left(\frac{1}{2}-z_3\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_3\right)c\,\hat{\mathbf{z}}$	(4*c*)	N II
$\mathbf{B}_{13}$	$=$	$x_4\,\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+z_4\,\mathbf{a}_3$	$=$	$x_4a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+z_4c\,\hat{\mathbf{z}}$	(4*c*)	N III
$\mathbf{B}_{14}$	$=$	$\left(\frac{1}{2}-x_4\right)\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2+\left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_4\right)c\,\hat{\mathbf{z}}$	(4*c*)	N III
$\mathbf{B}_{15}$	$=$	$-x_4\,\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2-z_4\,\mathbf{a}_3$	$=$	$-x_4a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}-z_4c\,\hat{\mathbf{z}}$	(4*c*)	N III
$\mathbf{B}_{16}$	$=$	$\left(\frac{1}{2}+x_4\right)\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+\left(\frac{1}{2}-z_4\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_4\right)c\,\hat{\mathbf{z}}$	(4*c*)	N III
$\mathbf{B}_{17}$	$=$	$x_5\,\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+z_5\,\mathbf{a}_3$	$=$	$x_5a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+z_5c\,\hat{\mathbf{z}}$	(4*c*)	N IV
$\mathbf{B}_{18}$	$=$	$\left(\frac{1}{2}-x_5\right)\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	(4*c*)	N IV
$\mathbf{B}_{19}$	$=$	$-x_5\,\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2-z_5\,\mathbf{a}_3$	$=$	$-x_5a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}-z_5c\,\hat{\mathbf{z}}$	(4*c*)	N IV
$\mathbf{B}_{20}$	$=$	$\left(\frac{1}{2}+x_5\right)\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+\left(\frac{1}{2}-z_5\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_5\right)c\,\hat{\mathbf{z}}$	(4*c*)	N IV
$\mathbf{B}_{21}$	$=$	$x_6\,\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+z_6\,\mathbf{a}_3$	$=$	$x_6a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+z_6c\,\hat{\mathbf{z}}$	(4*c*)	N V
$\mathbf{B}_{22}$	$=$	$\left(\frac{1}{2}-x_6\right)\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2+\left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	(4*c*)	N V
$\mathbf{B}_{23}$	$=$	$-x_6\,\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2-z_6\,\mathbf{a}_3$	$=$	$-x_6a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}-z_6c\,\hat{\mathbf{z}}$	(4*c*)	N V
$\mathbf{B}_{24}$	$=$	$\left(\frac{1}{2}+x_6\right)\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+\left(\frac{1}{2}-z_6\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_6\right)c\,\hat{\mathbf{z}}$	(4*c*)	N V
$\mathbf{B}_{25}$	$=$	$x_7\,\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+z_7\,\mathbf{a}_3$	$=$	$x_7a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+z_7c\,\hat{\mathbf{z}}$	(4*c*)	N VI
$\mathbf{B}_{26}$	$=$	$\left(\frac{1}{2}-x_7\right)\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2+\left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}-x_7\right)a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	(4*c*)	N VI
$\mathbf{B}_{27}$	$=$	$-x_7\,\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2-z_7\,\mathbf{a}_3$	$=$	$-x_7a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}-z_7c\,\hat{\mathbf{z}}$	(4*c*)	N VI
$\mathbf{B}_{28}$	$=$	$\left(\frac{1}{2}+x_7\right)\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+\left(\frac{1}{2}-z_7\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_7\right)c\,\hat{\mathbf{z}}$	(4*c*)	N VI

References:

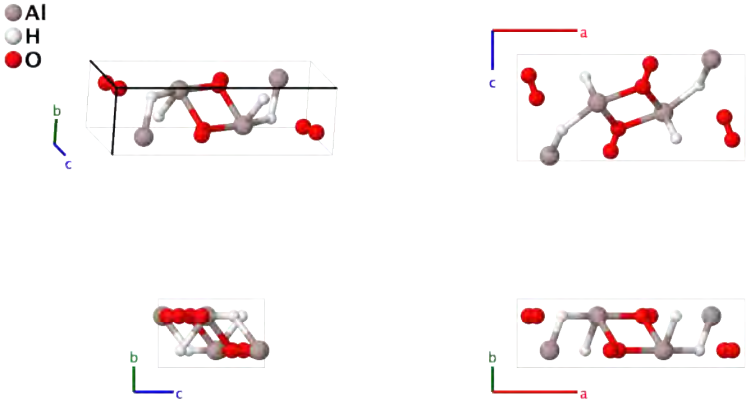
- I. Agrell, *The Crystal Structure of Cu(N_3)$_2$*, Acta Chem. Scand. **21**, 2647–2658 (1967), doi:10.3891/acta.chem.scand.21-2647.

Geometry files:

- CIF: pp. 1647
- POSCAR: pp. 1647

Diaspore ($AlOOH$, $E0_2$) Structure: ABC2_oP16_62_c_c_2c

http://aflow.org/prototype-encyclopedia/ABC2_oP16_62_c_c_2c



Prototype	:	$AlHO_2$
AFLOW prototype label	:	ABC2_oP16_62_c_c_2c
***Strukturbericht* designation**	:	$E0_2$
Pearson symbol	:	oP16
Space group number	:	62
Space group symbol	:	*Pnma*
AFLOW prototype command	:	`aflow --proto=ABC2_oP16_62_c_c_2c --params=`$a, b/a, c/a, x_1, z_1, x_2, z_2, x_3, z_3, x_4, z_4$

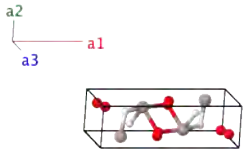
Other compounds with this structure

- FeO(OH) (goethite) and MnO(OH) (groutite)

- (Hermann, 1937), following (Verway, 1935) designated this as *Strukturbericht* $E0_2$, but had no knowledge of the positions of the hydrogen atoms, which we take from (Hill, 1979) and which do not change the space group. Hill gave the Wyckoff positions of the atoms in terms of the *Pbnm* setting of space group #62. We used FINDSYM to transform this structure to the standard *Pnma* setting.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$ =	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	($4c$)	Al

$\mathbf{B}_2$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \frac{3}{4}\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right)a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\,\hat{\mathbf{z}}$	(4*c*)	Al
$\mathbf{B}_3$	=	$-x_1\mathbf{a}_1 + \frac{3}{4}\mathbf{a}_2 - z_1\mathbf{a}_3$	=	$-x_1a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_1c\,\hat{\mathbf{z}}$	(4*c*)	Al
$\mathbf{B}_4$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \frac{1}{4}\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right)a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right)c\,\hat{\mathbf{z}}$	(4*c*)	Al
$\mathbf{B}_5$	=	$x_2\mathbf{a}_1 + \frac{1}{4}\mathbf{a}_2 + z_2\mathbf{a}_3$	=	$x_2a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_2c\,\hat{\mathbf{z}}$	(4*c*)	H
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \frac{3}{4}\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	(4*c*)	H
$\mathbf{B}_7$	=	$-x_2\mathbf{a}_1 + \frac{3}{4}\mathbf{a}_2 - z_2\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_2c\,\hat{\mathbf{z}}$	(4*c*)	H
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \frac{1}{4}\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	(4*c*)	H
$\mathbf{B}_9$	=	$x_3\mathbf{a}_1 + \frac{1}{4}\mathbf{a}_2 + z_3\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \frac{3}{4}\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_{11}$	=	$-x_3\mathbf{a}_1 + \frac{3}{4}\mathbf{a}_2 - z_3\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \frac{1}{4}\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_{13}$	=	$x_4\mathbf{a}_1 + \frac{1}{4}\mathbf{a}_2 + z_4\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \frac{3}{4}\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_{15}$	=	$-x_4\mathbf{a}_1 + \frac{3}{4}\mathbf{a}_2 - z_4\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + \frac{3}{4}b\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \frac{1}{4}\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(4*c*)	O II

References:

- R. J. Hill, *Crystal Structure Refinement and Electron Density Distribution in Diaspore*, Phys. Chem. Miner. **5**, 179–200 (1979), doi:10.1007/BF00307552.
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).
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Found in:

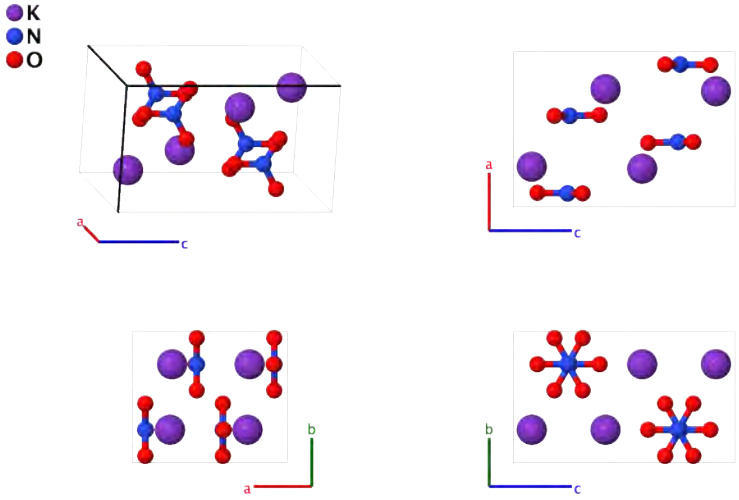
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1647
- POSCAR: pp. 1648

α-Potassium Nitrate (KNO_3) I Structure: ABC3_oP20_62_c_c_cd

http://aflow.org/prototype-encyclopedia/ABC3_oP20_62_c_c_cd.K.N.O3

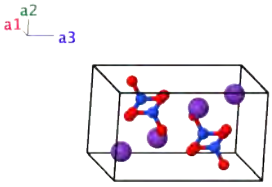


Prototype	:	KNO_3
AFLOW prototype label	:	ABC3_oP20_62_c_c_cd
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP20
Space group number	:	62
Space group symbol	:	*Pnma*
AFLOW prototype command	:	`aflow --proto=ABC3_oP20_62_c_c_cd --params=`$a, b/a, c/a, x_1, z_1, x_2, z_2, x_3, z_3, x_4, y_4, z_4$

- Two possible structures have been identified for α-KNO_3: (Nimmo, 1973) proposed the current structure, which we call "Structure I", in space group *Pnma* #62. (Adiwidjaja, 2003) found this structure, but also noted that it could be described by a doubling of the current unit cell into a superstructure with space group $Cmc2_1$ #36, which we call "Structure II". It is unclear to us which structure is correct, so we present both of them.
- (Nimmo, 1973) determined the lattice constants and Wyckoff positions of the type I structure at 25 °C in the *Pmcn* setting of space group #62. We present the data in the standard *Pnma* setting.
- This structure has the same space group and Wyckoff positions as the MgB_4 (A4B_oP20_62_2cd_c) structure and the $G0_{10}$ $(NH_4)NO_3$ (ABC3_oP20_62_c_c_cd.N.NH4.O) structure, but the structures are all different. We give the previous two structures the noted AFLOW designations, and give this one the designation ABC3_oP20_62_c_c_cd.K.N.O3.
- On heating, α-KNO_3 transforms into β-KNO_3 at 128 °C. When heated above 200 °C and then cooled, the β-phase transforms into the metastable ferroelectric γ-KNO_3 phase, which can remain down to room temperature.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	($4c$)	K
$\mathbf{B}_2$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	($4c$)	K
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	($4c$)	K
$\mathbf{B}_4$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right) c\,\hat{\mathbf{z}}$	($4c$)	K
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	($4c$)	N
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	($4c$)	N
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	($4c$)	N
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right) c\,\hat{\mathbf{z}}$	($4c$)	N
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	($4c$)	O I
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	($4c$)	O I
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	($4c$)	O I
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right) c\,\hat{\mathbf{z}}$	($4c$)	O I
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	($8d$)	O II
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	($8d$)	O II
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right) b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	($8d$)	O II
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\,\hat{\mathbf{z}}$	($8d$)	O II
$\mathbf{B}_{17}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	($8d$)	O II
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\,\hat{\mathbf{z}}$	($8d$)	O II
$\mathbf{B}_{19}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right) b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	($8d$)	O II
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	($8d$)	O II

References:

- J. K. Nimmo and B. W. Lucas, *A neutron diffraction determination of the crystal structure of α-phase potassium nitrate at 25°C and 100°C*, J. Phys. C: Solid State Phys. **6**, 201–211 (1973), doi:10.1088/0022-3719/6/2/001.
- G. Adiwidjaja and D. Pohl, *Superstructure of α-phase potassium nitrate*, Acta Crystallogr. C **59**, i139–i140 (2003), doi:10.1107/S0108270103025277.

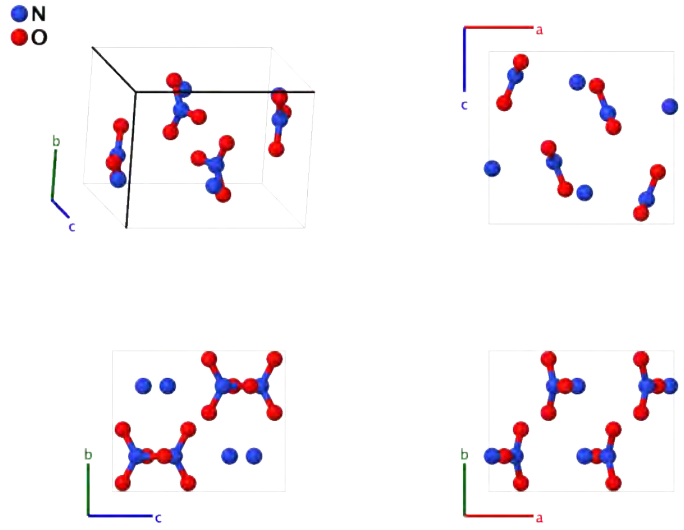
Found in:

- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:
- CIF: pp. 1648
- POSCAR: pp. 1648

NH_4NO_3 III ($G0_{10}$) Structure: ABC3_oP20_62_c_c_cd

http://aflow.org/prototype-encyclopedia/ABC3_oP20_62_c_c_cd.N.NH4.O



Prototype	:	$N(NH_4)O_3$
AFLOW prototype label	:	ABC3_oP20_62_c_c_cd
***Strukturbericht* designation**	:	$G0_{10}$
Pearson symbol	:	oP20
Space group number	:	62
Space group symbol	:	*Pnma*
AFLOW prototype command	:	`aflow --proto=ABC3_oP20_62_c_c_cd --params=`$a, b/a, c/a, x_1, z_1, x_2, z_2, x_3, z_3, x_4, y_4, z_4$

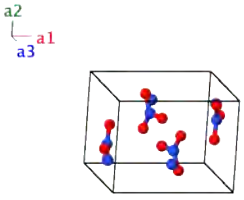
- Ammonium Nitrate exists in a variety of forms, (Hermann, 1937) depending on the temperature:

Phase	Temperature °C	*Strukturbericht*	Page	
I	125 – 170	$G0_8$	AB_cP2_221_a_b.NH4.NO3	
II	84 – 125	$G0_9$	ABC3_tP10_100_b_a_bc	
III	32 – 84	$G0_{10}$	ABC3_oP20_62_c_c_cd.N.NH4.O	(this structure)
IV	-17 – 32	$G0_{11}$	A4B2C3_oP18_59_ef_ab_af	
V	< −17	Gwihabaite	A4B2C3_tP72_77_8d_ab2c2d_6d2	

- Data for this structure was taken at 46 °C.
- It is likely that the NH_4 ions are free to rotate (Kracek, 1937).
- The positions of the hydrogen atoms in the NH_4 ions were not determined, so we only provide the positions of the nitrogen atoms (labeled as NH_4).
- (Goodwin, 1947) gave the structure in the *Pbnm* setting of space group #62. We used FINDSYM to change this to the standard *Pnma* setting.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(4*c*)	N
$\mathbf{B}_2$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	N
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	(4*c*)	N
$\mathbf{B}_4$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	N
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*c*)	NH_4
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	(4*c*)	NH_4
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(4*c*)	NH_4
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right) c\,\hat{\mathbf{z}}$	(4*c*)	NH_4
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right) c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right) b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{17}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{19}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right) b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	(8*d*)	O II

References:

- T. H. Goodwin and J. Whetstone, *The crystal structure of ammonium nitrate III, and atomic scattering factors in ionic crystals*, J. Chem. Soc. pp. 1455–1461 (1947), doi:10.1039/JR9470001455.
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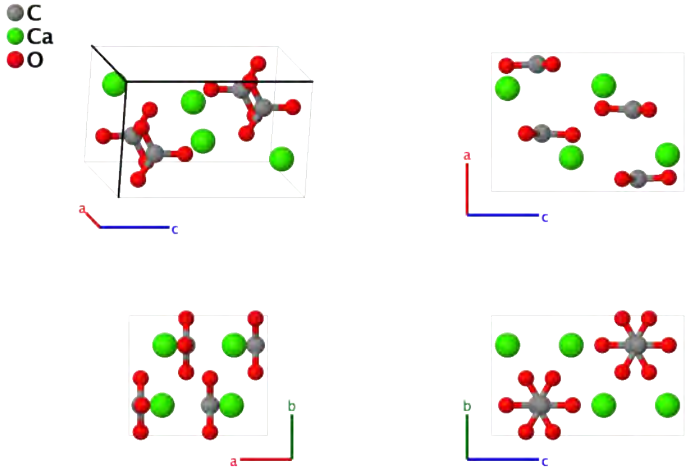
- F. C. Kracek, S. B. Hendricks, and E. Posnjak, *Group Rotation in Solid Ammonium and Calcium Nitrates*, Nature **128**, 410–411 (1931), doi:10.1038/128410b0.

Geometry files:
- CIF: pp. 1648
- POSCAR: pp. 1649

Aragonite ($CaCO_3$, $G0_2$) Structure: ABC3_oP20_62_c_c_cd

http://aflow.org/prototype-encyclopedia/ABC3_oP20_62_c_c_cd



Prototype	:	$CaCO_3$
AFLOW prototype label	:	ABC3_oP20_62_c_c_cd
***Strukturbericht* designation**	:	$G0_2$
Pearson symbol	:	oP20
Space group number	:	62
Space group symbol	:	*Pnma*
AFLOW prototype command	:	`aflow --proto=ABC3_oP20_62_c_c_cd --params=`$a, b/a, c/a, x_1, z_1, x_2, z_2, x_3, z_3, x_4, y_4, z_4$

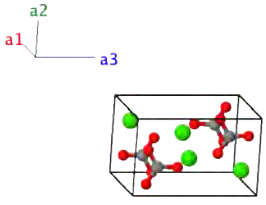
Other compounds with this structure

- $BaCO_3$ (witherite), $(Ba,Ca)CO_3$ (bromlite), $(Cu,Mg,Mn)CO_3$ (manganocalcite), $MnCO_3$ (rhodochrosite), $PbCO_3$ (cerussite), and $SrCO_3$ (strontianite)

- (Ewald, 1931) originally listed this as *Strukturbericht* $G2$, but it was designated $G0_2$ in (Gottfried, 1937).

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(4*c*)	C
$\mathbf{B}_2$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	C
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	(4*c*)	C
$\mathbf{B}_4$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	C
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*c*)	Ca
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	(4*c*)	Ca
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(4*c*)	Ca
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right) c\,\hat{\mathbf{z}}$	(4*c*)	Ca
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right) c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right) b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{17}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{19}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right) b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*d*)	O II
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right) b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	(8*d*)	O II

References:

- J. P. R. de Villiers, *Crystal Structures of Aragonite, Strontianite, and Witherite*, Am. Mineral. **56**, 758–767 (1971).
- P. P. Ewald and C. Hermann, eds., *Strukturbericht 1913-1928* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1931).
- C. Gottfried and F. Schossberger, eds., *Strukturbericht Band III 1933-1935* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Found in:

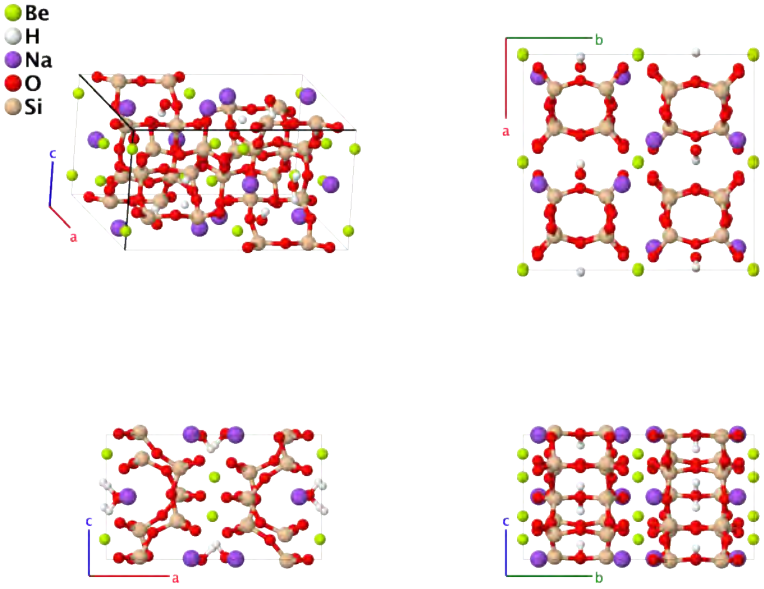
- mindat.org, *Aragonite*, http://www.mindat.org/min-307.html. Mineral Database.

Geometry files:

- CIF: pp. 1649
- POSCAR: pp. 1649

Epididymite ($BeHNaO_8Si_3$, $S4_7$) Structure: ABCD8E3_oP112_62_d_2c_d_4c6d_3d

http://aflow.org/prototype-encyclopedia/ABCD8E3_oP112_62_d_2c_d_4c6d_3d

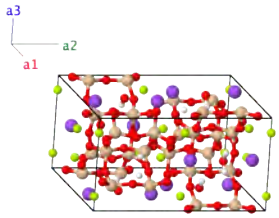


Prototype	:	$BeHNaO_8Si_3$
AFLOW prototype label	:	ABCD8E3_oP112_62_d_2c_d_4c6d_3d
***Strukturbericht* designation**	:	$S4_7$
Pearson symbol	:	oP112
Space group number	:	62
Space group symbol	:	*Pnma*
AFLOW prototype command	:	`aflow --proto=ABCD8E3_oP112_62_d_2c_d_4c6d_3d --params=`$a, b/a, c/a, x_1, z_1, x_2, z_2, x_3, z_3, x_4, z_4, x_5, z_5, x_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}, x_{17}, y_{17}, z_{17}$

- (Ito, 1934)'s original determination of the epididymite structure, designated as $S4_7$ by (Gottfried, 1937), was flawed, even in the author's estimation. In particular, both the Si-O and Si-Be distances are very small, with Si-Be less than 1Å. In addition, Ito and many later authors assumed that the hydrogen atoms formed OH radicals. (Diego Gatta, 2008) found that the hydrogen atoms were actually a part of a water molecule. Since the initial structure is not even suitable for starting first-principles calculations, while the new structure maintains the *Pnma* #62 space group, we have dropped the former and designate the later as $S4_7$.
- Epididymite and its dimorph, eudidymite, are two forms of hydrated sodium beryllium silicate which are stable under ambient conditions (Diego Gatta, 2008).

Simple Orthorhombic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= a\,\hat{\mathbf{x}}\\ \mathbf{a}_2 &= b\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	$x_1 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	$(4c)$	H I
$\mathbf{B}_2$	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	$(4c)$	H I
$\mathbf{B}_3$	$-x_1\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	$-x_1 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	$(4c)$	H I
$\mathbf{B}_4$	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_1\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right) c\,\hat{\mathbf{z}}$	$(4c)$	H I
$\mathbf{B}_5$	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	$x_2 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(4c)$	H II
$\mathbf{B}_6$	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	$(4c)$	H II
$\mathbf{B}_7$	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	$-x_2 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	$(4c)$	H II
$\mathbf{B}_8$	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right) c\,\hat{\mathbf{z}}$	$(4c)$	H II
$\mathbf{B}_9$	$x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$x_3 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(4c)$	O I
$\mathbf{B}_{10}$	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_3\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	$(4c)$	O I
$\mathbf{B}_{11}$	$-x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	$-x_3 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(4c)$	O I
$\mathbf{B}_{12}$	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_3\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right) c\,\hat{\mathbf{z}}$	$(4c)$	O I
$\mathbf{B}_{13}$	$x_4\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$x_4 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(4c)$	O II
$\mathbf{B}_{14}$	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	$(4c)$	O II
$\mathbf{B}_{15}$	$-x_4\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	$-x_4 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	$(4c)$	O II
$\mathbf{B}_{16}$	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\,\hat{\mathbf{z}}$	$(4c)$	O II
$\mathbf{B}_{17}$	$x_5\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	$x_5 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(4c)$	O III
$\mathbf{B}_{18}$	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_5\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right) c\,\hat{\mathbf{z}}$	$(4c)$	O III
$\mathbf{B}_{19}$	$-x_5\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	$-x_5 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	$(4c)$	O III
$\mathbf{B}_{20}$	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right) c\,\hat{\mathbf{z}}$	$(4c)$	O III
$\mathbf{B}_{21}$	$x_6\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	$x_6 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(4c)$	O IV
$\mathbf{B}_{22}$	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	$(4c)$	O IV
$\mathbf{B}_{23}$	$-x_6\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	$-x_6 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	$(4c)$	O IV
$\mathbf{B}_{24}$	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right) c\,\hat{\mathbf{z}}$	$(4c)$	O IV
$\mathbf{B}_{25}$	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	$x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	$(8d)$	Be
$\mathbf{B}_{26}$	$\left(\frac{1}{2} - x_7\right)\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_7\right) a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right) c\,\hat{\mathbf{z}}$	$(8d)$	Be
$\mathbf{B}_{27}$	$-x_7\,\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\mathbf{a}_2 - z_7\,\mathbf{a}_3$	$-x_7 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right) b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	$(8d)$	Be

$$
\begin{array}{lcccll}
\mathbf{B}_{28} & = & \left(\frac{1}{2}+x_7\right)\mathbf{a}_1+\left(\frac{1}{2}-y_7\right)\mathbf{a}_2+\left(\frac{1}{2}-z_7\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_7\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_7\right)c\,\hat{\mathbf{z}} & (8d)\ \text{Be}\\
\mathbf{B}_{29} & = & -x_7\,\mathbf{a}_1-y_7\,\mathbf{a}_2-z_7\,\mathbf{a}_3 & = & -x_7a\,\hat{\mathbf{x}}-y_7b\,\hat{\mathbf{y}}-z_7c\,\hat{\mathbf{z}} & (8d)\ \text{Be}\\
\mathbf{B}_{30} & = & \left(\frac{1}{2}+x_7\right)\mathbf{a}_1+y_7\,\mathbf{a}_2+\left(\frac{1}{2}-z_7\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{x}}+y_7b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_7\right)c\,\hat{\mathbf{z}} & (8d)\ \text{Be}\\
\mathbf{B}_{31} & = & x_7\,\mathbf{a}_1+\left(\frac{1}{2}-y_7\right)\mathbf{a}_2+z_7\,\mathbf{a}_3 & = & x_7a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_7\right)b\,\hat{\mathbf{y}}+z_7c\,\hat{\mathbf{z}} & (8d)\ \text{Be}\\
\mathbf{B}_{32} & = & \left(\frac{1}{2}-x_7\right)\mathbf{a}_1+\left(\frac{1}{2}+y_7\right)\mathbf{a}_2+\left(\frac{1}{2}+z_7\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_7\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_7\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}} & (8d)\ \text{Be}\\
\mathbf{B}_{33} & = & x_8\,\mathbf{a}_1+y_8\,\mathbf{a}_2+z_8\,\mathbf{a}_3 & = & x_8a\,\hat{\mathbf{x}}+y_8b\,\hat{\mathbf{y}}+z_8c\,\hat{\mathbf{z}} & (8d)\ \text{Na}\\
\mathbf{B}_{34} & = & \left(\frac{1}{2}-x_8\right)\mathbf{a}_1-y_8\,\mathbf{a}_2+\left(\frac{1}{2}+z_8\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_8\right)a\,\hat{\mathbf{x}}-y_8b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_8\right)c\,\hat{\mathbf{z}} & (8d)\ \text{Na}\\
\mathbf{B}_{35} & = & -x_8\,\mathbf{a}_1+\left(\frac{1}{2}+y_8\right)\mathbf{a}_2-z_8\,\mathbf{a}_3 & = & -x_8a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_8\right)b\,\hat{\mathbf{y}}-z_8c\,\hat{\mathbf{z}} & (8d)\ \text{Na}\\
\mathbf{B}_{36} & = & \left(\frac{1}{2}+x_8\right)\mathbf{a}_1+\left(\frac{1}{2}-y_8\right)\mathbf{a}_2+\left(\frac{1}{2}-z_8\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_8\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_8\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_8\right)c\,\hat{\mathbf{z}} & (8d)\ \text{Na}\\
\mathbf{B}_{37} & = & -x_8\,\mathbf{a}_1-y_8\,\mathbf{a}_2-z_8\,\mathbf{a}_3 & = & -x_8a\,\hat{\mathbf{x}}-y_8b\,\hat{\mathbf{y}}-z_8c\,\hat{\mathbf{z}} & (8d)\ \text{Na}\\
\mathbf{B}_{38} & = & \left(\frac{1}{2}+x_8\right)\mathbf{a}_1+y_8\,\mathbf{a}_2+\left(\frac{1}{2}-z_8\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_8\right)a\,\hat{\mathbf{x}}+y_8b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_8\right)c\,\hat{\mathbf{z}} & (8d)\ \text{Na}\\
\mathbf{B}_{39} & = & x_8\,\mathbf{a}_1+\left(\frac{1}{2}-y_8\right)\mathbf{a}_2+z_8\,\mathbf{a}_3 & = & x_8a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_8\right)b\,\hat{\mathbf{y}}+z_8c\,\hat{\mathbf{z}} & (8d)\ \text{Na}\\
\mathbf{B}_{40} & = & \left(\frac{1}{2}-x_8\right)\mathbf{a}_1+\left(\frac{1}{2}+y_8\right)\mathbf{a}_2+\left(\frac{1}{2}+z_8\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_8\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_8\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_8\right)c\,\hat{\mathbf{z}} & (8d)\ \text{Na}\\
\mathbf{B}_{41} & = & x_9\,\mathbf{a}_1+y_9\,\mathbf{a}_2+z_9\,\mathbf{a}_3 & = & x_9a\,\hat{\mathbf{x}}+y_9b\,\hat{\mathbf{y}}+z_9c\,\hat{\mathbf{z}} & (8d)\ \text{O V}\\
\mathbf{B}_{42} & = & \left(\frac{1}{2}-x_9\right)\mathbf{a}_1-y_9\,\mathbf{a}_2+\left(\frac{1}{2}+z_9\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_9\right)a\,\hat{\mathbf{x}}-y_9b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_9\right)c\,\hat{\mathbf{z}} & (8d)\ \text{O V}\\
\mathbf{B}_{43} & = & -x_9\,\mathbf{a}_1+\left(\frac{1}{2}+y_9\right)\mathbf{a}_2-z_9\,\mathbf{a}_3 & = & -x_9a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_9\right)b\,\hat{\mathbf{y}}-z_9c\,\hat{\mathbf{z}} & (8d)\ \text{O V}\\
\mathbf{B}_{44} & = & \left(\frac{1}{2}+x_9\right)\mathbf{a}_1+\left(\frac{1}{2}-y_9\right)\mathbf{a}_2+\left(\frac{1}{2}-z_9\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_9\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_9\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_9\right)c\,\hat{\mathbf{z}} & (8d)\ \text{O V}\\
\mathbf{B}_{45} & = & -x_9\,\mathbf{a}_1-y_9\,\mathbf{a}_2-z_9\,\mathbf{a}_3 & = & -x_9a\,\hat{\mathbf{x}}-y_9b\,\hat{\mathbf{y}}-z_9c\,\hat{\mathbf{z}} & (8d)\ \text{O V}\\
\mathbf{B}_{46} & = & \left(\frac{1}{2}+x_9\right)\mathbf{a}_1+y_9\,\mathbf{a}_2+\left(\frac{1}{2}-z_9\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_9\right)a\,\hat{\mathbf{x}}+y_9b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_9\right)c\,\hat{\mathbf{z}} & (8d)\ \text{O V}\\
\mathbf{B}_{47} & = & x_9\,\mathbf{a}_1+\left(\frac{1}{2}-y_9\right)\mathbf{a}_2+z_9\,\mathbf{a}_3 & = & x_9a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_9\right)b\,\hat{\mathbf{y}}+z_9c\,\hat{\mathbf{z}} & (8d)\ \text{O V}\\
\mathbf{B}_{48} & = & \left(\frac{1}{2}-x_9\right)\mathbf{a}_1+\left(\frac{1}{2}+y_9\right)\mathbf{a}_2+\left(\frac{1}{2}+z_9\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_9\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_9\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_9\right)c\,\hat{\mathbf{z}} & (8d)\ \text{O V}\\
\mathbf{B}_{49} & = & x_{10}\,\mathbf{a}_1+y_{10}\,\mathbf{a}_2+z_{10}\,\mathbf{a}_3 & = & x_{10}a\,\hat{\mathbf{x}}+y_{10}b\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}} & (8d)\ \text{O VI}\\
\mathbf{B}_{50} & = & \left(\frac{1}{2}-x_{10}\right)\mathbf{a}_1-y_{10}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{10}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{x}}-y_{10}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{10}\right)c\,\hat{\mathbf{z}} & (8d)\ \text{O VI}\\
\mathbf{B}_{51} & = & -x_{10}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{10}\right)\mathbf{a}_2-z_{10}\,\mathbf{a}_3 & = & -x_{10}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{10}\right)b\,\hat{\mathbf{y}}-z_{10}c\,\hat{\mathbf{z}} & (8d)\ \text{O VI}\\
\mathbf{B}_{52} & = & \left(\frac{1}{2}+x_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{10}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{10}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{10}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{10}\right)c\,\hat{\mathbf{z}} & (8d)\ \text{O VI}\\
\mathbf{B}_{53} & = & -x_{10}\,\mathbf{a}_1-y_{10}\,\mathbf{a}_2-z_{10}\,\mathbf{a}_3 & = & -x_{10}a\,\hat{\mathbf{x}}-y_{10}b\,\hat{\mathbf{y}}-z_{10}c\,\hat{\mathbf{z}} & (8d)\ \text{O VI}\\
\mathbf{B}_{54} & = & \left(\frac{1}{2}+x_{10}\right)\mathbf{a}_1+y_{10}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{10}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{x}}+y_{10}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{10}\right)c\,\hat{\mathbf{z}} & (8d)\ \text{O VI}\\
\mathbf{B}_{55} & = & x_{10}\,\mathbf{a}_1+\left(\frac{1}{2}-y_{10}\right)\mathbf{a}_2+z_{10}\,\mathbf{a}_3 & = & x_{10}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{10}\right)b\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}} & (8d)\ \text{O VI}\\
\mathbf{B}_{56} & = & \left(\frac{1}{2}-x_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{10}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{10}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{10}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{10}\right)c\,\hat{\mathbf{z}} & (8d)\ \text{O VI}\\
\mathbf{B}_{57} & = & x_{11}\,\mathbf{a}_1+y_{11}\,\mathbf{a}_2+z_{11}\,\mathbf{a}_3 & = & x_{11}a\,\hat{\mathbf{x}}+y_{11}b\,\hat{\mathbf{y}}+z_{11}c\,\hat{\mathbf{z}} & (8d)\ \text{O VII}
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\begin{array}{lcllcll}
\mathbf{B}_{58} & = & \left(\frac{1}{2}-x_{11}\right)\mathbf{a}_1-y_{11}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{11}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{11}\right)a\,\hat{\mathbf{x}}-y_{11}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{11}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O VII}\\
\mathbf{B}_{59} & = & -x_{11}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{11}\right)\mathbf{a}_2-z_{11}\,\mathbf{a}_3 & = & -x_{11}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{11}\right)b\,\hat{\mathbf{y}}-z_{11}c\,\hat{\mathbf{z}} & (8d) & \text{O VII}\\
\mathbf{B}_{60} & = & \left(\frac{1}{2}+x_{11}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{11}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{11}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{11}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{11}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O VII}\\
\mathbf{B}_{61} & = & -x_{11}\,\mathbf{a}_1-y_{11}\,\mathbf{a}_2-z_{11}\,\mathbf{a}_3 & = & -x_{11}a\,\hat{\mathbf{x}}-y_{11}b\,\hat{\mathbf{y}}-z_{11}c\,\hat{\mathbf{z}} & (8d) & \text{O VII}\\
\mathbf{B}_{62} & = & \left(\frac{1}{2}+x_{11}\right)\mathbf{a}_1+y_{11}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{11}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{11}\right)a\,\hat{\mathbf{x}}+y_{11}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{11}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O VII}\\
\mathbf{B}_{63} & = & x_{11}\,\mathbf{a}_1+\left(\frac{1}{2}-y_{11}\right)\mathbf{a}_2+z_{11}\,\mathbf{a}_3 & = & x_{11}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{11}\right)b\,\hat{\mathbf{y}}+z_{11}c\,\hat{\mathbf{z}} & (8d) & \text{O VII}\\
\mathbf{B}_{64} & = & \left(\frac{1}{2}-x_{11}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{11}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{11}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{11}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{11}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O VII}\\
\mathbf{B}_{65} & = & x_{12}\,\mathbf{a}_1+y_{12}\,\mathbf{a}_2+z_{12}\,\mathbf{a}_3 & = & x_{12}a\,\hat{\mathbf{x}}+y_{12}b\,\hat{\mathbf{y}}+z_{12}c\,\hat{\mathbf{z}} & (8d) & \text{O VIII}\\
\mathbf{B}_{66} & = & \left(\frac{1}{2}-x_{12}\right)\mathbf{a}_1-y_{12}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{12}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{12}\right)a\,\hat{\mathbf{x}}-y_{12}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{12}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O VIII}\\
\mathbf{B}_{67} & = & -x_{12}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{12}\right)\mathbf{a}_2-z_{12}\,\mathbf{a}_3 & = & -x_{12}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{12}\right)b\,\hat{\mathbf{y}}-z_{12}c\,\hat{\mathbf{z}} & (8d) & \text{O VIII}\\
\mathbf{B}_{68} & = & \left(\frac{1}{2}+x_{12}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{12}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{12}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{12}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{12}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{12}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O VIII}\\
\mathbf{B}_{69} & = & -x_{12}\,\mathbf{a}_1-y_{12}\,\mathbf{a}_2-z_{12}\,\mathbf{a}_3 & = & -x_{12}a\,\hat{\mathbf{x}}-y_{12}b\,\hat{\mathbf{y}}-z_{12}c\,\hat{\mathbf{z}} & (8d) & \text{O VIII}\\
\mathbf{B}_{70} & = & \left(\frac{1}{2}+x_{12}\right)\mathbf{a}_1+y_{12}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{12}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{12}\right)a\,\hat{\mathbf{x}}+y_{12}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{12}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O VIII}\\
\mathbf{B}_{71} & = & x_{12}\,\mathbf{a}_1+\left(\frac{1}{2}-y_{12}\right)\mathbf{a}_2+z_{12}\,\mathbf{a}_3 & = & x_{12}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{12}\right)b\,\hat{\mathbf{y}}+z_{12}c\,\hat{\mathbf{z}} & (8d) & \text{O VIII}\\
\mathbf{B}_{72} & = & \left(\frac{1}{2}-x_{12}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{12}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{12}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{12}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{12}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{12}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O VIII}\\
\mathbf{B}_{73} & = & x_{13}\,\mathbf{a}_1+y_{13}\,\mathbf{a}_2+z_{13}\,\mathbf{a}_3 & = & x_{13}a\,\hat{\mathbf{x}}+y_{13}b\,\hat{\mathbf{y}}+z_{13}c\,\hat{\mathbf{z}} & (8d) & \text{O IX}\\
\mathbf{B}_{74} & = & \left(\frac{1}{2}-x_{13}\right)\mathbf{a}_1-y_{13}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{13}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{13}\right)a\,\hat{\mathbf{x}}-y_{13}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{13}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O IX}\\
\mathbf{B}_{75} & = & -x_{13}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{13}\right)\mathbf{a}_2-z_{13}\,\mathbf{a}_3 & = & -x_{13}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{13}\right)b\,\hat{\mathbf{y}}-z_{13}c\,\hat{\mathbf{z}} & (8d) & \text{O IX}\\
\mathbf{B}_{76} & = & \left(\frac{1}{2}+x_{13}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{13}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{13}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{13}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{13}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{13}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O IX}\\
\mathbf{B}_{77} & = & -x_{13}\,\mathbf{a}_1-y_{13}\,\mathbf{a}_2-z_{13}\,\mathbf{a}_3 & = & -x_{13}a\,\hat{\mathbf{x}}-y_{13}b\,\hat{\mathbf{y}}-z_{13}c\,\hat{\mathbf{z}} & (8d) & \text{O IX}\\
\mathbf{B}_{78} & = & \left(\frac{1}{2}+x_{13}\right)\mathbf{a}_1+y_{13}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{13}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{13}\right)a\,\hat{\mathbf{x}}+y_{13}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{13}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O IX}\\
\mathbf{B}_{79} & = & x_{13}\,\mathbf{a}_1+\left(\frac{1}{2}-y_{13}\right)\mathbf{a}_2+z_{13}\,\mathbf{a}_3 & = & x_{13}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{13}\right)b\,\hat{\mathbf{y}}+z_{13}c\,\hat{\mathbf{z}} & (8d) & \text{O IX}\\
\mathbf{B}_{80} & = & \left(\frac{1}{2}-x_{13}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{13}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{13}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{13}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{13}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{13}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O IX}\\
\mathbf{B}_{81} & = & x_{14}\,\mathbf{a}_1+y_{14}\,\mathbf{a}_2+z_{14}\,\mathbf{a}_3 & = & x_{14}a\,\hat{\mathbf{x}}+y_{14}b\,\hat{\mathbf{y}}+z_{14}c\,\hat{\mathbf{z}} & (8d) & \text{O X}\\
\mathbf{B}_{82} & = & \left(\frac{1}{2}-x_{14}\right)\mathbf{a}_1-y_{14}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{14}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{14}\right)a\,\hat{\mathbf{x}}-y_{14}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{14}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O X}\\
\mathbf{B}_{83} & = & -x_{14}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{14}\right)\mathbf{a}_2-z_{14}\,\mathbf{a}_3 & = & -x_{14}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{14}\right)b\,\hat{\mathbf{y}}-z_{14}c\,\hat{\mathbf{z}} & (8d) & \text{O X}\\
\mathbf{B}_{84} & = & \left(\frac{1}{2}+x_{14}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{14}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{14}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{14}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{14}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{14}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O X}\\
\mathbf{B}_{85} & = & -x_{14}\,\mathbf{a}_1-y_{14}\,\mathbf{a}_2-z_{14}\,\mathbf{a}_3 & = & -x_{14}a\,\hat{\mathbf{x}}-y_{14}b\,\hat{\mathbf{y}}-z_{14}c\,\hat{\mathbf{z}} & (8d) & \text{O X}\\
\mathbf{B}_{86} & = & \left(\frac{1}{2}+x_{14}\right)\mathbf{a}_1+y_{14}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{14}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{14}\right)a\,\hat{\mathbf{x}}+y_{14}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{14}\right)c\,\hat{\mathbf{z}} & (8d) & \text{O X}\\
\mathbf{B}_{87} & = & x_{14}\,\mathbf{a}_1+\left(\frac{1}{2}-y_{14}\right)\mathbf{a}_2+z_{14}\,\mathbf{a}_3 & = & x_{14}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{14}\right)b\,\hat{\mathbf{y}}+z_{14}c\,\hat{\mathbf{z}} & (8d) & \text{O X}
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$\mathbf{B}_{88}$	=	$\left(\frac{1}{2}-x_{14}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{14}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{14}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{14}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{14}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{14}\right)c\,\hat{\mathbf{z}}$	(8*d*)	O X
$\mathbf{B}_{89}$	=	$x_{15}\,\mathbf{a}_1+y_{15}\,\mathbf{a}_2+z_{15}\,\mathbf{a}_3$	=	$x_{15}a\,\hat{\mathbf{x}}+y_{15}b\,\hat{\mathbf{y}}+z_{15}c\,\hat{\mathbf{z}}$	(8*d*)	Si I
$\mathbf{B}_{90}$	=	$\left(\frac{1}{2}-x_{15}\right)\mathbf{a}_1-y_{15}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{15}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{15}\right)a\,\hat{\mathbf{x}}-y_{15}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{15}\right)c\,\hat{\mathbf{z}}$	(8*d*)	Si I
$\mathbf{B}_{91}$	=	$-x_{15}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{15}\right)\mathbf{a}_2-z_{15}\,\mathbf{a}_3$	=	$-x_{15}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{15}\right)b\,\hat{\mathbf{y}}-z_{15}c\,\hat{\mathbf{z}}$	(8*d*)	Si I
$\mathbf{B}_{92}$	=	$\left(\frac{1}{2}+x_{15}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{15}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{15}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{15}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{15}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{15}\right)c\,\hat{\mathbf{z}}$	(8*d*)	Si I
$\mathbf{B}_{93}$	=	$-x_{15}\,\mathbf{a}_1-y_{15}\,\mathbf{a}_2-z_{15}\,\mathbf{a}_3$	=	$-x_{15}a\,\hat{\mathbf{x}}-y_{15}b\,\hat{\mathbf{y}}-z_{15}c\,\hat{\mathbf{z}}$	(8*d*)	Si I
$\mathbf{B}_{94}$	=	$\left(\frac{1}{2}+x_{15}\right)\mathbf{a}_1+y_{15}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{15}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{15}\right)a\,\hat{\mathbf{x}}+y_{15}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{15}\right)c\,\hat{\mathbf{z}}$	(8*d*)	Si I
$\mathbf{B}_{95}$	=	$x_{15}\,\mathbf{a}_1+\left(\frac{1}{2}-y_{15}\right)\mathbf{a}_2+z_{15}\,\mathbf{a}_3$	=	$x_{15}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{15}\right)b\,\hat{\mathbf{y}}+z_{15}c\,\hat{\mathbf{z}}$	(8*d*)	Si I
$\mathbf{B}_{96}$	=	$\left(\frac{1}{2}-x_{15}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{15}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{15}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{15}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{15}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{15}\right)c\,\hat{\mathbf{z}}$	(8*d*)	Si I
$\mathbf{B}_{97}$	=	$x_{16}\,\mathbf{a}_1+y_{16}\,\mathbf{a}_2+z_{16}\,\mathbf{a}_3$	=	$x_{16}a\,\hat{\mathbf{x}}+y_{16}b\,\hat{\mathbf{y}}+z_{16}c\,\hat{\mathbf{z}}$	(8*d*)	Si II
$\mathbf{B}_{98}$	=	$\left(\frac{1}{2}-x_{16}\right)\mathbf{a}_1-y_{16}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{16}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{16}\right)a\,\hat{\mathbf{x}}-y_{16}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{16}\right)c\,\hat{\mathbf{z}}$	(8*d*)	Si II
$\mathbf{B}_{99}$	=	$-x_{16}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{16}\right)\mathbf{a}_2-z_{16}\,\mathbf{a}_3$	=	$-x_{16}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{16}\right)b\,\hat{\mathbf{y}}-z_{16}c\,\hat{\mathbf{z}}$	(8*d*)	Si II
$\mathbf{B}_{100}$	=	$\left(\frac{1}{2}+x_{16}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{16}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{16}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{16}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{16}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{16}\right)c\,\hat{\mathbf{z}}$	(8*d*)	Si II
$\mathbf{B}_{101}$	=	$-x_{16}\,\mathbf{a}_1-y_{16}\,\mathbf{a}_2-z_{16}\,\mathbf{a}_3$	=	$-x_{16}a\,\hat{\mathbf{x}}-y_{16}b\,\hat{\mathbf{y}}-z_{16}c\,\hat{\mathbf{z}}$	(8*d*)	Si II
$\mathbf{B}_{102}$	=	$\left(\frac{1}{2}+x_{16}\right)\mathbf{a}_1+y_{16}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{16}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{16}\right)a\,\hat{\mathbf{x}}+y_{16}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{16}\right)c\,\hat{\mathbf{z}}$	(8*d*)	Si II
$\mathbf{B}_{103}$	=	$x_{16}\,\mathbf{a}_1+\left(\frac{1}{2}-y_{16}\right)\mathbf{a}_2+z_{16}\,\mathbf{a}_3$	=	$x_{16}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{16}\right)b\,\hat{\mathbf{y}}+z_{16}c\,\hat{\mathbf{z}}$	(8*d*)	Si II
$\mathbf{B}_{104}$	=	$\left(\frac{1}{2}-x_{16}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{16}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{16}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{16}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{16}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{16}\right)c\,\hat{\mathbf{z}}$	(8*d*)	Si II
$\mathbf{B}_{105}$	=	$x_{17}\,\mathbf{a}_1+y_{17}\,\mathbf{a}_2+z_{17}\,\mathbf{a}_3$	=	$x_{17}a\,\hat{\mathbf{x}}+y_{17}b\,\hat{\mathbf{y}}+z_{17}c\,\hat{\mathbf{z}}$	(8*d*)	Si III
$\mathbf{B}_{106}$	=	$\left(\frac{1}{2}-x_{17}\right)\mathbf{a}_1-y_{17}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{17}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{17}\right)a\,\hat{\mathbf{x}}-y_{17}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{17}\right)c\,\hat{\mathbf{z}}$	(8*d*)	Si III
$\mathbf{B}_{107}$	=	$-x_{17}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{17}\right)\mathbf{a}_2-z_{17}\,\mathbf{a}_3$	=	$-x_{17}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{17}\right)b\,\hat{\mathbf{y}}-z_{17}c\,\hat{\mathbf{z}}$	(8*d*)	Si III
$\mathbf{B}_{108}$	=	$\left(\frac{1}{2}+x_{17}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{17}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{17}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{17}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{17}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{17}\right)c\,\hat{\mathbf{z}}$	(8*d*)	Si III
$\mathbf{B}_{109}$	=	$-x_{17}\,\mathbf{a}_1-y_{17}\,\mathbf{a}_2-z_{17}\,\mathbf{a}_3$	=	$-x_{17}a\,\hat{\mathbf{x}}-y_{17}b\,\hat{\mathbf{y}}-z_{17}c\,\hat{\mathbf{z}}$	(8*d*)	Si III
$\mathbf{B}_{110}$	=	$\left(\frac{1}{2}+x_{17}\right)\mathbf{a}_1+y_{17}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{17}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{17}\right)a\,\hat{\mathbf{x}}+y_{17}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{17}\right)c\,\hat{\mathbf{z}}$	(8*d*)	Si III
$\mathbf{B}_{111}$	=	$x_{17}\,\mathbf{a}_1+\left(\frac{1}{2}-y_{17}\right)\mathbf{a}_2+z_{17}\,\mathbf{a}_3$	=	$x_{17}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{17}\right)b\,\hat{\mathbf{y}}+z_{17}c\,\hat{\mathbf{z}}$	(8*d*)	Si III
$\mathbf{B}_{112}$	=	$\left(\frac{1}{2}-x_{17}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{17}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{17}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{17}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{17}\right)b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{17}\right)c\,\hat{\mathbf{z}}$	(8*d*)	Si III

References:

- G. Diego Gatta, N. Rotiroti, G. J. McIntyre, A. Guastoni, and F. Nestola, *New insights into the crystal chemistry of epididymite and eudidymite from Malosa, Malawi: A single-crystal neutron diffraction study*, Am. Mineral. **93**, 1158–1165 (2008), doi:10.2138/am.2008.2965.
- T. Ito, *The Structure of Epididymite ($HNaBeSi_3O_8$)*, Zeitschrift für Kristallographie - Crystalline Materials **88**, 142–149 (1934), doi:10.1524/zkri.1934.88.1.142.
- C. Gottfried and F. Schossberger, eds., *Strukturbericht Band III 1933-1935* (Akademische Verlagsgesellschaft M. B. H.,

Leipzig, 1937).

Found in:

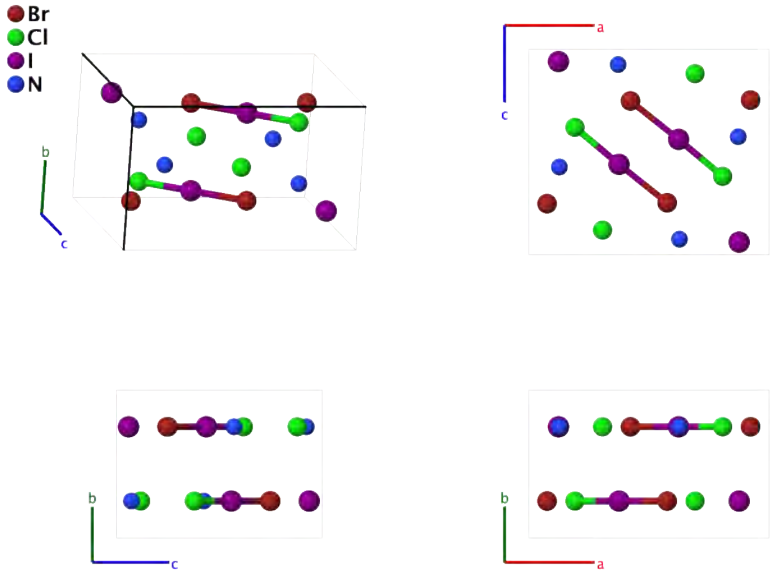
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1649
- POSCAR: pp. 1650

NH_4ClBrI ($F5_{14}$) Structure: ABCD_oP16_62_c_c_c_c

http://aflow.org/prototype-encyclopedia/ABCD_oP16_62_c_c_c_c

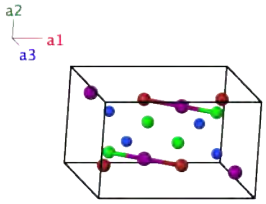


Prototype	:	$ClBrI(NH_4)$
AFLOW prototype label	:	ABCD_oP16_62_c_c_c_c
***Strukturbericht* designation**	:	$F5_{14}$
Pearson symbol	:	oP16
Space group number	:	62
Space group symbol	:	*Pnma*
AFLOW prototype command	:	`aflow --proto=ABCD_oP16_62_c_c_c_c --params=`$a, b/a, c/a, x_1, z_1, x_2, z_2, x_3, z_3, x_4, z_4$

- (Gottfried, 1937) originally described NH_4ClBrI as isostructural with NH_4I_3, $D0_{16}$, based on the early work of (Mooney, 1935). However, (Gottfried, 1940) used the newer work of (Mooney, 1937) to create a separate *Strukturbericht* designation for NH_4ClBrI, $F5_{14}$, even though this structure is still obviously related to NH_4I_3.
- The positions of the hydrogen atoms in the NH_4 were not determined, so we only provide the positions of the nitrogen atoms (labeled as NH_4).
- The original structure is given in the *Pmcn* setting of space group #62, which was transformed to the standard *Pnma* setting by FINDSYM. However, note that the coordinate transformations for the (4*c*) Wyckoff positions in (Gottfried, 1940) are erroneous, they give a structure in space group *Pmmn* #59. (Gottfried, 1937) has the correct tranformations in the $D0_{16}$ section.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(4c)	Br
$\mathbf{B}_2$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	(4c)	Br
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	(4c)	Br
$\mathbf{B}_4$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right) c\,\hat{\mathbf{z}}$	(4c)	Br
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4c)	Cl
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	(4c)	Cl
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(4c)	Cl
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right) c\,\hat{\mathbf{z}}$	(4c)	Cl
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(4c)	I
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	(4c)	I
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(4c)	I
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right) c\,\hat{\mathbf{z}}$	(4c)	I
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(4c)	NH_4
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	(4c)	NH_4
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(4c)	NH_4
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\,\hat{\mathbf{z}}$	(4c)	NH_4

References:

- R. C. L. Mooney, *The Crystal Structure of Ammonium Chlorobromoiodide and the Configuration of the Chlorobromoiodide Group*, Zeitschrift für Kristallographie - Crystalline Materials **98**, 324–333 (1938), doi:10.1524/zkri.1938.98.1.324.
- R. C. L. Mooney, *The Crystal Structure of Ammonium Chlorobromoiodide*, Phys. Rev. **47**, 807–808 (1935), doi:10.1103/PhysRev.47.785.
- C. Gottfried and F. Schossberger, eds., *Strukturbericht Band III 1933-1935* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Found in:

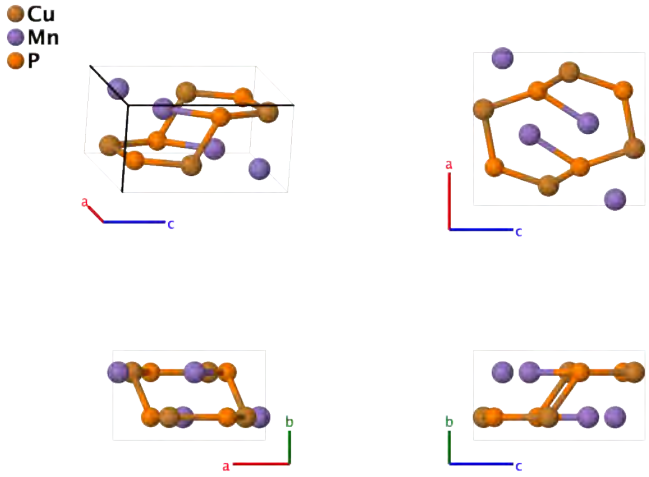
- C. Gottfried, ed., *Strukturbericht Band V 1937* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1940).

Geometry files:

- CIF: pp. 1650
- POSCAR: pp. 1651

MnCuP Structure: ABC_oP12_62_c_c_c

http://aflow.org/prototype-encyclopedia/ABC_oP12_62_c_c_c



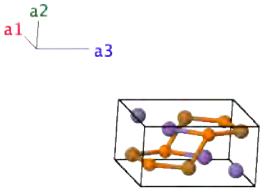
Prototype	:	CuMnP
AFLOW prototype label	:	ABC_oP12_62_c_c_c
***Strukturbericht* designation**	:	None
Pearson symbol	:	oP12
Space group number	:	62
Space group symbol	:	*Pnma*
AFLOW prototype command	:	`aflow --proto=ABC_oP12_62_c_c_c --params=`$a, b/a, c/a, x_1, z_1, x_2, z_2, x_3, z_3$

Other compounds with this structure

- CuMnAs and $CuMnP_xAs_{1-x}$

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	($4c$)	Cu
$\mathbf{B}_2$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	($4c$)	Cu
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	($4c$)	Cu

$\mathbf{B}_4$	=	$\left(\frac{1}{2}+x_1\right)\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+\left(\frac{1}{2}-z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_1\right)a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_1\right)c\,\hat{\mathbf{z}}$	(4*c*)	Cu
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+z_2\,\mathbf{a}_3$	=	$x_2a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+z_2c\,\hat{\mathbf{z}}$	(4*c*)	Mn
$\mathbf{B}_6$	=	$\left(\frac{1}{2}-x_2\right)\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2+\left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_2\right)a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_2\right)c\,\hat{\mathbf{z}}$	(4*c*)	Mn
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2-z_2\,\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}-z_2c\,\hat{\mathbf{z}}$	(4*c*)	Mn
$\mathbf{B}_8$	=	$\left(\frac{1}{2}+x_2\right)\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+\left(\frac{1}{2}-z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_2\right)a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_2\right)c\,\hat{\mathbf{z}}$	(4*c*)	Mn
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+z_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+z_3c\,\hat{\mathbf{z}}$	(4*c*)	P
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2}-x_3\right)\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2+\left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_3\right)c\,\hat{\mathbf{z}}$	(4*c*)	P
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2-z_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}-z_3c\,\hat{\mathbf{z}}$	(4*c*)	P
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2}+x_3\right)\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+\left(\frac{1}{2}-z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_3\right)c\,\hat{\mathbf{z}}$	(4*c*)	P

References:

- J. Mündelein and H.-U. Schuster, *Darstellung und Kristallstruktur der Verbindungen MnCuX (X = P, As, P_xAs_{1-x})*, Z. Naturforsch. B **47**, 925–928 (1992), doi:10.1515/znb-1992-0705.

Found in:

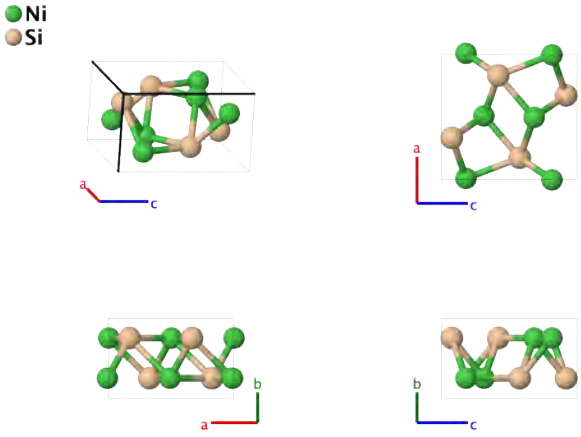
- F. Máca, J. Mašek, O. Stelmakhovych, X. Martí, H. Reichlová, K. Uhlířová, P. Beran, P. Wadley, V. Novák, and T. Jungwirth, *Room-temperature antiferromagnetism in CuMnAs*, J. Magn. Magn. Mater. **324**, 1606–1612 (2012), doi:10.1016/j.jmmm.2011.12.017.

Geometry files:

- CIF: pp. 1651
- POSCAR: pp. 1651

η-NiSi (B_d) Structure: AB_oP8_62_c_c

http://aflow.org/prototype-encyclopedia/AB_oP8_62_c_c.NiSi

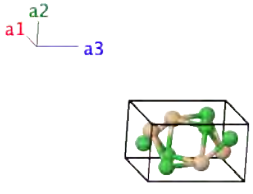


Prototype	:	NiSi
AFLOW prototype label	:	AB_oP8_62_c_c
***Strukturbericht* designation**	:	B_d
Pearson symbol	:	oP8
Space group number	:	62
Space group symbol	:	*Pnma*
AFLOW prototype command	:	`aflow --proto=AB_oP8_62_c_c --params=`$a, b/a, c/a, x_1, z_1, x_2, z_2$

- (Toman, 1951) described this in the *Pbnm* setting of space group #62. We have used FINDSYM to transform this to the standard *Pnma* setting. Note that GeS (AB_oP8_62_c_c, GeS), MnP (AB_oP8_62_c_c, MnP), FeB (AB_oP8_62_c_c, FeB), and SnS (AB_oP8_62_c_c, SnS) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters (`--params`) specified in their corresponding CIF files.

Simple Orthorhombic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	($4c$)	Ni
$\mathbf{B}_2$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	($4c$)	Ni
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	($4c$)	Ni

$$\mathbf{B}_4 = \left(\tfrac{1}{2}+x_1\right)\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2 + \left(\tfrac{1}{2}-z_1\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_1\right)a\,\hat{\mathbf{x}} + \tfrac{1}{4}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_1\right)c\,\hat{\mathbf{z}} \quad (4c) \quad \text{Ni}$$

$$\mathbf{B}_5 = x_2\,\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2 + z_2\,\mathbf{a}_3 = x_2 a\,\hat{\mathbf{x}} + \tfrac{1}{4}b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}} \quad (4c) \quad \text{Si}$$

$$\mathbf{B}_6 = \left(\tfrac{1}{2}-x_2\right)\mathbf{a}_1 + \tfrac{3}{4}\mathbf{a}_2 + \left(\tfrac{1}{2}+z_2\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_2\right)a\,\hat{\mathbf{x}} + \tfrac{3}{4}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_2\right)c\,\hat{\mathbf{z}} \quad (4c) \quad \text{Si}$$

$$\mathbf{B}_7 = -x_2\,\mathbf{a}_1 + \tfrac{3}{4}\mathbf{a}_2 - z_2\,\mathbf{a}_3 = -x_2 a\,\hat{\mathbf{x}} + \tfrac{3}{4}b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}} \quad (4c) \quad \text{Si}$$

$$\mathbf{B}_8 = \left(\tfrac{1}{2}+x_2\right)\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2 + \left(\tfrac{1}{2}-z_2\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_2\right)a\,\hat{\mathbf{x}} + \tfrac{1}{4}b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_2\right)c\,\hat{\mathbf{z}} \quad (4c) \quad \text{Si}$$

References:

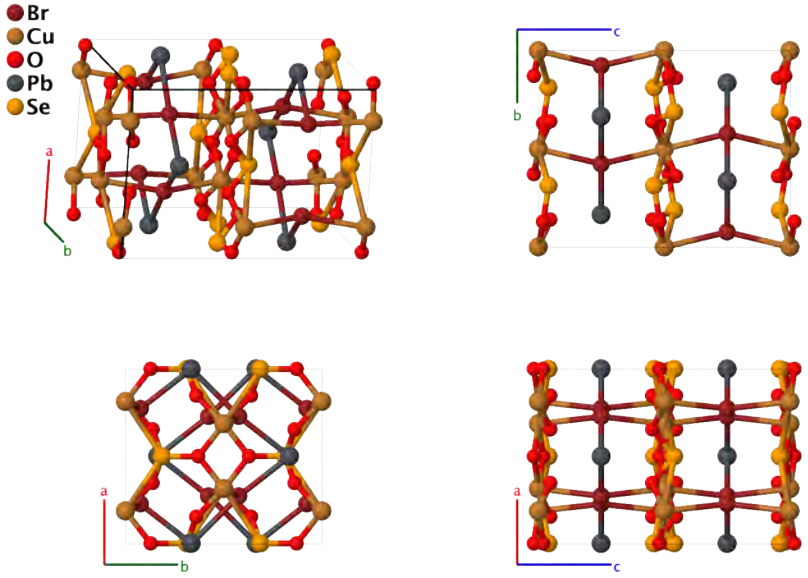
- K. Toman, *The structure of NiSi*, Acta Cryst. **4**, 462–464 (1951), doi:10.1107/S0365110X51001458.

Geometry files:

- CIF: pp. 1651
- POSCAR: pp. 1652

$Cu_2Pb(SeO_3)_2Br_2$ Structure: A2B2C6DE2_oC52_63_g_e_fh_c_f

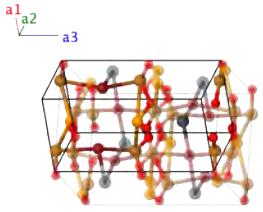
http://aflow.org/prototype-encyclopedia/A2B2C6DE2_oC52_63_g_e_fh_c_f



Prototype	:	$Br_2Cu_2O_6PbSe_2$
AFLOW prototype label	:	A2B2C6DE2_oC52_63_g_e_fh_c_f
***Strukturbericht* designation**	:	None
Pearson symbol	:	oC52
Space group number	:	63
Space group symbol	:	*Cmcm*
AFLOW prototype command	:	`aflow --proto=A2B2C6DE2_oC52_63_g_e_fh_c_f --params=`$a, b/a, c/a, y_1, x_2, y_3, z_3, y_4, z_4, x_5, y_5, x_6, y_6, z_6$

Base-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_1 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(4c)$	Pb
$\mathbf{B}_2$	=	$y_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_1 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(4c)$	Pb
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2$	=	$x_2 a\,\hat{\mathbf{x}}$	$(8e)$	Cu

$\mathbf{B}_4$	=	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(8*e*)	Cu
$\mathbf{B}_5$	=	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2$	=	$-x_2 a\,\hat{\mathbf{x}}$	(8*e*)	Cu
$\mathbf{B}_6$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(8*e*)	Cu
$\mathbf{B}_7$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(8*f*)	O I
$\mathbf{B}_8$	=	$y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$-y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(8*f*)	O I
$\mathbf{B}_9$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(8*f*)	O I
$\mathbf{B}_{10}$	=	$y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-y_3 b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(8*f*)	O I
$\mathbf{B}_{11}$	=	$-y_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*f*)	Se
$\mathbf{B}_{12}$	=	$y_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$-y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(8*f*)	Se
$\mathbf{B}_{13}$	=	$-y_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(8*f*)	Se
$\mathbf{B}_{14}$	=	$y_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-y_4 b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*f*)	Se
$\mathbf{B}_{15}$	=	$(x_5 - y_5)\,\mathbf{a}_1 + (x_5 + y_5)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(8*g*)	Br
$\mathbf{B}_{16}$	=	$(-x_5 + y_5)\,\mathbf{a}_1 + (-x_5 - y_5)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(8*g*)	Br
$\mathbf{B}_{17}$	=	$(-x_5 - y_5)\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(8*g*)	Br
$\mathbf{B}_{18}$	=	$(x_5 + y_5)\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(8*g*)	Br
$\mathbf{B}_{19}$	=	$(x_6 - y_6)\,\mathbf{a}_1 + (x_6 + y_6)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(16*h*)	O II
$\mathbf{B}_{20}$	=	$(-x_6 + y_6)\,\mathbf{a}_1 + (-x_6 - y_6)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\,\hat{\mathbf{z}}$	(16*h*)	O II
$\mathbf{B}_{21}$	=	$(-x_6 - y_6)\,\mathbf{a}_1 + (-x_6 + y_6)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\,\hat{\mathbf{z}}$	(16*h*)	O II
$\mathbf{B}_{22}$	=	$(x_6 + y_6)\,\mathbf{a}_1 + (x_6 - y_6)\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(16*h*)	O II
$\mathbf{B}_{23}$	=	$(-x_6 + y_6)\,\mathbf{a}_1 + (-x_6 - y_6)\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(16*h*)	O II
$\mathbf{B}_{24}$	=	$(x_6 - y_6)\,\mathbf{a}_1 + (x_6 + y_6)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\,\hat{\mathbf{z}}$	(16*h*)	O II
$\mathbf{B}_{25}$	=	$(x_6 + y_6)\,\mathbf{a}_1 + (x_6 - y_6)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\,\hat{\mathbf{z}}$	(16*h*)	O II
$\mathbf{B}_{26}$	=	$(-x_6 - y_6)\,\mathbf{a}_1 + (-x_6 + y_6)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(16*h*)	O II

References:

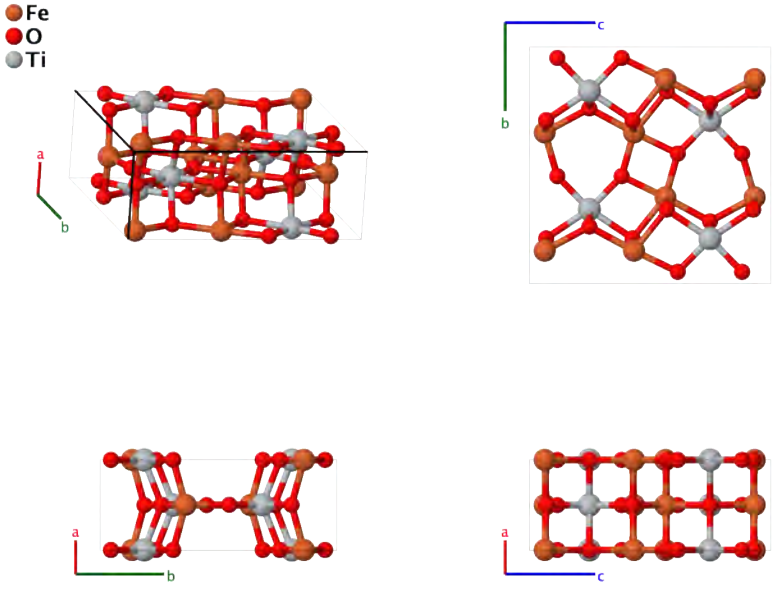
- O. I. Siidra, M. S. Kozin, W. Depmeier, R. A. Kayukov, and V. M. Kovrugin, *Copper-lead selenite bromides: a new large family of compounds partly having Cu^{2+} substructures derivable from kagome nets*, Acta Crystallogr. Sect. B Struct. Sci. **74**, 712–724 (2018), doi:10.1107/S2052520618016542.

Geometry files:

- CIF: pp. 1652
- POSCAR: pp. 1652

Pseudobrookite (Fe_2TiO_5, $E4_1$) Structure: A2B5C_oC32_63_f_c2f_c

http://aflow.org/prototype-encyclopedia/A2B5C_oC32_63_f_c2f_c



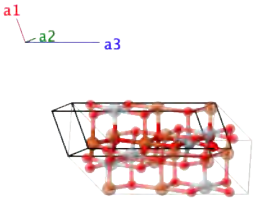
Prototype	:	Fe_2O_5Ti
AFLOW prototype label	:	A2B5C_oC32_63_f_c2f_c
***Strukturbericht* designation**	:	$E4_1$
Pearson symbol	:	oC32
Space group number	:	63
Space group symbol	:	*Cmcm*
AFLOW prototype command	:	`aflow --proto=A2B5C_oC32_63_f_c2f_c --params=`$a, b/a, c/a, y_1, y_2, y_3, z_3, y_4, z_4, y_5, z_5$

Other compounds with this structure

- $Fe_{1+x}Ti_{2-x}O_5$ and Ti_2MgO_5

Base-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_1 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(4c)$	O I
$\mathbf{B}_2$	=	$y_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_1 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(4c)$	O I
$\mathbf{B}_3$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_2 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(4c)$	Ti
$\mathbf{B}_4$	=	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_2 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(4c)$	Ti
$\mathbf{B}_5$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(8f)$	Fe
$\mathbf{B}_6$	=	$y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$-y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	$(8f)$	Fe
$\mathbf{B}_7$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	$(8f)$	Fe
$\mathbf{B}_8$	=	$y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-y_3 b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(8f)$	Fe
$\mathbf{B}_9$	=	$-y_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{10}$	=	$y_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$-y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{11}$	=	$-y_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{12}$	=	$y_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-y_4 b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{13}$	=	$-y_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{14}$	=	$y_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$-y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{15}$	=	$-y_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{16}$	=	$y_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-y_5 b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	$(8f)$	O III

References:

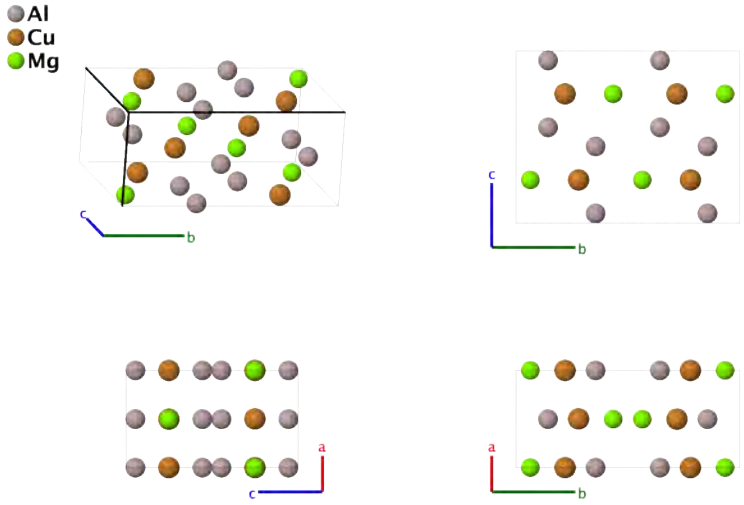
- W. Q. Guo, S. Malus, D. H. Ryan, and Z. Altounian, *Crystal structure and cation distributions in the $FeTi_2O_5$-Fe_2TiO_5 solid solution series*, J. Phys.: Condens. Matter **11**, 6337–6346 (1999), doi:10.1088/0953-8984/11/33/304.

Geometry files:

- CIF: pp. 1652
- POSCAR: pp. 1653

$MgCuAl_2$ ($E1_a$) Structure: A2BC_oC16_63_f_c_c

http://aflow.org/prototype-encyclopedia/A2BC_oC16_63_f_c_c



Prototype	:	Al_2CuMg
AFLOW prototype label	:	A2BC_oC16_63_f_c_c
***Strukturbericht* designation**	:	$E1_a$
Pearson symbol	:	oC16
Space group number	:	63
Space group symbol	:	*Cmcm*
AFLOW prototype command	:	`aflow --proto=A2BC_oC16_63_f_c_c --params=`$a, b/a, c/a, y_1, y_2, y_3, z_3$

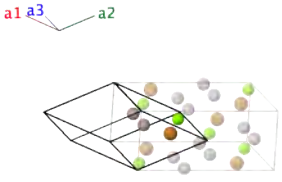
Other compounds with this structure

- $YNiAl_2$, $TaBCo_2$, $ScNiAl_2$, $LaPdIn_2$, RE-$CuPd_2$ (RE = Ce, Pr, Nd, Sm), $CaNiGa_2$, $PPdNi_2$, $EuIrSn_2$, $EuPdIn_2$, $EuPdSn_2$, $YbAuIn_2$, and $YbPdIn_2$

- This is often referred to as an 'S-phase precipitate'. It can be considered as the ternary version of the Re_3B structure.

Base-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$\mathbf{B}_1$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_1 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(4c)$	Cu
$\mathbf{B}_2$	=	$y_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_1 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(4c)$	Cu
$\mathbf{B}_3$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_2 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(4c)$	Mg
$\mathbf{B}_4$	=	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_2 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(4c)$	Mg
$\mathbf{B}_5$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(8f)$	Al
$\mathbf{B}_6$	=	$y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\,\mathbf{a}_3$	=	$-y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	$(8f)$	Al
$\mathbf{B}_7$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\,\mathbf{a}_3$	=	$y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	$(8f)$	Al
$\mathbf{B}_8$	=	$y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-y_3 b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(8f)$	Al

References:

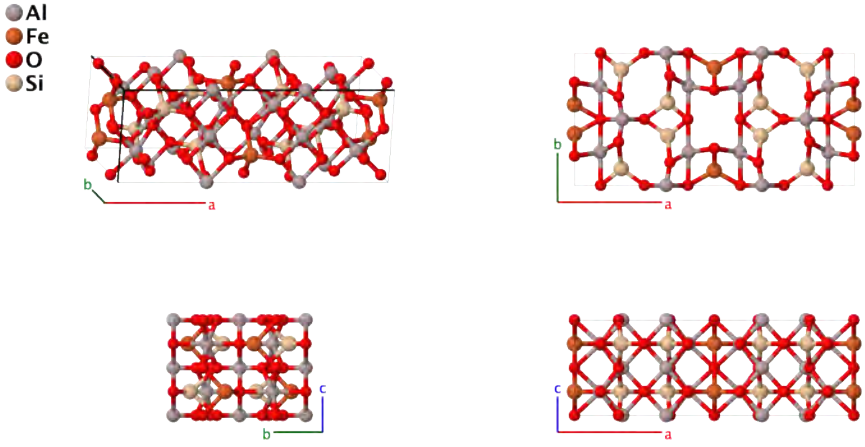
- B. Heying, R.-D. Hoffmann, and R. Pöttgen, *Structure Refinement of the S-Phase Precipitate $MgCuAl_2$*, Z. Naturforsch. B **60**, 491–494 (2005), doi:10.1515/znb-2005-0502.

Geometry files:

- CIF: pp. 1653
- POSCAR: pp. 1653

$S0_4$ (Staurolite, $Fe(OH)_2Al_4Si_2O_{10}$) (*obsolete*) Structure: A4BC12D2_oC76_63_eg_c_f3gh_g

http://aflow.org/prototype-encyclopedia/A4BC12D2_oC76_63_eg_c_f3gh_g

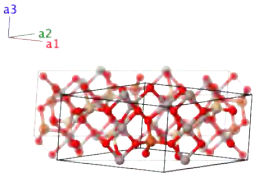


Prototype	:	$Al_4FeO_{12}Si_2$
AFLOW prototype label	:	A4BC12D2_oC76_63_eg_c_f3gh_g
***Strukturbericht* designation**	:	$S0_4$
Pearson symbol	:	oC76
Space group number	:	63
Space group symbol	:	*Cmcm*
AFLOW prototype command	:	`aflow --proto=A4BC12D2_oC76_63_eg_c_f3gh_g --params=`$a, b/a, c/a, y_1, x_2, y_3, z_3, x_4, y_4, x_5, y_5, x_6, y_6, x_7, y_7, x_8, y_8, x_9, y_9, z_9$

- This orthorhombic structure of staurolite determined by (Náray-Szabó, 1929) was given the *Strukturbericht* designation $S0_4$ by (Hermann, 1937). (Smith, 1968) showed that the structure is actually monoclinic with $\beta \approx 90°$, rather than orthorhombic and corrected the chemical composition of the mineral. Accordingly, we have marked this version of the structure as obsolete, but retain it for historical interest.
- One sixth of the oxygen sites listed here should actually be OH ions, but the positions of the hydrogen atoms were not given in (Náray-Szabó, 1929) and are labeled as O.
- We take our data from (Hermann, 1937), who presented it in the *Ccmm* representation of space group #63. We used FINDSYM to transform the structure to the standard *Cmcm* representation.
- (Hermann, 1937) also gave this the label $H5_4$ in the index.

Base-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_1 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(4c)$	Fe
$\mathbf{B}_2$	=	$y_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_1 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(4c)$	Fe
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2$	=	$x_2 a\,\hat{\mathbf{x}}$	$(8e)$	Al I
$\mathbf{B}_4$	=	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(8e)$	Al I
$\mathbf{B}_5$	=	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2$	=	$-x_2 a\,\hat{\mathbf{x}}$	$(8e)$	Al I
$\mathbf{B}_6$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(8e)$	Al I
$\mathbf{B}_7$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(8f)$	O I
$\mathbf{B}_8$	=	$y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$-y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	$(8f)$	O I
$\mathbf{B}_9$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	$(8f)$	O I
$\mathbf{B}_{10}$	=	$y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-y_3 b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(8f)$	O I
$\mathbf{B}_{11}$	=	$(x_4 - y_4)\,\mathbf{a}_1 + (x_4 + y_4)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(8g)$	Al II
$\mathbf{B}_{12}$	=	$(-x_4 + y_4)\,\mathbf{a}_1 + (-x_4 - y_4)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(8g)$	Al II
$\mathbf{B}_{13}$	=	$(-x_4 - y_4)\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(8g)$	Al II
$\mathbf{B}_{14}$	=	$(x_4 + y_4)\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(8g)$	Al II
$\mathbf{B}_{15}$	=	$(x_5 - y_5)\,\mathbf{a}_1 + (x_5 + y_5)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(8g)$	O II
$\mathbf{B}_{16}$	=	$(-x_5 + y_5)\,\mathbf{a}_1 + (-x_5 - y_5)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(8g)$	O II
$\mathbf{B}_{17}$	=	$(-x_5 - y_5)\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(8g)$	O II
$\mathbf{B}_{18}$	=	$(x_5 + y_5)\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(8g)$	O II
$\mathbf{B}_{19}$	=	$(x_6 - y_6)\,\mathbf{a}_1 + (x_6 + y_6)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(8g)$	O III
$\mathbf{B}_{20}$	=	$(-x_6 + y_6)\,\mathbf{a}_1 + (-x_6 - y_6)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(8g)$	O III
$\mathbf{B}_{21}$	=	$(-x_6 - y_6)\,\mathbf{a}_1 + (-x_6 + y_6)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + y_6 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(8g)$	O III
$\mathbf{B}_{22}$	=	$(x_6 + y_6)\,\mathbf{a}_1 + (x_6 - y_6)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} - y_6 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(8g)$	O III
$\mathbf{B}_{23}$	=	$(x_7 - y_7)\,\mathbf{a}_1 + (x_7 + y_7)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(8g)$	O IV
$\mathbf{B}_{24}$	=	$(-x_7 + y_7)\,\mathbf{a}_1 + (-x_7 - y_7)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(8g)$	O IV
$\mathbf{B}_{25}$	=	$(-x_7 - y_7)\,\mathbf{a}_1 + (-x_7 + y_7)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(8g)$	O IV
$\mathbf{B}_{26}$	=	$(x_7 + y_7)\,\mathbf{a}_1 + (x_7 - y_7)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(8g)$	O IV
$\mathbf{B}_{27}$	=	$(x_8 - y_8)\,\mathbf{a}_1 + (x_8 + y_8)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(8g)$	Si
$\mathbf{B}_{28}$	=	$(-x_8 + y_8)\,\mathbf{a}_1 + (-x_8 - y_8)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(8g)$	Si
$\mathbf{B}_{29}$	=	$(-x_8 - y_8)\,\mathbf{a}_1 + (-x_8 + y_8)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(8g)$	Si
$\mathbf{B}_{30}$	=	$(x_8 + y_8)\,\mathbf{a}_1 + (x_8 - y_8)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(8g)$	Si
$\mathbf{B}_{31}$	=	$(x_9 - y_9)\,\mathbf{a}_1 + (x_9 + y_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	$(16h)$	O V
$\mathbf{B}_{32}$	=	$(-x_9 + y_9)\,\mathbf{a}_1 + (-x_9 - y_9)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right)c\,\hat{\mathbf{z}}$	$(16h)$	O V
$\mathbf{B}_{33}$	=	$(-x_9 - y_9)\,\mathbf{a}_1 + (-x_9 + y_9)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_9\right)\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_9\right)c\,\hat{\mathbf{z}}$	$(16h)$	O V
$\mathbf{B}_{34}$	=	$(x_9 + y_9)\,\mathbf{a}_1 + (x_9 - y_9)\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}}$	$(16h)$	O V

$$\mathbf{B}_{35} = (-x_9 + y_9)\,\mathbf{a}_1 + (-x_9 - y_9)\,\mathbf{a}_2 - z_9\,\mathbf{a}_3 = -x_9 a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}} \quad (16h) \quad \text{O V}$$

$$\mathbf{B}_{36} = (x_9 - y_9)\,\mathbf{a}_1 + (x_9 + y_9)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_9\right)\mathbf{a}_3 = x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_9\right) c\,\hat{\mathbf{z}} \quad (16h) \quad \text{O V}$$

$$\mathbf{B}_{37} = (x_9 + y_9)\,\mathbf{a}_1 + (x_9 - y_9)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_3 = x_9 a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_9\right) c\,\hat{\mathbf{z}} \quad (16h) \quad \text{O V}$$

$$\mathbf{B}_{38} = (-x_9 - y_9)\,\mathbf{a}_1 + (-x_9 + y_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 = -x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}} \quad (16h) \quad \text{O V}$$

References:

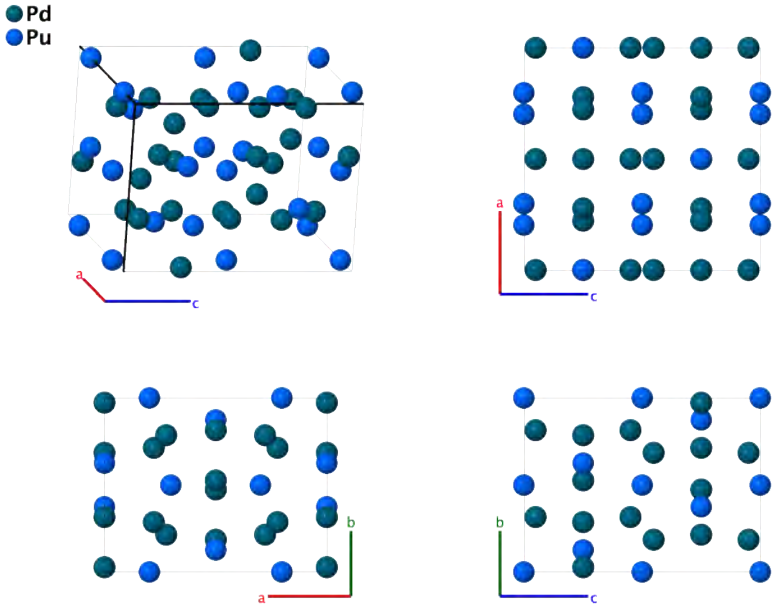
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).
- St. Náray-Szabó, *The structure of staurolite*, Zeitschrift für Kristallographie - Crystalline Materials **71**, 103–116 (1929), doi:10.1524/zkri.1929.71.1.103.
- J. D. H. Donnay and G. Donnay, *The staurolite story*, Tschermaks Min. Petr. Mitt. **31**, 1–15 (1983), doi:10.1007/BF01084757.
- J. V. Smith, *The crystal structure of staurolite*, Am. Mineral. **53**, 1139–1155 (1968).

Geometry files:

- CIF: pp. 1653
- POSCAR: pp. 1654

Pd_5Pu_3 Structure: A5B3_oC32_63_cfg_ce

http://aflow.org/prototype-encyclopedia/A5B3_oC32_63_cfg_ce



Prototype	:	Pd_5Pu_3
AFLOW prototype label	:	A5B3_oC32_63_cfg_ce
***Strukturbericht* designation**	:	None
Pearson symbol	:	oC32
Space group number	:	63
Space group symbol	:	*Cmcm*
AFLOW prototype command	:	`aflow --proto=A5B3_oC32_63_cfg_ce --params=`$a, b/a, c/a, y_1, y_2, x_3, y_4, z_4, x_5, y_5$

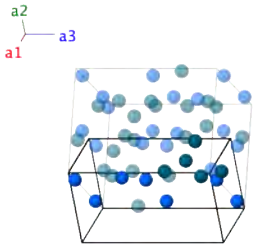
Other compounds with this structure

- Ga_5Zr_3, In_5R_3 (R = Ce, Dy, Er, Gd, Ho, La, Lu, Nd, Pr, Sm, Tb, Th, Y), Pb_5Ba_3, Pd_5R_3 (R = Sc, Y, Gd-Lu), Rh_5Zr_3, Ga_5U_3, $(Mg_xSn_{1-x})_5La_3$, Sn_5La_3, Sn_5Sr_3, and Sn_5Yb_3

- Although (Massalski, 1990) lists Pd_5Pu_3 as the prototype for many structures, it is not shown in the assessed Pd-Pu phase diagram, which is based on data from 1967.
- (Cromer, 1976) states that this phase may be isostructural with Ga_5Zr_3, but at the time of publication the exact structure of that phase had not been solved.

Base-centered Orthorhombic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2}\,b\,\hat{\mathbf{y}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,b\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_1 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(4c)$	Pd I
$\mathbf{B}_2$	=	$y_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_1 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(4c)$	Pd I
$\mathbf{B}_3$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_2 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(4c)$	Pu I
$\mathbf{B}_4$	=	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_2 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(4c)$	Pu I
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2$	=	$x_3 a\,\hat{\mathbf{x}}$	$(8e)$	Pu II
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(8e)$	Pu II
$\mathbf{B}_7$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2$	=	$-x_3 a\,\hat{\mathbf{x}}$	$(8e)$	Pu II
$\mathbf{B}_8$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(8e)$	Pu II
$\mathbf{B}_9$	=	$-y_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(8f)$	Pd II
$\mathbf{B}_{10}$	=	$y_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$-y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	$(8f)$	Pd II
$\mathbf{B}_{11}$	=	$-y_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	$(8f)$	Pd II
$\mathbf{B}_{12}$	=	$y_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-y_4 b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	$(8f)$	Pd II
$\mathbf{B}_{13}$	=	$(x_5 - y_5)\,\mathbf{a}_1 + (x_5 + y_5)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(8g)$	Pd III
$\mathbf{B}_{14}$	=	$(-x_5 + y_5)\,\mathbf{a}_1 + (-x_5 - y_5)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(8g)$	Pd III
$\mathbf{B}_{15}$	=	$(-x_5 - y_5)\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(8g)$	Pd III
$\mathbf{B}_{16}$	=	$(x_5 + y_5)\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(8g)$	Pd III

References:

- D. T. Cromer, *Plutonium-palladium Pu_3Pd_5*, Acta Crystallogr. Sect. B Struct. Sci. **32**, 1930–1932 (1976), doi:10.1107/S0567740876006778.
- T. B. Massalski, H. Okamoto, P. R. Subramanian, and L. Kacprzak, eds., *Binary Alloy Phase Diagrams*, vol. 1 (ASM International, Materials Park, Ohio, USA, 1990), 2^{nd} edn.

Found in:

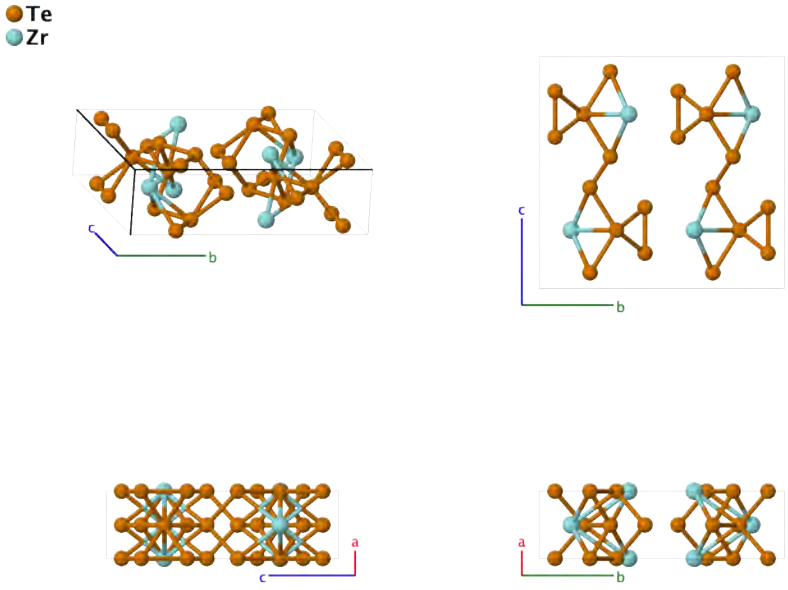
- A. Provino, N. S. Sangeetha, S. K. Dhar, V. Smetana, K. A. Gschneidner Jr., V. K. Pecharsky, P. Manfrinetti, and A.-V. Mudring, *New R_3Pd_5 Compounds (R = Sc, Y, Gd-Lu): Formation and Stability, Crystal Structure, and Antiferromagnetism*, Cryst. Growth Des. **16**, 6001–6015 (2016), doi:10.1021/acs.cgd.6b01045.

Geometry files:

- CIF: pp. 1654
- POSCAR: pp. 1654

$ZrTe_5$ Structure: A5B_oC24_63_c2f_c

http://aflow.org/prototype-encyclopedia/A5B_oC24_63_c2f_c



Prototype	:	Te_5Zr
AFLOW prototype label	:	A5B_oC24_63_c2f_c
***Strukturbericht* designation**	:	None
Pearson symbol	:	oC24
Space group number	:	63
Space group symbol	:	*Cmcm*
AFLOW prototype command	:	`aflow --proto=A5B_oC24_63_c2f_c --params=`$a, b/a, c/a, y_1, y_2, y_3, z_3, y_4, z_4$

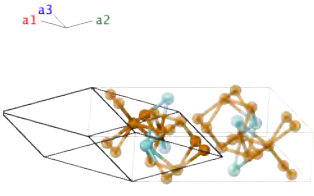
Other compounds with this structure

- $HfTe_5$

- We use the data taken at room temperature.

Base-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_1 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(4c)$	Te I
$\mathbf{B}_2$	=	$y_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_1 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(4c)$	Te I
$\mathbf{B}_3$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_2 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(4c)$	Zr
$\mathbf{B}_4$	=	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_2 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(4c)$	Zr
$\mathbf{B}_5$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(8f)$	Te II
$\mathbf{B}_6$	=	$y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$-y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	$(8f)$	Te II
$\mathbf{B}_7$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	$(8f)$	Te II
$\mathbf{B}_8$	=	$y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-y_3 b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(8f)$	Te II
$\mathbf{B}_9$	=	$-y_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(8f)$	Te III
$\mathbf{B}_{10}$	=	$y_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$-y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	$(8f)$	Te III
$\mathbf{B}_{11}$	=	$-y_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	$(8f)$	Te III
$\mathbf{B}_{12}$	=	$y_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-y_4 b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	$(8f)$	Te III

References:

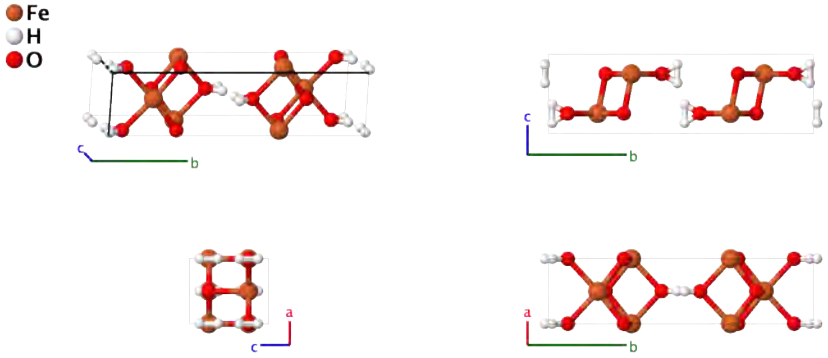
- H. Fjellvåg and A. Kjekshus, *Structural Properties of $ZrTe_5$ and $HfTe_5$ as Seen by Power Diffraction*, Solid State Commun. **60**, 91–93 (1986), doi:10.1016/0038-1098(86)90536-3.

Geometry files:

- CIF: pp. 1655
- POSCAR: pp. 1655

Lepidocrocite (γ-FeO(OH), $E0_4$) Structure: AB2C2_oC20_63_c_f_2c

http://aflow.org/prototype-encyclopedia/AB2C2_oC20_63_c_f_2c



Prototype	:	$FeHO_2$
AFLOW prototype label	:	AB2C2_oC20_63_c_f_2c
***Strukturbericht* designation**	:	$E0_4$
Pearson symbol	:	oC20
Space group number	:	63
Space group symbol	:	*Cmcm*
AFLOW prototype command	:	`aflow --proto=AB2C2_oC20_63_c_f_2c --params=`$a, b/a, c/a, y_1, y_2, y_3, y_4, z_4$

- (Gottfried, 1937) gave γ-FeO(OH) the *Strukturbericht* designation $E0_4$, but did not determine the positions of the hydrogens. Using a deuterated form of FeO(OH) (Christensen, 1982) found that the hydrogen atoms are located on the ($8f$) Wyckoff sites, but these sites are only 41.7% occupied. We use this structure as the prototype.

Base-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$

Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_1 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($4c$)	Fe
$\mathbf{B}_2$	=	$y_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_1 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	($4c$)	Fe
$\mathbf{B}_3$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_2 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($4c$)	O I
$\mathbf{B}_4$	=	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_2 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	($4c$)	O I
$\mathbf{B}_5$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_3 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($4c$)	O II
$\mathbf{B}_6$	=	$y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_3 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	($4c$)	O II

$\mathbf{B}_7$	=	$-y_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(8f)$	H
$\mathbf{B}_8$	=	$y_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$-y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right) c\,\hat{\mathbf{z}}$	$(8f)$	H
$\mathbf{B}_9$	=	$-y_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right) c\,\hat{\mathbf{z}}$	$(8f)$	H
$\mathbf{B}_{10}$	=	$y_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-y_4 b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	$(8f)$	H

References:

- A. Nørlund Christensen, M. S. Lehmann, and P. Convert, *Deuteration of Crystalline Hydroxides, Hydrogen Bonds of γ-AlOO(H,D) and γ-FeOO(H,D)*, Acta Chem. Scand. **36a**, 303–308 (1982), doi:10.3891/acta.chem.scand.36a-0303.
- C. Gottfried and F. Schossberger, eds., *Strukturbericht Band III 1933-1935* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Found in:

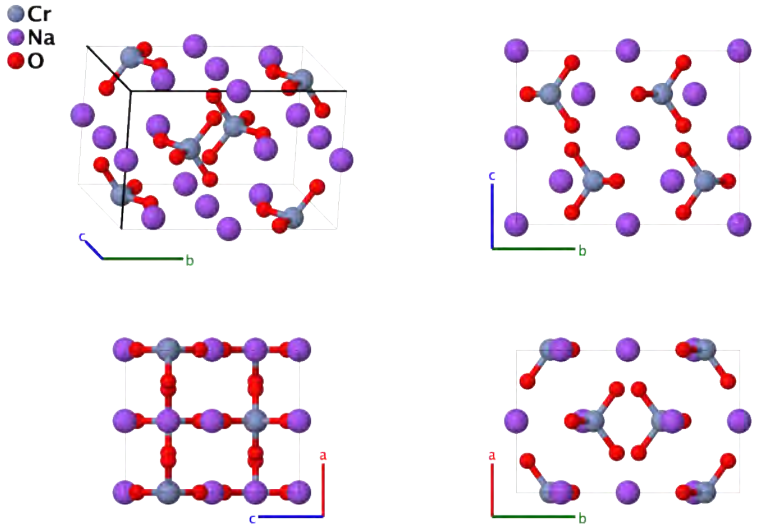
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1655
- POSCAR: pp. 1655

Na_2CrO_4 ($H1_8$) Structure: AB2C4_oC28_63_c_bc_fg

http://aflow.org/prototype-encyclopedia/AB2C4_oC28_63_c_bc_fg



Prototype	:	$CrNa_2O_4$
AFLOW prototype label	:	AB2C4_oC28_63_c_bc_fg
***Strukturbericht* designation**	:	$H1_8$
Pearson symbol	:	oC28
Space group number	:	63
Space group symbol	:	*Cmcm*
AFLOW prototype command	:	`aflow --proto=AB2C4_oC28_63_c_bc_fg --params=`$a, b/a, c/a, y_2, y_3, y_4, z_4, x_5, y_5$

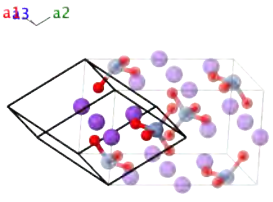
Other compounds with this structure

- Li_2SO_4, $LiFeP_4$, Na_2FeO_4, Na_2SO_4 (III), $NaCaVO_4$, $NaMnPO_4$, $NaVCdO_4$, and Tl_2SeO_4

- This structure was originally determined by (Miller, 1936), who placed it in space group *Pnna* #52, and (Gottfried, 1938) uses this data for $H1_8$. Subsequently (Niggli, 1954) rather acerbically pointed out that Miller's coordinates were consistent with the more compact *Cmcm* #63 space group. This does not change the positions of the atoms in the conventional cell, so we use the compact structure as our prototype for *Strukturbericht* designation $H1_8$.
- This structure is stable up to 413 °C (Amirathanlingam, 1982).

Base-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	($4b$)	Na I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4b$)	Na I
$\mathbf{B}_3$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_2 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($4c$)	Cr
$\mathbf{B}_4$	=	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_2 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	($4c$)	Cr
$\mathbf{B}_5$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_3 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($4c$)	Na II
$\mathbf{B}_6$	=	$y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_3 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	($4c$)	Na II
$\mathbf{B}_7$	=	$-y_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	($8f$)	O I
$\mathbf{B}_8$	=	$y_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$-y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	($8f$)	O I
$\mathbf{B}_9$	=	$-y_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	($8f$)	O I
$\mathbf{B}_{10}$	=	$y_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-y_4 b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	($8f$)	O I
$\mathbf{B}_{11}$	=	$(x_5 - y_5)\,\mathbf{a}_1 + (x_5 + y_5)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($8g$)	O II
$\mathbf{B}_{12}$	=	$(-x_5 + y_5)\,\mathbf{a}_1 + (-x_5 - y_5)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	($8g$)	O II
$\mathbf{B}_{13}$	=	$(-x_5 - y_5)\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($8g$)	O II
$\mathbf{B}_{14}$	=	$(x_5 + y_5)\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	($8g$)	O II

References:

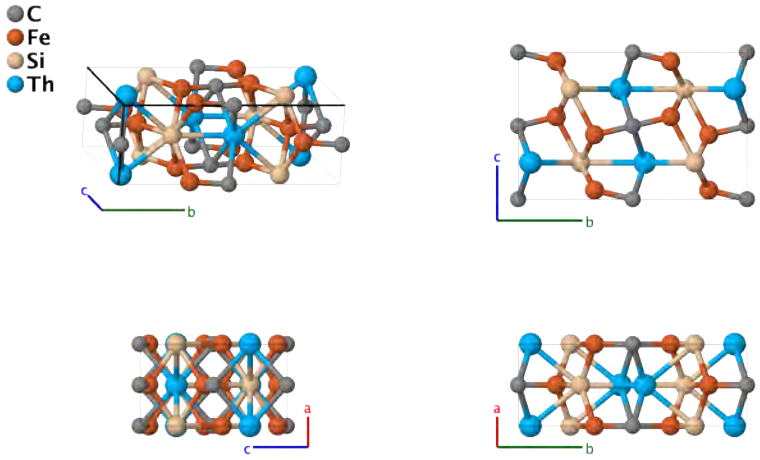
- A. Niggli, *Die Raumgruppe von Na_2CrO_4*, Acta Cryst. **7**, 776 (1954), doi:10.1107/S0365110X54002368.
- J. J. Miller, *The Crystal Structure of Anhydrous Sodium Chromate, Na_2CrO_4*, Zeitschrift für Kristallographie - Crystalline Materials **94**, 131–136 (1936), doi:10.1524/zkri.1936.94.1.131.
- C. Gottfried, ed., *Strukturbericht Band IV 1936* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1938).
- V. Amirathanlingam and K. S. Venkateswarlu, *The Thermal Expansion and Crystallographic Phase Transformation of Na_2CrO_4*, Thermochim. Acta **58**, 107–109 (1982), doi:10.1016/0040-6031(82)87145-1.

Geometry files:

- CIF: pp. 1656
- POSCAR: pp. 1656

$ThFe_2SiC$ Structure: AB2CD_oC20_63_b_f_c_c

http://aflow.org/prototype-encyclopedia/AB2CD_oC20_63_b_f_c_c



Prototype	:	CFe_2SiTh
AFLOW prototype label	:	AB2CD_oC20_63_b_f_c_c
***Strukturbericht* designation**	:	None
Pearson symbol	:	oC20
Space group number	:	63
Space group symbol	:	*Cmcm*
AFLOW prototype command	:	`aflow --proto=AB2CD_oC20_63_b_f_c_c --params=`$a, b/a, c/a, y_2, y_3, y_4, z_4$

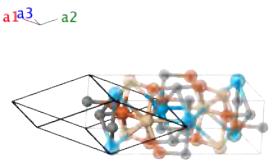
Other compounds with this structure

- AFe_2SiC, (A = Y, Sm, Gd, Tb, Ho, Er, Tm, Lu, U)

- This structure is a "filled" version of the Re_3B structure, with carbon atoms sitting in the ($4b$) Wyckoff positions. This is the quaternary version of the structure. The ternary version, where one of the ($4c$) Wyckoff positions has the same atom type as the ($8f$) site, is designated the V_3AsC structure.
- (Witte, 1994) does not designate a prototype for this structure, but as they only give full crystallographic data for the $ThFe_2SiC$ structure we will use this as the prototype.
- (Witte, 1994) lists the occupation of each site as Th (100.2%), Fe (97.7%), Si (96.0%), and C (97.0%), indicating that there are vacancies on the Fe, Si, and C sites, and that some of the Th atoms are either interstitial or occuping the other ($4c$) site or the ($8f$) site.

Base-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	$(4b)$	C
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(4b)$	C
$\mathbf{B}_3$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_2 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(4c)$	Si
$\mathbf{B}_4$	=	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_2 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(4c)$	Si
$\mathbf{B}_5$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_3 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(4c)$	Th
$\mathbf{B}_6$	=	$y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_3 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(4c)$	Th
$\mathbf{B}_7$	=	$-y_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(8f)$	Fe
$\mathbf{B}_8$	=	$y_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\,\mathbf{a}_3$	=	$-y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	$(8f)$	Fe
$\mathbf{B}_9$	=	$-y_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\,\mathbf{a}_3$	=	$y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	$(8f)$	Fe
$\mathbf{B}_{10}$	=	$y_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-y_4 b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	$(8f)$	Fe

References:

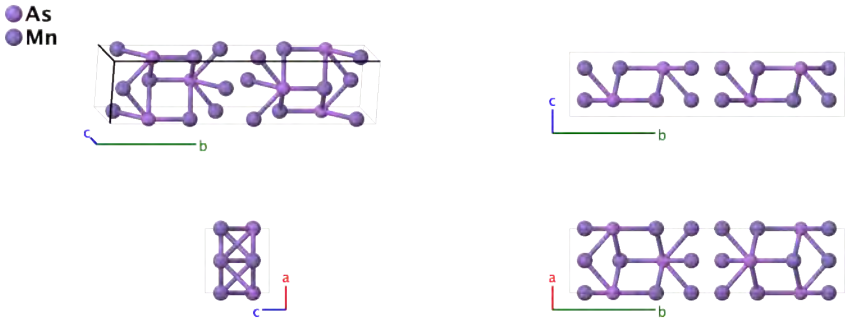
- A. M. Witte and W. Jeitschko, *Carbides with Filled Re_3B-Type Structure*, J. Solid State Chem. **112**, 232–236 (1994), doi:10.1006/jssc.1994.1297.

Geometry files:

- CIF: pp. 1656
- POSCAR: pp. 1656

Mn_3As ($D0_d$) Structure: AB3_oC16_63_c_3c

http://aflow.org/prototype-encyclopedia/AB3_oC16_63_c_3c



Prototype	:	$AsMn_3$
AFLOW prototype label	:	AB3_oC16_63_c_3c
***Strukturbericht* designation**	:	$D0_d$
Pearson symbol	:	oC16
Space group number	:	63
Space group symbol	:	*Cmcm*
AFLOW prototype command	:	`aflow --proto=AB3_oC16_63_c_3c --params=`$a, b/a, c/a, y_1, y_2, y_3, y_4$

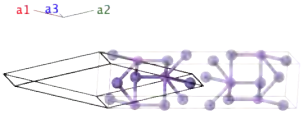
Other compounds with this structure

- Te_3Nd, $DyGe_3$

- (Nowotny, 1951) set the structure of Mn_3As in space group *Pmmn* #59. This was repeated by (Villars, 1991) and (Brandes, 1992). However, (Carrillo-Cabrera, 1983) showed that the structure actually reduces to space group *Cmcm* #63, and this was recognized by (Parthé, 1993). We follow the latter two works and assign the $D0_d$ structure to space group *Cmcm*.
- (Carrillo-Cabrera, 1983) placed the structure in setting *Bmmb* of space group #63, but we have shifted it to the standard *Cmcm* setting.

Base-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_1 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*c*)	As
$\mathbf{B}_2$	=	$y_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_1 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(4*c*)	As
$\mathbf{B}_3$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_2 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*c*)	Mn I
$\mathbf{B}_4$	=	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_2 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(4*c*)	Mn I

$\mathbf{B}_5$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_3 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*c*)	Mn II
$\mathbf{B}_6$	=	$y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_3 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(4*c*)	Mn II
$\mathbf{B}_7$	=	$-y_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_4 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*c*)	Mn III
$\mathbf{B}_8$	=	$y_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_4 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(4*c*)	Mn III

References:

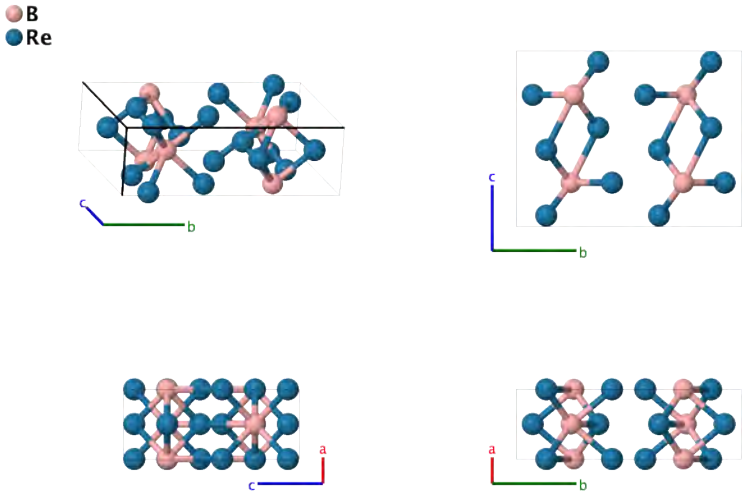
- W. Carrillo-Cabrera, *The Crystal Structure of $TiCu_2P$ and Its Relationship to the Structure of Mn_3As*, Acta Chem. Scand. **37a**, 93–98 (1983), doi:10.3891/acta.chem.scand.37a-0093.
- H. Nowotny, R. Funk, and J. Pesl, *Kristallchemische Untersuchungen in den Systemen Mn-As, V-Sb, Ti-Sb*, Monatsh. Chem. Verw. Teile Anderer Wiss. **82**, 513–525 (1951), doi:10.1007/BF00900849.
- E. Parthé, L. Gelato, B. Chabot, M. Penso, K. Cenzual, and R. Gladyshevskii, in *Standardized Data and Crystal Chemical Characterization of Inorganic Structure Types* (Springer-Verlag, Berlin, Heidelberg, 1993), *Gmelin Handbook of Inorganic and Organometallic Chemistry*, vol. 2, chap. Crystal Chemical Characterization of Inorganic Structure Types, 8 edn., doi:10.1007/978-3-662-02909-1_3.
- E. A. Brandes and G. B. Brook, eds., *Smithells Metals Reference Book* (Butterworth Heinemann, Oxford, Auckland, Boston, Johannesburg, Melbourne, New Delhi, 1992), seventh edn. Strukturbericht designations start on page 6-36 (163 in PDF), see table 6.3 on page 6-63 (190).
- P. Villars and L. D. Calvert, eds., *Pearson's Handbook of Crystallographic Data for Intermetallic Phases* (ASM International, Materials Park, Ohio, 1991), 2^{nd} edn.

Geometry files:

- CIF: pp. 1657
- POSCAR: pp. 1657

Re_3B Structure: AB3_oC16_63_c_cf

http://aflow.org/prototype-encyclopedia/AB3_oC16_63_c_cf

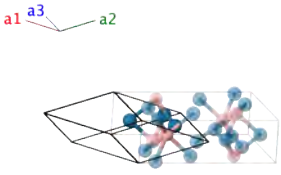


Prototype	:	BRe_3
AFLOW prototype label	:	AB3_oC16_63_c_cf
***Strukturbericht* designation**	:	None
Pearson symbol	:	oC16
Space group number	:	63
Space group symbol	:	*Cmcm*
AFLOW prototype command	:	`aflow --proto=AB3_oC16_63_c_cf --params=`$a, b/a, c/a, y_1, y_2, y_3, z_3$

- This is the parent structure of the ternary $MgCuAl_2$ ($E1_a$) structure. There are also "filled" versions of this structure, V_3AsC and $ThFe_2SiC$.

Base-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_1 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*c*)	B
$\mathbf{B}_2$	=	$y_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_1 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(4*c*)	B
$\mathbf{B}_3$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_2 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*c*)	Re I

$\mathbf{B}_4$	=	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_2 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(4c)$	Re I
$\mathbf{B}_5$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(8f)$	Re II
$\mathbf{B}_6$	=	$y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$-y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	$(8f)$	Re II
$\mathbf{B}_7$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	$(8f)$	Re II
$\mathbf{B}_8$	=	$y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-y_3 b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(8f)$	Re II

References:

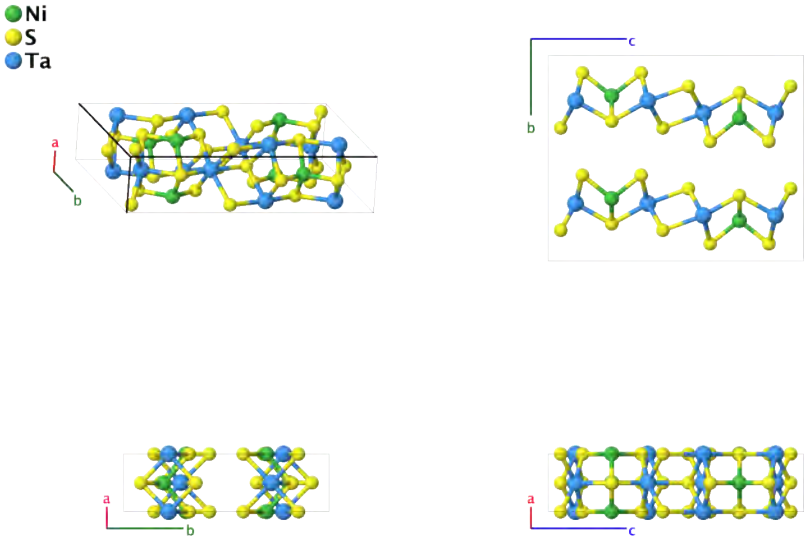
- B. Aronsson, M. Bäckman, and S. Rundqvist, *The Crystal Structure of Re_3B*, Acta Chem. Scand. **14**, 1001–1005 (1960), doi:10.3891/acta.chem.scand.14-1001.

Geometry files:

- CIF: pp. 1657
- POSCAR: pp. 1657

Ta_2NiS_5 Structure: AB5C2_oC32_63_c_c2f_f

http://aflow.org/prototype-encyclopedia/AB5C2_oC32_63_c_c2f_f



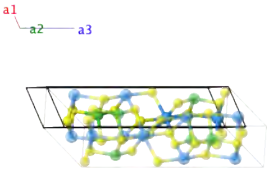
Prototype	:	NiS_5Ta_2
AFLOW prototype label	:	AB5C2_oC32_63_c_c2f_f
***Strukturbericht* designation**	:	None
Pearson symbol	:	oC32
Space group number	:	63
Space group symbol	:	*Cmcm*
AFLOW prototype command	:	`aflow --proto=AB5C2_oC32_63_c_c2f_f --params=`$a, b/a, c/a, y_1, y_2, y_3, z_3, y_4, z_4, y_5, z_5$

Other compounds with this structure

- Ta_2NiSe_5 (above 328 K)

Base-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_1 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(4c)$	Ni

$\mathbf{B}_2$	=	$y_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_1 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(4c)$	Ni
$\mathbf{B}_3$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_2 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(4c)$	S I
$\mathbf{B}_4$	=	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_2 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(4c)$	S I
$\mathbf{B}_5$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(8f)$	S II
$\mathbf{B}_6$	=	$y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\,\mathbf{a}_3$	=	$-y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	$(8f)$	S II
$\mathbf{B}_7$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\,\mathbf{a}_3$	=	$y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	$(8f)$	S II
$\mathbf{B}_8$	=	$y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-y_3 b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(8f)$	S II
$\mathbf{B}_9$	=	$-y_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(8f)$	S III
$\mathbf{B}_{10}$	=	$y_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\,\mathbf{a}_3$	=	$-y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	$(8f)$	S III
$\mathbf{B}_{11}$	=	$-y_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\,\mathbf{a}_3$	=	$y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	$(8f)$	S III
$\mathbf{B}_{12}$	=	$y_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-y_4 b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	$(8f)$	S III
$\mathbf{B}_{13}$	=	$-y_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$y_5 b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(8f)$	Ta
$\mathbf{B}_{14}$	=	$y_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\,\mathbf{a}_3$	=	$-y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	$(8f)$	Ta
$\mathbf{B}_{15}$	=	$-y_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\,\mathbf{a}_3$	=	$y_5 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	$(8f)$	Ta
$\mathbf{B}_{16}$	=	$y_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-y_5 b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	$(8f)$	Ta

References:

- S. A. Sunshine and J. A. Ibers, *Structure and physical properties of the new layered ternary chalcogenides tantalum nickel sulfide (Ta_2NiS_5) and tantalum nickel selenide (Ta_2NiSe_5)*, Inorg. Chem. **24**, 3611–3614 (1985), doi:10.1021/ic00216a027.

Found in:

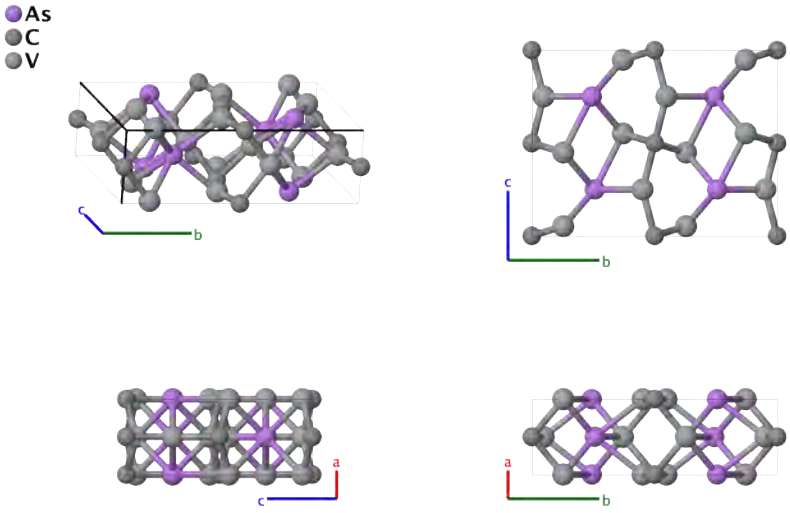
- F. J. Di Salvo, C. H. Chen, R. M. Fleming, J. V. Waszczak, R. G. Dunn, S. A. Sunshine, and J. A. Ibers, *Physical and structural properties of the new layered compounds Ta_2NiS_5 and Ta_2NiSe_5*, J. Less-Common Met. **116**, 51–61 (1986), doi:10.1016/0022-5088(86)90216-X.

Geometry files:

- CIF: pp. 1658
- POSCAR: pp. 1658

V_3AsC Structure: ABC3_oC20_63_c_b_cf

http://aflow.org/prototype-encyclopedia/ABC3_oC20_63_c_b_cf



Prototype	:	$AsCV_3$
AFLOW prototype label	:	ABC3_oC20_63_c_b_cf
***Strukturbericht* designation**	:	None
Pearson symbol	:	oC20
Space group number	:	63
Space group symbol	:	*Cmcm*
AFLOW prototype command	:	`aflow --proto=ABC3_oC20_63_c_b_cf --params=`$a, b/a, c/a, y_2, y_3, y_4, z_4$

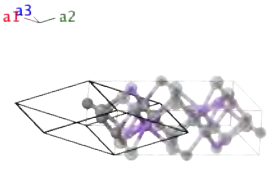
Other compounds with this structure

- V_3PC, V_3PN, V_3AsN, Cr_3PC, Cr_3AsC, Zr_3AlN, and $UScS_3$

- This structure is a "filled" version of the Re_3B structure, with carbon atoms sitting in the (4*b*) Wyckoff positions. This is the ternary version of the structure. The quaternary version, where all Wyckoff positions contain different species of atoms, is listed as the $ThFe_2SiC$ structure.

Base-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	(4*b*)	C
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*b*)	C
$\mathbf{B}_3$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_2 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*c*)	As
$\mathbf{B}_4$	=	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_2 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(4*c*)	As
$\mathbf{B}_5$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$y_3 b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*c*)	V I
$\mathbf{B}_6$	=	$y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-y_3 b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(4*c*)	V I
$\mathbf{B}_7$	=	$-y_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*f*)	V II
$\mathbf{B}_8$	=	$y_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$-y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(8*f*)	V II
$\mathbf{B}_9$	=	$-y_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$y_4 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(8*f*)	V II
$\mathbf{B}_{10}$	=	$y_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-y_4 b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*f*)	V II

References:

- H. Boller and H. Nowotny, *Zum Dreistoff: Vanadin-Arsen-Kohlenstoff*, Monatsh. Chem. Verw. Teile Anderer Wiss. **98**, 2127–2132 (1967), doi:10.1007/BF01167176.

Found in:

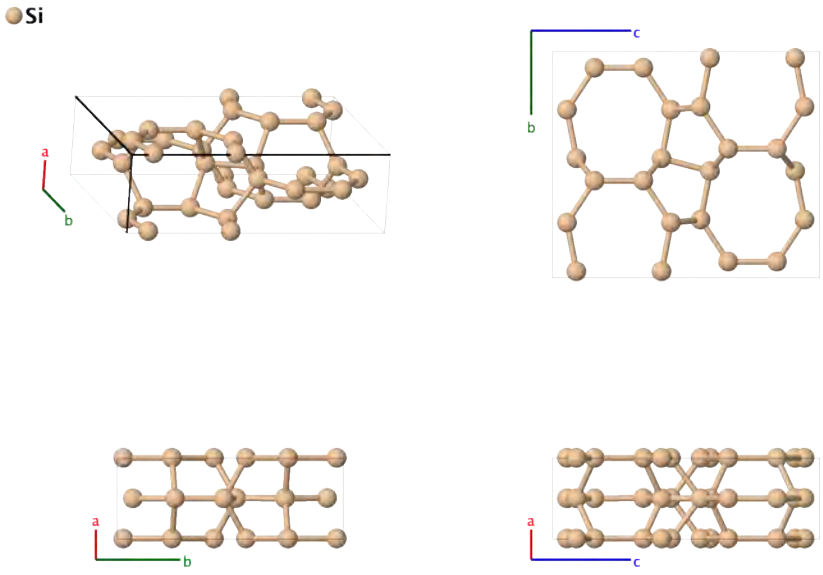
- A. M. Witte and W. Jeitschko, *Carbides with Filled Re_3B-Type Structure*, J. Solid State Chem. **112**, 232–236 (1994), doi:10.1006/jssc.1994.1297.

Geometry files:

- CIF: pp. 1658
- POSCAR: pp. 1659

Si_{24} Clathrate Structure: A_oC24_63_3f

http://aflow.org/prototype-encyclopedia/A_oC24_63_3f

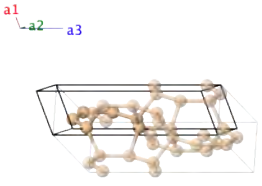


Prototype	:	Si_{24}
AFLOW prototype label	:	A_oC24_63_3f
***Strukturbericht* designation**	:	None
Pearson symbol	:	oC24
Space group number	:	63
Space group symbol	:	*Cmcm*
AFLOW prototype command	:	`aflow --proto=A_oC24_63_3f --params=`$a, b/a, c/a, y_1, z_1, y_2, z_2, y_3, z_3$

- Unlike the Si_{34} and Si_{46} clathrates, this is an experimentally determined structure, prepared by removing sodium atoms from an Na_4Si_{24} predecessor.
- There is no consistency in the naming of these clathrate structures. Si_{34} and Si_{46} were named by their authors based on the size of the primitive unit cell. Here, (Kim, 2015) has chosen to name the structure based on the size of the conventional cell. For now, at least, we will follow the authors' naming schemes for these structures.

Base-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$y_1 b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	$(8f)$	Si I
$\mathbf{B}_2$	=	$y_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$-y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right) c\,\hat{\mathbf{z}}$	$(8f)$	Si I
$\mathbf{B}_3$	=	$-y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	=	$y_1 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right) c\,\hat{\mathbf{z}}$	$(8f)$	Si I
$\mathbf{B}_4$	=	$y_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$-y_1 b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	$(8f)$	Si I
$\mathbf{B}_5$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(8f)$	Si II
$\mathbf{B}_6$	=	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$-y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	$(8f)$	Si II
$\mathbf{B}_7$	=	$-y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$y_2 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right) c\,\hat{\mathbf{z}}$	$(8f)$	Si II
$\mathbf{B}_8$	=	$y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$-y_2 b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	$(8f)$	Si II
$\mathbf{B}_9$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(8f)$	Si III
$\mathbf{B}_{10}$	=	$y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$-y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right) c\,\hat{\mathbf{z}}$	$(8f)$	Si III
$\mathbf{B}_{11}$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right) c\,\hat{\mathbf{z}}$	$(8f)$	Si III
$\mathbf{B}_{12}$	=	$y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-y_3 b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(8f)$	Si III

References:

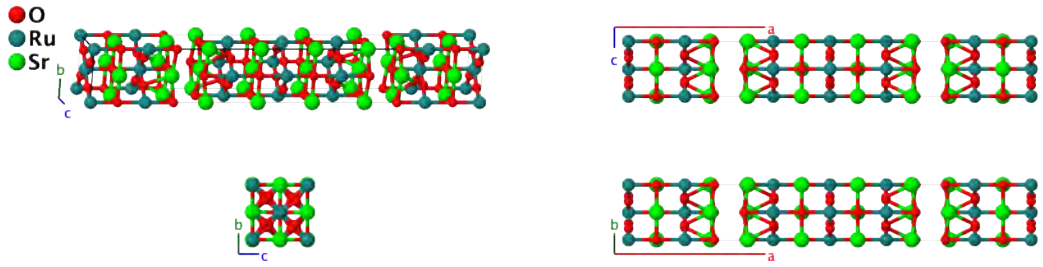
- D. Y. Kim, S. Stefanoski, O. O. Kurakevych, and T. A. Strobel, *Synthesis of an open-framework allotrope of silicon*, Nat. Mater. **14**, 169–173 (2015), doi:10.1038/NMAT4140.

Geometry files:

- CIF: pp. 1659
- POSCAR: pp. 1659

Base-centered orthorhombic $Sr_4Ru_3O_{10}$ Structure: A10B3C4_oC68_64_2dfg_ad_2d

http://aflow.org/prototype-encyclopedia/A10B3C4_oC68_64_2dfg_ad_2d

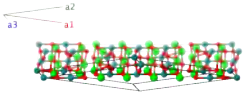


Prototype	:	$O_{10}Ru_3Sr_4$
AFLOW prototype label	:	A10B3C4_oC68_64_2dfg_ad_2d
***Strukturbericht* designation**	:	None
Pearson symbol	:	oC68
Space group number	:	64
Space group symbol	:	*Cmca*
AFLOW prototype command	:	`aflow --proto=A10B3C4_oC68_64_2dfg_ad_2d --params=`$a, b/a, c/a, x_2, x_3, x_4, x_5, x_6, y_7, z_7, x_8, y_8, z_8$

- (Crawford, 2002) placed $Sr_4Ru_3O_{10}$ in space group *Pbam* #55, oP68, but found that it was very close to the current structure. Indeed, we derived this structure from the original structure by allowing a small amount of uncertainty in the atomic positions. In both cases, the structure consists of triple-layer ruthenate separated by 2.37 Å. In the original structure, there are two of these layers in the primitive cell which are not equivalent. In the current structure there is only one triple-layer in the *primitive base-centered cell*, and so the layers in the *conventional* cell are equivalent.

Base-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(4*a*)	Ru I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*a*)	Ru I
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2$	=	$x_2 a\,\hat{\mathbf{x}}$	(8*d*)	O I
$\mathbf{B}_4$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(8*d*)	O I
$\mathbf{B}_5$	=	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2$	=	$-x_2 a\,\hat{\mathbf{x}}$	(8*d*)	O I
$\mathbf{B}_6$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(8*d*)	O I
$\mathbf{B}_7$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2$	=	$x_3 a\,\hat{\mathbf{x}}$	(8*d*)	O II
$\mathbf{B}_8$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(8*d*)	O II

$\mathbf{B}_{9}$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2$	=	$-x_3 a\,\hat{\mathbf{x}}$	$(8d)$	O II
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2}+x_3\right)\mathbf{a}_1 + \left(\frac{1}{2}+x_3\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(8d)$	O II
$\mathbf{B}_{11}$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2$	=	$x_4 a\,\hat{\mathbf{x}}$	$(8d)$	Ru II
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2}-x_4\right)\mathbf{a}_1 + \left(\frac{1}{2}-x_4\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(8d)$	Ru II
$\mathbf{B}_{13}$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2$	=	$-x_4 a\,\hat{\mathbf{x}}$	$(8d)$	Ru II
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2}+x_4\right)\mathbf{a}_1 + \left(\frac{1}{2}+x_4\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(8d)$	Ru II
$\mathbf{B}_{15}$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2$	=	$x_5 a\,\hat{\mathbf{x}}$	$(8d)$	Sr I
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2}-x_5\right)\mathbf{a}_1 + \left(\frac{1}{2}-x_5\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(8d)$	Sr I
$\mathbf{B}_{17}$	=	$-x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2$	=	$-x_5 a\,\hat{\mathbf{x}}$	$(8d)$	Sr I
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2}+x_5\right)\mathbf{a}_1 + \left(\frac{1}{2}+x_5\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(8d)$	Sr I
$\mathbf{B}_{19}$	=	$x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2$	=	$x_6 a\,\hat{\mathbf{x}}$	$(8d)$	Sr II
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2}-x_6\right)\mathbf{a}_1 + \left(\frac{1}{2}-x_6\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(8d)$	Sr II
$\mathbf{B}_{21}$	=	$-x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2$	=	$-x_6 a\,\hat{\mathbf{x}}$	$(8d)$	Sr II
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2}+x_6\right)\mathbf{a}_1 + \left(\frac{1}{2}+x_6\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(8d)$	Sr II
$\mathbf{B}_{23}$	=	$-y_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$y_7 b\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2}+y_7\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_7\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{25}$	=	$\left(\frac{1}{2}-y_7\right)\mathbf{a}_1 + \left(\frac{1}{2}+y_7\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_7\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + y_7 b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_7\right)c\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{26}$	=	$y_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$-y_7 b\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	$(8f)$	O III
$\mathbf{B}_{27}$	=	$(x_8-y_8)\,\mathbf{a}_1 + (x_8+y_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	$(16g)$	O IV
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2}-x_8+y_8\right)\mathbf{a}_1 + \left(\frac{1}{2}-x_8-y_8\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_8\right)a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_8\right)c\,\hat{\mathbf{z}}$	$(16g)$	O IV
$\mathbf{B}_{29}$	=	$\left(\frac{1}{2}-x_8-y_8\right)\mathbf{a}_1 + \left(\frac{1}{2}-x_8+y_8\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_8\right)a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_8\right)c\,\hat{\mathbf{z}}$	$(16g)$	O IV
$\mathbf{B}_{30}$	=	$(x_8+y_8)\,\mathbf{a}_1 + (x_8-y_8)\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	$(16g)$	O IV
$\mathbf{B}_{31}$	=	$(-x_8+y_8)\,\mathbf{a}_1 + (-x_8-y_8)\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	$(16g)$	O IV
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2}+x_8-y_8\right)\mathbf{a}_1 + \left(\frac{1}{2}+x_8+y_8\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_8\right)a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_8\right)c\,\hat{\mathbf{z}}$	$(16g)$	O IV
$\mathbf{B}_{33}$	=	$\left(\frac{1}{2}+x_8+y_8\right)\mathbf{a}_1 + \left(\frac{1}{2}+x_8-y_8\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_8\right)a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_8\right)c\,\hat{\mathbf{z}}$	$(16g)$	O IV
$\mathbf{B}_{34}$	=	$(-x_8-y_8)\,\mathbf{a}_1 + (-x_8+y_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	$(16g)$	O IV

References:

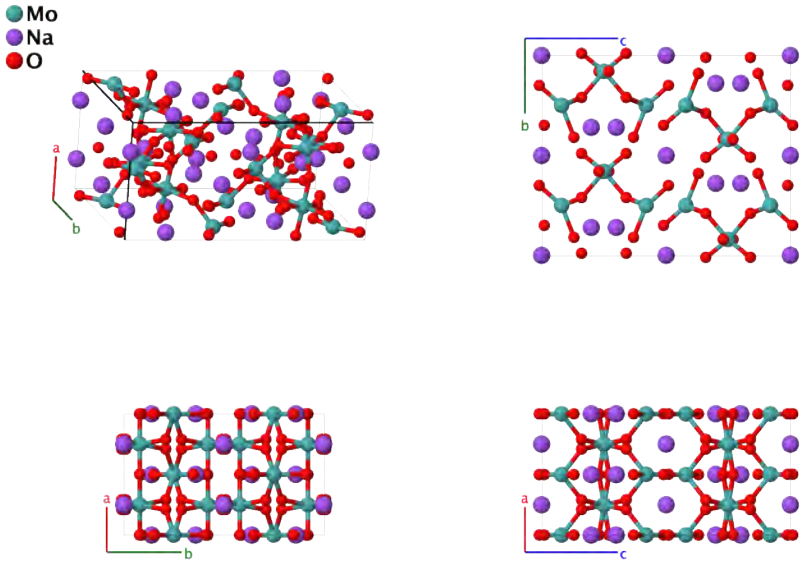
- M. K. Crawford, R. L. Harlow, W. Marshall, Z. Li, G. Cao, R. L. Lindstrom, Q. Huang, and J. W. Lynn, *Structure and magnetism of single crystal $Sr_4Ru_3O_{10}$: A ferromagnetic triple-layer ruthenate*, Phys. Rev. B **65**, 214412 (2002), doi:10.1103/PhysRevB.65.214412.

Geometry files:

- CIF: pp. 1659
- POSCAR: pp. 1660

$Na_2Mo_2O_7$ Structure: A2B2C7_oC88_64_ef_df_3f2g

http://aflow.org/prototype-encyclopedia/A2B2C7_oC88_64_ef_df_3f2g



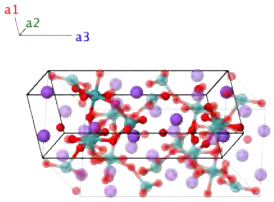
Prototype	:	$Mo_2Na_2O_7$
AFLOW prototype label	:	A2B2C7_oC88_64_ef_df_3f2g
***Strukturbericht* designation**	:	None
Pearson symbol	:	oC88
Space group number	:	64
Space group symbol	:	*Cmca*
AFLOW prototype command	:	`aflow --proto=A2B2C7_oC88_64_ef_df_3f2g --params=`$a, b/a, c/a, x_1, y_2, y_3, z_3, y_4, z_4, y_5, z_5, y_6, z_6, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9$

Other compounds with this structure

- $Na_2W_2O_7$

Base-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + x_1\,\mathbf{a}_2$	=	$x_1 a\,\hat{\mathbf{x}}$	($8d$)	Na I

$$
\begin{array}{lcccccc}
\mathbf{B}_2 & = & \left(\frac{1}{2}-x_1\right)\mathbf{a}_1+\left(\frac{1}{2}-x_1\right)\mathbf{a}_2+\frac{1}{2}\mathbf{a}_3 & = & \left(\frac{1}{2}-x_1\right)a\,\hat{\mathbf{x}}+\frac{1}{2}c\,\hat{\mathbf{z}} & (8d) & \text{Na I}\\
\mathbf{B}_3 & = & -x_1\,\mathbf{a}_1-x_1\,\mathbf{a}_2 & = & -x_1 a\,\hat{\mathbf{x}} & (8d) & \text{Na I}\\
\mathbf{B}_4 & = & \left(\frac{1}{2}+x_1\right)\mathbf{a}_1+\left(\frac{1}{2}+x_1\right)\mathbf{a}_2+\frac{1}{2}\mathbf{a}_3 & = & \left(\frac{1}{2}+x_1\right)a\,\hat{\mathbf{x}}+\frac{1}{2}c\,\hat{\mathbf{z}} & (8d) & \text{Na I}\\
\mathbf{B}_5 & = & \left(\frac{1}{4}-y_2\right)\mathbf{a}_1+\left(\frac{1}{4}+y_2\right)\mathbf{a}_2+\frac{1}{4}\mathbf{a}_3 & = & \frac{1}{4}a\,\hat{\mathbf{x}}+y_2 b\,\hat{\mathbf{y}}+\frac{1}{4}c\,\hat{\mathbf{z}} & (8e) & \text{Mo I}\\
\mathbf{B}_6 & = & \left(\frac{1}{4}+y_2\right)\mathbf{a}_1+\left(\frac{1}{4}-y_2\right)\mathbf{a}_2+\frac{3}{4}\mathbf{a}_3 & = & \frac{1}{4}a\,\hat{\mathbf{x}}-y_2 b\,\hat{\mathbf{y}}+\frac{3}{4}c\,\hat{\mathbf{z}} & (8e) & \text{Mo I}\\
\mathbf{B}_7 & = & \left(\frac{3}{4}+y_2\right)\mathbf{a}_1+\left(\frac{3}{4}-y_2\right)\mathbf{a}_2+\frac{3}{4}\mathbf{a}_3 & = & \frac{3}{4}a\,\hat{\mathbf{x}}-y_2 b\,\hat{\mathbf{y}}+\frac{3}{4}c\,\hat{\mathbf{z}} & (8e) & \text{Mo I}\\
\mathbf{B}_8 & = & \left(\frac{3}{4}-y_2\right)\mathbf{a}_1+\left(\frac{3}{4}+y_2\right)\mathbf{a}_2+\frac{1}{4}\mathbf{a}_3 & = & \frac{3}{4}a\,\hat{\mathbf{x}}+y_2 b\,\hat{\mathbf{y}}+\frac{1}{4}c\,\hat{\mathbf{z}} & (8e) & \text{Mo I}\\
\mathbf{B}_9 & = & -y_3\,\mathbf{a}_1+y_3\,\mathbf{a}_2+z_3\,\mathbf{a}_3 & = & y_3 b\,\hat{\mathbf{y}}+z_3 c\,\hat{\mathbf{z}} & (8f) & \text{Mo II}\\
\mathbf{B}_{10} & = & \left(\frac{1}{2}+y_3\right)\mathbf{a}_1+\left(\frac{1}{2}-y_3\right)\mathbf{a}_2+\left(\frac{1}{2}+z_3\right)\mathbf{a}_3 & = & \frac{1}{2}a\,\hat{\mathbf{x}}-y_3 b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_3\right)c\,\hat{\mathbf{z}} & (8f) & \text{Mo II}\\
\mathbf{B}_{11} & = & \left(\frac{1}{2}-y_3\right)\mathbf{a}_1+\left(\frac{1}{2}+y_3\right)\mathbf{a}_2+\left(\frac{1}{2}-z_3\right)\mathbf{a}_3 & = & \frac{1}{2}a\,\hat{\mathbf{x}}+y_3 b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_3\right)c\,\hat{\mathbf{z}} & (8f) & \text{Mo II}\\
\mathbf{B}_{12} & = & y_3\,\mathbf{a}_1-y_3\,\mathbf{a}_2-z_3\,\mathbf{a}_3 & = & -y_3 b\,\hat{\mathbf{y}}-z_3 c\,\hat{\mathbf{z}} & (8f) & \text{Mo II}\\
\mathbf{B}_{13} & = & -y_4\,\mathbf{a}_1+y_4\,\mathbf{a}_2+z_4\,\mathbf{a}_3 & = & y_4 b\,\hat{\mathbf{y}}+z_4 c\,\hat{\mathbf{z}} & (8f) & \text{Na II}\\
\mathbf{B}_{14} & = & \left(\frac{1}{2}+y_4\right)\mathbf{a}_1+\left(\frac{1}{2}-y_4\right)\mathbf{a}_2+\left(\frac{1}{2}+z_4\right)\mathbf{a}_3 & = & \frac{1}{2}a\,\hat{\mathbf{x}}-y_4 b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_4\right)c\,\hat{\mathbf{z}} & (8f) & \text{Na II}\\
\mathbf{B}_{15} & = & \left(\frac{1}{2}-y_4\right)\mathbf{a}_1+\left(\frac{1}{2}+y_4\right)\mathbf{a}_2+\left(\frac{1}{2}-z_4\right)\mathbf{a}_3 & = & \frac{1}{2}a\,\hat{\mathbf{x}}+y_4 b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_4\right)c\,\hat{\mathbf{z}} & (8f) & \text{Na II}\\
\mathbf{B}_{16} & = & y_4\,\mathbf{a}_1-y_4\,\mathbf{a}_2-z_4\,\mathbf{a}_3 & = & -y_4 b\,\hat{\mathbf{y}}-z_4 c\,\hat{\mathbf{z}} & (8f) & \text{Na II}\\
\mathbf{B}_{17} & = & -y_5\,\mathbf{a}_1+y_5\,\mathbf{a}_2+z_5\,\mathbf{a}_3 & = & y_5 b\,\hat{\mathbf{y}}+z_5 c\,\hat{\mathbf{z}} & (8f) & \text{O I}\\
\mathbf{B}_{18} & = & \left(\frac{1}{2}+y_5\right)\mathbf{a}_1+\left(\frac{1}{2}-y_5\right)\mathbf{a}_2+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3 & = & \frac{1}{2}a\,\hat{\mathbf{x}}-y_5 b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}} & (8f) & \text{O I}\\
\mathbf{B}_{19} & = & \left(\frac{1}{2}-y_5\right)\mathbf{a}_1+\left(\frac{1}{2}+y_5\right)\mathbf{a}_2+\left(\frac{1}{2}-z_5\right)\mathbf{a}_3 & = & \frac{1}{2}a\,\hat{\mathbf{x}}+y_5 b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_5\right)c\,\hat{\mathbf{z}} & (8f) & \text{O I}\\
\mathbf{B}_{20} & = & y_5\,\mathbf{a}_1-y_5\,\mathbf{a}_2-z_5\,\mathbf{a}_3 & = & -y_5 b\,\hat{\mathbf{y}}-z_5 c\,\hat{\mathbf{z}} & (8f) & \text{O I}\\
\mathbf{B}_{21} & = & -y_6\,\mathbf{a}_1+y_6\,\mathbf{a}_2+z_6\,\mathbf{a}_3 & = & y_6 b\,\hat{\mathbf{y}}+z_6 c\,\hat{\mathbf{z}} & (8f) & \text{O II}\\
\mathbf{B}_{22} & = & \left(\frac{1}{2}+y_6\right)\mathbf{a}_1+\left(\frac{1}{2}-y_6\right)\mathbf{a}_2+\left(\frac{1}{2}+z_6\right)\mathbf{a}_3 & = & \frac{1}{2}a\,\hat{\mathbf{x}}-y_6 b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}} & (8f) & \text{O II}\\
\mathbf{B}_{23} & = & \left(\frac{1}{2}-y_6\right)\mathbf{a}_1+\left(\frac{1}{2}+y_6\right)\mathbf{a}_2+\left(\frac{1}{2}-z_6\right)\mathbf{a}_3 & = & \frac{1}{2}a\,\hat{\mathbf{x}}+y_6 b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_6\right)c\,\hat{\mathbf{z}} & (8f) & \text{O II}\\
\mathbf{B}_{24} & = & y_6\,\mathbf{a}_1-y_6\,\mathbf{a}_2-z_6\,\mathbf{a}_3 & = & -y_6 b\,\hat{\mathbf{y}}-z_6 c\,\hat{\mathbf{z}} & (8f) & \text{O II}\\
\mathbf{B}_{25} & = & -y_7\,\mathbf{a}_1+y_7\,\mathbf{a}_2+z_7\,\mathbf{a}_3 & = & y_7 b\,\hat{\mathbf{y}}+z_7 c\,\hat{\mathbf{z}} & (8f) & \text{O III}\\
\mathbf{B}_{26} & = & \left(\frac{1}{2}+y_7\right)\mathbf{a}_1+\left(\frac{1}{2}-y_7\right)\mathbf{a}_2+\left(\frac{1}{2}+z_7\right)\mathbf{a}_3 & = & \frac{1}{2}a\,\hat{\mathbf{x}}-y_7 b\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}} & (8f) & \text{O III}\\
\mathbf{B}_{27} & = & \left(\frac{1}{2}-y_7\right)\mathbf{a}_1+\left(\frac{1}{2}+y_7\right)\mathbf{a}_2+\left(\frac{1}{2}-z_7\right)\mathbf{a}_3 & = & \frac{1}{2}a\,\hat{\mathbf{x}}+y_7 b\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_7\right)c\,\hat{\mathbf{z}} & (8f) & \text{O III}\\
\mathbf{B}_{28} & = & y_7\,\mathbf{a}_1-y_7\,\mathbf{a}_2-z_7\,\mathbf{a}_3 & = & -y_7 b\,\hat{\mathbf{y}}-z_7 c\,\hat{\mathbf{z}} & (8f) & \text{O III}\\
\mathbf{B}_{29} & = & (x_8-y_8)\,\mathbf{a}_1+(x_8+y_8)\,\mathbf{a}_2+z_8\,\mathbf{a}_3 & = & x_8 a\,\hat{\mathbf{x}}+y_8 b\,\hat{\mathbf{y}}+z_8 c\,\hat{\mathbf{z}} & (16g) & \text{O IV}
\end{array}
$$

$\mathbf{B}_{30}$	=	$\left(\frac{1}{2} - x_8 + y_8\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_8 - y_8\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_8\right)a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right)c\,\hat{\mathbf{z}}$	(16*g*)	O IV
$\mathbf{B}_{31}$	=	$\left(\frac{1}{2} - x_8 - y_8\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_8 + y_8\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_8\right)a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right)c\,\hat{\mathbf{z}}$	(16*g*)	O IV
$\mathbf{B}_{32}$	=	$(x_8 + y_8)\,\mathbf{a}_1 + (x_8 - y_8)\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(16*g*)	O IV
$\mathbf{B}_{33}$	=	$(-x_8 + y_8)\,\mathbf{a}_1 + (-x_8 - y_8)\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(16*g*)	O IV
$\mathbf{B}_{34}$	=	$\left(\frac{1}{2} + x_8 - y_8\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_8 + y_8\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_8\right)a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right)c\,\hat{\mathbf{z}}$	(16*g*)	O IV
$\mathbf{B}_{35}$	=	$\left(\frac{1}{2} + x_8 + y_8\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_8 - y_8\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_8\right)a\,\hat{\mathbf{x}} - y_8 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right)c\,\hat{\mathbf{z}}$	(16*g*)	O IV
$\mathbf{B}_{36}$	=	$(-x_8 - y_8)\,\mathbf{a}_1 + (-x_8 + y_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} + y_8 b\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(16*g*)	O IV
$\mathbf{B}_{37}$	=	$(x_9 - y_9)\,\mathbf{a}_1 + (x_9 + y_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(16*g*)	O V
$\mathbf{B}_{38}$	=	$\left(\frac{1}{2} - x_9 + y_9\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_9 - y_9\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_9\right)a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right)c\,\hat{\mathbf{z}}$	(16*g*)	O V
$\mathbf{B}_{39}$	=	$\left(\frac{1}{2} - x_9 - y_9\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_9 + y_9\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_9\right)a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_9\right)c\,\hat{\mathbf{z}}$	(16*g*)	O V
$\mathbf{B}_{40}$	=	$(x_9 + y_9)\,\mathbf{a}_1 + (x_9 - y_9)\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}}$	(16*g*)	O V
$\mathbf{B}_{41}$	=	$(-x_9 + y_9)\,\mathbf{a}_1 + (-x_9 - y_9)\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}}$	(16*g*)	O V
$\mathbf{B}_{42}$	=	$\left(\frac{1}{2} + x_9 - y_9\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_9 + y_9\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_9\right)a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_9\right)c\,\hat{\mathbf{z}}$	(16*g*)	O V
$\mathbf{B}_{43}$	=	$\left(\frac{1}{2} + x_9 + y_9\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_9 - y_9\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_9\right)a\,\hat{\mathbf{x}} - y_9 b\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right)c\,\hat{\mathbf{z}}$	(16*g*)	O V
$\mathbf{B}_{44}$	=	$(-x_9 - y_9)\,\mathbf{a}_1 + (-x_9 + y_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{x}} + y_9 b\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(16*g*)	O V

References:

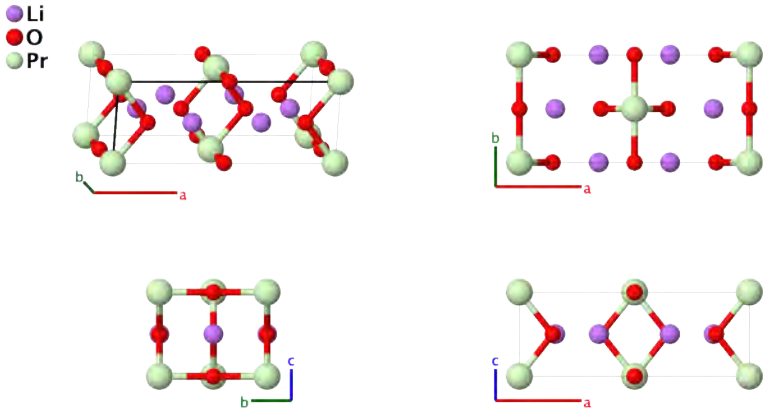
- I. Lindqvist, *Crystal Structure Studies on Anhydrous Sodium Molybdates and Tungstates*, Acta Chem. Scand. **4**, 1066–1074 (1950), doi:10.3891/acta.chem.scand.04-1066.

Geometry files:

- CIF: pp. 1660
- POSCAR: pp. 1660

Li_2PrO_3 Structure: A2B3C_oC12_65_h_bh_a

http://aflow.org/prototype-encyclopedia/A2B3C_oC12_65_h_bh_a



Prototype	:	Li_2O_3Pr
AFLOW prototype label	:	A2B3C_oC12_65_h_bh_a
***Strukturbericht* designation**	:	None
Pearson symbol	:	oC12
Space group number	:	65
Space group symbol	:	*Cmmm*
AFLOW prototype command	:	`aflow --proto=A2B3C_oC12_65_h_bh_a --params=`$a, b/a, c/a, x_3, x_4$

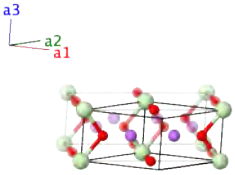
Other compounds with this structure

- Na_2PrO_3

- Data was taken at 20K.

Base-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(2*a*)	Pr
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	(2*b*)	O I
$\mathbf{B}_3$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*h*)	Li
$\mathbf{B}_4$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*h*)	Li

$\mathbf{B}_5$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + \frac{1}{2} c\,\hat{\mathbf{z}}$	(4*h*)	O II
$\mathbf{B}_6$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + \frac{1}{2} c\,\hat{\mathbf{z}}$	(4*h*)	O II

References:

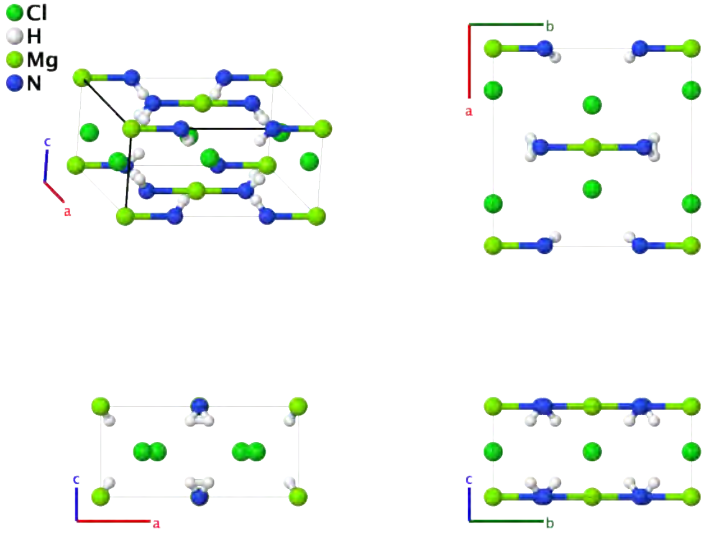
- Y. Hinatsu and Y. Doi, *Crystal structures and magnetic properties of alkali-metal lanthanide oxides A_2LnO_3 (A = Li, Na; Ln = Ce, Pr, Tb)*, J. Alloys Compd. **418**, 155–160 (2006), doi:10.1016/j.jallcom.2005.08.100.

Geometry files:

- CIF: pp. 1660
- POSCAR: pp. 1661

$Mg(NH_3)_2Cl_2$ ($E1_3$) Structure: A2B8CD2_oC26_65_h_r_a_i

http://aflow.org/prototype-encyclopedia/A2B8CD2_oC26_65_h_r_a_i



Prototype	:	$Cl_2H_6MgN_2$
AFLOW prototype label	:	A2B8CD2_oC26_65_h_r_a_i
***Strukturbericht* designation**	:	$E1_3$
Pearson symbol	:	oC26
Space group number	:	65
Space group symbol	:	*Cmmm*
AFLOW prototype command	:	`aflow --proto=A2B8CD2_oC26_65_h_r_a_i --params=`$a, b/a, c/a, x_2, y_3, x_4, y_4, z_4$

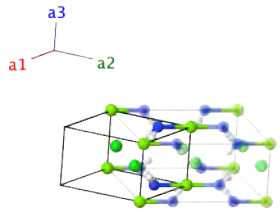
Other compounds with this structure

- $Cd(NH_3)_2Cl_2$, $Mg(NH_3)_2Cl_2$, $Hg(NH_3)_2Cl_2$, $Ni(NH_3)_2Br_2$, $Ni(NH_3)_2Cl_2$, and $Ni(NH_3)_2I_2$

- (Gottfried, 1938) gave the $E1_3$ designation to $Cd(NH_3)_2Cl_2$, and gave coordinates in the *Cmm*2 #35 space group. However, the cited reference, (MacGillavry, 1936) noted that their coordinates allowed several different space groups, with *Cmmm* #65 having the highest symmetry. We therefore follow most authors and use the *Cmmm* representation.
- (MacGillavry, 1936) could not determine the positions of the hydrogen atoms, but (Leineweber, 1999) was able to do this using the isostructural compound $Mg(NH_3)_2Cl_2$. Accordingly, we use $Mg(NH_3)_2Cl_2$ for the prototype of $E1_3$.
- Twelve hydrogen atoms are statistically distributed among the (16*r*) positions, so each site has a 75% probability of being occupied.

Base-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(2*a*)	Mg
$\mathbf{B}_2$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*h*)	Cl
$\mathbf{B}_3$	=	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*h*)	Cl
$\mathbf{B}_4$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2$	=	$y_3 b\,\hat{\mathbf{y}}$	(4*i*)	N
$\mathbf{B}_5$	=	$y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2$	=	$-y_3 b\,\hat{\mathbf{y}}$	(4*i*)	N
$\mathbf{B}_6$	=	$(x_4 - y_4)\,\mathbf{a}_1 + (x_4 + y_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(16*r*)	H
$\mathbf{B}_7$	=	$(-x_4 + y_4)\,\mathbf{a}_1 + (-x_4 - y_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(16*r*)	H
$\mathbf{B}_8$	=	$(-x_4 - y_4)\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(16*r*)	H
$\mathbf{B}_9$	=	$(x_4 + y_4)\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(16*r*)	H
$\mathbf{B}_{10}$	=	$(-x_4 + y_4)\,\mathbf{a}_1 + (-x_4 - y_4)\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(16*r*)	H
$\mathbf{B}_{11}$	=	$(x_4 - y_4)\,\mathbf{a}_1 + (x_4 + y_4)\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(16*r*)	H
$\mathbf{B}_{12}$	=	$(x_4 + y_4)\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(16*r*)	H
$\mathbf{B}_{13}$	=	$(-x_4 - y_4)\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(16*r*)	H

References:

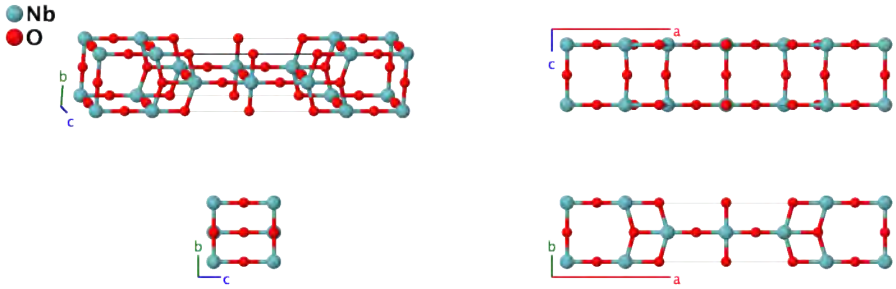
- A. Leineweber, M. W. Friedriszik, and H. Jacobs, *Preparation and Crystal Structures of $Mg(NH_3)_2Cl_2$, $Mg(NH_3)_2Br_2$, and $Mg(NH_3)_2I_2$*, J. Solid State Chem. **147**, 229–234 (1999), doi:10.1006/jssc.1999.8238.
- C. H. MacGillavry and J. M. Bijvoet, *Die Kristallstruktur von $Zn(NH_3)_2Cl_2$ und $Zn(NH_3)_2Br_2$*, Zeitschrift für Kristallographie - Crystalline Materials **94**, 249–255 (1936), doi:10.1524/zkri.1936.94.1.249.
- C. Gottfried, ed., *Strukturbericht Band IV 1936* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1938).

Geometry files:

- CIF: pp. 1661
- POSCAR: pp. 1661

Nb_3O_7F Structure: A3B8_oC22_65_ag_bd2gh

http://aflow.org/prototype-encyclopedia/A3B8_oC22_65_ag_bd2gh



Prototype	:	FNb_3O_7
AFLOW prototype label	:	A3B8_oC22_65_ag_bd2gh
***Strukturbericht* designation**	:	None
Pearson symbol	:	oC22
Space group number	:	65
Space group symbol	:	*Cmmm*
AFLOW prototype command	:	`aflow --proto=A3B8_oC22_65_ag_bd2gh --params=`$a, b/a, c/a, x_4, x_5, x_6, x_7$

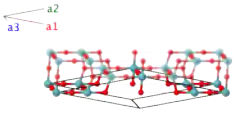
Other compounds with this structure

- $Nb_3O_7(OH)$

- (Andersson, 1964) was not able to distinguish between the oxygen and fluorine atoms, so it is assumed that the fluorine is distributed randomly on the oxygen sites.

Base-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(2*a*)	Nb I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	(2*b*)	O I
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	(2*d*)	O II
$\mathbf{B}_4$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2$	=	$x_4 a\,\hat{\mathbf{x}}$	(4*g*)	Nb II
$\mathbf{B}_5$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2$	=	$-x_4 a\,\hat{\mathbf{x}}$	(4*g*)	Nb II
$\mathbf{B}_6$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2$	=	$x_5 a\,\hat{\mathbf{x}}$	(4*g*)	O III
$\mathbf{B}_7$	=	$-x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2$	=	$-x_5 a\,\hat{\mathbf{x}}$	(4*g*)	O III
$\mathbf{B}_8$	=	$x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2$	=	$x_6 a\,\hat{\mathbf{x}}$	(4*g*)	O IV

$\mathbf{B}_9$	=	$-x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2$	=	$-x_6 a\,\hat{\mathbf{x}}$	($4g$)	O IV
$\mathbf{B}_{10}$	=	$x_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + \frac{1}{2} c\,\hat{\mathbf{z}}$	($4h$)	O V
$\mathbf{B}_{11}$	=	$-x_7\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} + \frac{1}{2} c\,\hat{\mathbf{z}}$	($4h$)	O V

References:

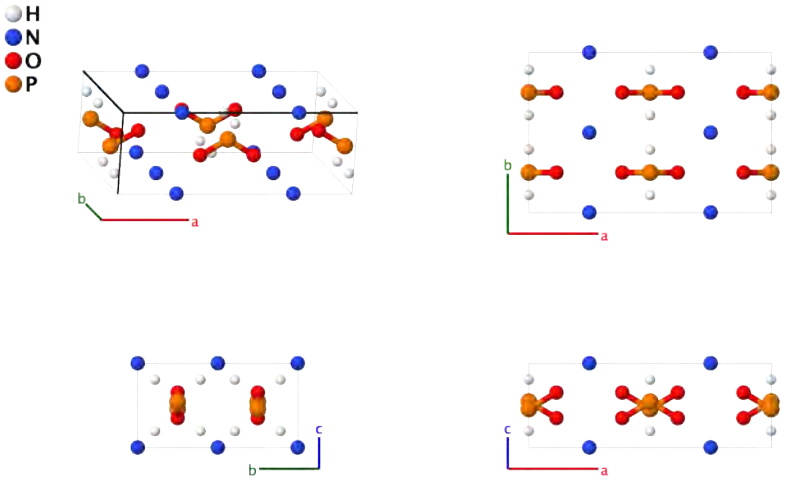
- S. Andersson, *The Crystal Structure of Nb_3O_7F*, Acta Chem. Scand. **18**, 2339–2344 (1964), doi:10.3891/acta.chem.scand.18-2339.

Geometry files:

- CIF: pp. 1661
- POSCAR: pp. 1662

$NH_4H_2PO_2$ ($F5_7$) Structure: A2BC2D_oC24_67_m_a_n_g

http://aflow.org/prototype-encyclopedia/A2BC2D_oC24_67_m_a_n_g

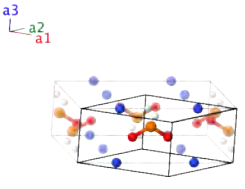


Prototype	:	$H_2(NH_4)O_2P$
AFLOW prototype label	:	A2BC2D_oC24_67_m_a_n_g
***Strukturbericht* designation**	:	$F5_7$
Pearson symbol	:	oC24
Space group number	:	67
Space group symbol	:	*Cmma*
AFLOW prototype command	:	`aflow --proto=A2BC2D_oC24_67_m_a_n_g --params=`$a, b/a, c/a, z_2, y_3, z_3, x_4, z_4$

- (Zachriasen, 1934) state that the H atoms in the ammonium ion must be along the lines between the nitrogen and oxygen atoms, but give no further information.
- The positions of the hydrogen atoms in the NH_4 ions were not determined, so we only provide the position of the nitrogen atoms (labeled as NH_4).
- The data for this structure was presented in the *Acmm* setting of space group #67. We transformed this to the standard *Cmma* setting using FINDSYM.

Base-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$\mathbf{B}_1$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2$	=	$\frac{1}{4}a\,\hat{\mathbf{x}}$	($4a$)	NH_4
$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2$	=	$\frac{3}{4}a\,\hat{\mathbf{x}}$	($4a$)	NH_4
$\mathbf{B}_3$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{4}b\,\hat{\mathbf{y}} + z_2c\,\hat{\mathbf{z}}$	($4g$)	P
$\mathbf{B}_4$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} - z_2c\,\hat{\mathbf{z}}$	($4g$)	P
$\mathbf{B}_5$	=	$-y_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$y_3b\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	($8m$)	H
$\mathbf{B}_6$	=	$\left(\frac{1}{2} + y_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - y_3b\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	($8m$)	H
$\mathbf{B}_7$	=	$\left(\frac{1}{2} - y_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + y_3b\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	($8m$)	H
$\mathbf{B}_8$	=	$y_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-y_3b\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	($8m$)	H
$\mathbf{B}_9$	=	$\left(\frac{3}{4} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{4} + x_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} - \frac{1}{4}b\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	($8n$)	O
$\mathbf{B}_{10}$	=	$\left(\frac{3}{4} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{4} - x_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} - \frac{1}{4}b\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	($8n$)	O
$\mathbf{B}_{11}$	=	$\left(\frac{1}{4} - x_4\right)\mathbf{a}_1 + \left(\frac{3}{4} - x_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	($8n$)	O
$\mathbf{B}_{12}$	=	$\left(\frac{1}{4} + x_4\right)\mathbf{a}_1 + \left(\frac{3}{4} + x_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	($8n$)	O

References:

- W. H. Zachariasen and R. C. L. Mooney, *The Structure of the Hypophosphite Group as Determined from the Crystal Lattice of Ammonium Hypophosphite*, J. Chem. Phys. **2**, 34–37 (1934), doi:10.1063/1.1749354.

Found in:

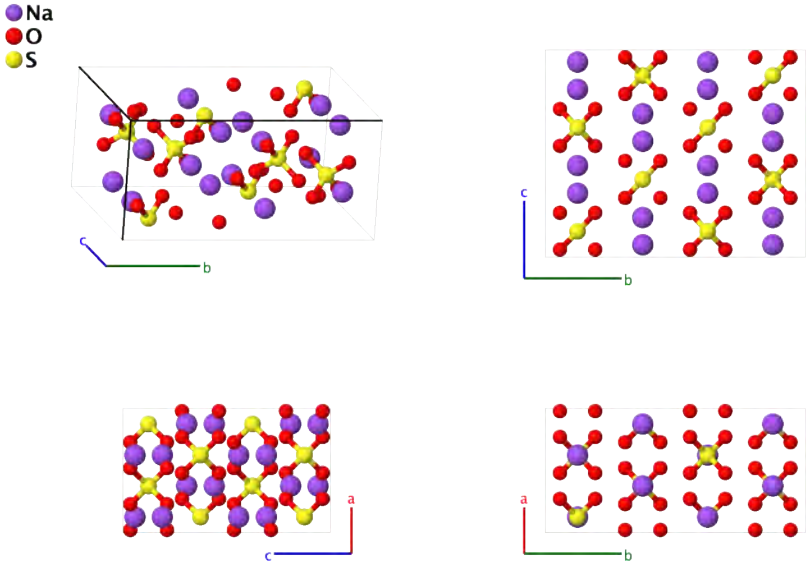
- C. Gottfried and F. Schossberger, eds., *Strukturbericht Band III 1933-1935* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1662
- POSCAR: pp. 1662

Thenardite [Na_2SO_4 (V), $H1_7$] Structure: A2B4C_oF56_70_g_h_a

http://aflow.org/prototype-encyclopedia/A2B4C_oF56_70_g_h_a



Prototype	:	Na_2O_4S
AFLOW prototype label	:	A2B4C_oF56_70_g_h_a
***Strukturbericht* designation**	:	$H1_7$
Pearson symbol	:	oF56
Space group number	:	70
Space group symbol	:	*Fddd*
AFLOW prototype command	:	`aflow --proto=A2B4C_oF56_70_g_h_a --params=`$a, b/a, c/a, z_2, x_3, y_3, z_3$

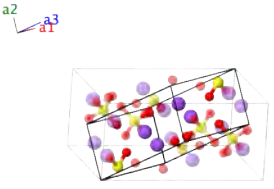
Other compounds with this structure

- Ag_2SO_4 and Cr_2SO_4

- Na_2SO_4 has eight known anhydrous phases. The thenardite phase is "reported to be stable between 32 °C and about 180 °C" (Nord, 1973), but the data reported here was taken on synthetic thenardite at 25 °C.

Face-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,b\,\hat{\mathbf{y}} + \frac{1}{2}\,c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{1}{2}\,c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{1}{2}\,b\,\hat{\mathbf{y}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{8}\,\mathbf{a}_1 + \frac{1}{8}\,\mathbf{a}_2 + \frac{1}{8}\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \frac{1}{8}b\,\hat{\mathbf{y}} + \frac{1}{8}c\,\hat{\mathbf{z}}$	(8*a*)	S
$\mathbf{B}_2$	=	$\frac{7}{8}\,\mathbf{a}_1 + \frac{7}{8}\,\mathbf{a}_2 + \frac{7}{8}\,\mathbf{a}_3$	=	$\frac{7}{8}a\,\hat{\mathbf{x}} + \frac{7}{8}b\,\hat{\mathbf{y}} + \frac{7}{8}c\,\hat{\mathbf{z}}$	(8*a*)	S
$\mathbf{B}_3$	=	$z_2\,\mathbf{a}_1 + z_2\,\mathbf{a}_2 + \left(\frac{1}{4} - z_2\right)\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \frac{1}{8}b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(16*g*)	Na
$\mathbf{B}_4$	=	$\left(\frac{1}{4} - z_2\right)\mathbf{a}_1 + \left(\frac{1}{4} - z_2\right)\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \frac{1}{8}b\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_2\right)c\,\hat{\mathbf{z}}$	(16*g*)	Na
$\mathbf{B}_5$	=	$-z_2\,\mathbf{a}_1 - z_2\,\mathbf{a}_2 + \left(\frac{3}{4} + z_2\right)\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} + \frac{3}{8}b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(16*g*)	Na
$\mathbf{B}_6$	=	$\left(\frac{3}{4} + z_2\right)\mathbf{a}_1 + \left(\frac{3}{4} + z_2\right)\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} + \frac{3}{8}b\,\hat{\mathbf{y}} + \left(\frac{3}{4} + z_2\right)c\,\hat{\mathbf{z}}$	(16*g*)	Na
$\mathbf{B}_7$	=	$(-x_3 + y_3 + z_3)\,\mathbf{a}_1 + (x_3 - y_3 + z_3)\,\mathbf{a}_2 + (x_3 + y_3 - z_3)\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(32*h*)	O
$\mathbf{B}_8$	=	$(x_3 - y_3 + z_3)\,\mathbf{a}_1 + (-x_3 + y_3 + z_3)\,\mathbf{a}_2 + \left(\frac{1}{2} - x_3 - y_3 - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - y_3\right)b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(32*h*)	O
$\mathbf{B}_9$	=	$(x_3 + y_3 - z_3)\,\mathbf{a}_1 + \left(\frac{1}{2} - x_3 - y_3 - z_3\right)\mathbf{a}_2 + (-x_3 + y_3 + z_3)\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_3\right)c\,\hat{\mathbf{z}}$	(32*h*)	O
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_3 - y_3 - z_3\right)\mathbf{a}_1 + (x_3 + y_3 - z_3)\,\mathbf{a}_2 + (x_3 - y_3 + z_3)\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_3\right)c\,\hat{\mathbf{z}}$	(32*h*)	O
$\mathbf{B}_{11}$	=	$(x_3 - y_3 - z_3)\,\mathbf{a}_1 + (-x_3 + y_3 - z_3)\,\mathbf{a}_2 + (-x_3 - y_3 + z_3)\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(32*h*)	O
$\mathbf{B}_{12}$	=	$(-x_3 + y_3 - z_3)\,\mathbf{a}_1 + (x_3 - y_3 - z_3)\,\mathbf{a}_2 + \left(\frac{1}{2} + x_3 + y_3 + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + y_3\right)b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(32*h*)	O
$\mathbf{B}_{13}$	=	$(-x_3 - y_3 + z_3)\,\mathbf{a}_1 + \left(\frac{1}{2} + x_3 + y_3 + z_3\right)\mathbf{a}_2 + (x_3 - y_3 - z_3)\,\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_3\right)c\,\hat{\mathbf{z}}$	(32*h*)	O
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} + x_3 + y_3 + z_3\right)\mathbf{a}_1 + (-x_3 - y_3 + z_3)\,\mathbf{a}_2 + (-x_3 + y_3 - z_3)\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + y_3\right)b\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_3\right)c\,\hat{\mathbf{z}}$	(32*h*)	O

References:

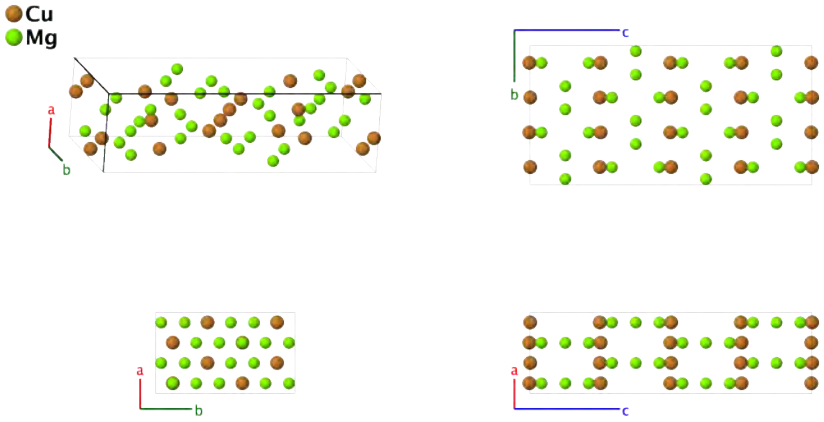
- A. G. Nord, *Refinement of the Crystal Structure of Thenardite, $Na_2SO_4(V)$*, Acta Chem. Scand. **27**, 814–822 (1973), doi:10.3891/acta.chem.scand.27-0814.

Geometry files:

- CIF: pp. 1662
- POSCAR: pp. 1663

Mg_2Cu (C_b) Structure: AB2_oF48_70_g_fg

http://aflow.org/prototype-encyclopedia/AB2_oF48_70_g_fg



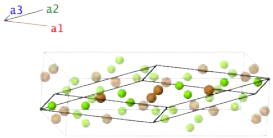
Prototype	:	$CuMg_2$
AFLOW prototype label	:	AB2_oF48_70_g_fg
***Strukturbericht* designation**	:	C_b
Pearson symbol	:	oF48
Space group number	:	70
Space group symbol	:	$Fddd$
AFLOW prototype command	:	`aflow --proto=AB2_oF48_70_g_fg --params=`$a, b/a, c/a, y_1, z_2, z_3$

Other compounds with this structure

- In_2Ir, $NbSn_2$, $TiBi_2$, FeGaGe, and SbSnTi

Face-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,b\,\hat{\mathbf{y}} + \frac{1}{2}\,c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{1}{2}\,c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{1}{2}\,b\,\hat{\mathbf{y}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$y_1\,\mathbf{a}_1 + \left(\frac{1}{4} - y_1\right)\mathbf{a}_2 + y_1\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + y_1 b\,\hat{\mathbf{y}} + \frac{1}{8}c\,\hat{\mathbf{z}}$	$(16f)$	Mg I
$\mathbf{B}_2$	=	$\left(\frac{1}{4} - y_1\right)\mathbf{a}_1 + y_1\,\mathbf{a}_2 + \left(\frac{1}{4} - y_1\right)\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - y_1\right)b\,\hat{\mathbf{y}} + \frac{1}{8}c\,\hat{\mathbf{z}}$	$(16f)$	Mg I
$\mathbf{B}_3$	=	$-y_1\,\mathbf{a}_1 + \left(\frac{3}{4} + y_1\right)\mathbf{a}_2 - y_1\,\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} - y_1 b\,\hat{\mathbf{y}} + \frac{3}{8}c\,\hat{\mathbf{z}}$	$(16f)$	Mg I
$\mathbf{B}_4$	=	$\left(\frac{3}{4} + y_1\right)\mathbf{a}_1 - y_1\,\mathbf{a}_2 + \left(\frac{3}{4} + y_1\right)\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} + \left(\frac{3}{4} + y_1\right)b\,\hat{\mathbf{y}} + \frac{3}{8}c\,\hat{\mathbf{z}}$	$(16f)$	Mg I
$\mathbf{B}_5$	=	$z_2\,\mathbf{a}_1 + z_2\,\mathbf{a}_2 + \left(\frac{1}{4} - z_2\right)\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \frac{1}{8}b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(16g)$	Cu
$\mathbf{B}_6$	=	$\left(\frac{1}{4} - z_2\right)\mathbf{a}_1 + \left(\frac{1}{4} - z_2\right)\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \frac{1}{8}b\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_2\right)c\,\hat{\mathbf{z}}$	$(16g)$	Cu

$\mathbf{B}_7$	=	$-z_2\,\mathbf{a}_1 - z_2\,\mathbf{a}_2 + \left(\frac{3}{4} + z_2\right)\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} + \frac{3}{8}b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(16*g*)	Cu
$\mathbf{B}_8$	=	$\left(\frac{3}{4} + z_2\right)\mathbf{a}_1 + \left(\frac{3}{4} + z_2\right)\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} + \frac{3}{8}b\,\hat{\mathbf{y}} + \left(\frac{3}{4} + z_2\right)c\,\hat{\mathbf{z}}$	(16*g*)	Cu
$\mathbf{B}_9$	=	$z_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_2 + \left(\frac{1}{4} - z_3\right)\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \frac{1}{8}b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(16*g*)	Mg II
$\mathbf{B}_{10}$	=	$\left(\frac{1}{4} - z_3\right)\mathbf{a}_1 + \left(\frac{1}{4} - z_3\right)\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \frac{1}{8}b\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_3\right)c\,\hat{\mathbf{z}}$	(16*g*)	Mg II
$\mathbf{B}_{11}$	=	$-z_3\,\mathbf{a}_1 - z_3\,\mathbf{a}_2 + \left(\frac{3}{4} + z_3\right)\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} + \frac{3}{8}b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(16*g*)	Mg II
$\mathbf{B}_{12}$	=	$\left(\frac{3}{4} + z_3\right)\mathbf{a}_1 + \left(\frac{3}{4} + z_3\right)\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} + \frac{3}{8}b\,\hat{\mathbf{y}} + \left(\frac{3}{4} + z_3\right)c\,\hat{\mathbf{z}}$	(16*g*)	Mg II

References:

- F. Gingl, P. Selvam, and K. Yvon, *Structure refinement of Mg_2Cu and a comparison of the Mg_2Cu, Mg_2Ni and Al_2Cu structure types*, Acta Crystallogr. Sect. B Struct. Sci. **49**, 201–203 (1993), doi:10.1107/S0108768192008723.

Found in:

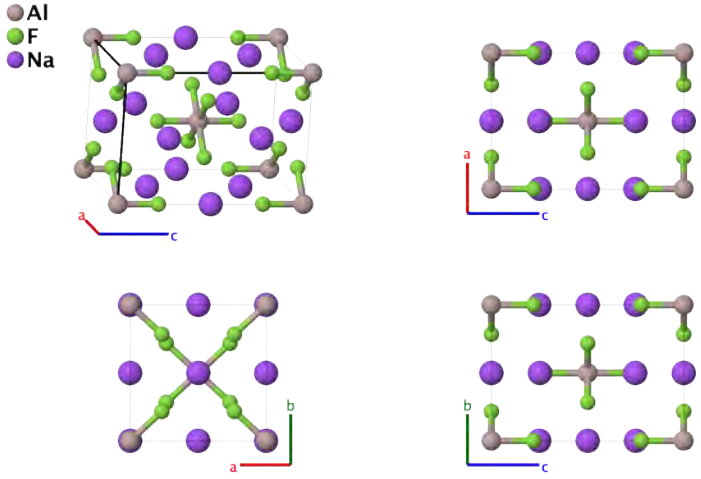
- V. Vreshch, *Crystal Structure of $CuMg_2$*, http://crystallography-online.com/structure/2100926 (2018). Crystallography online.com.

Geometry files:

- CIF: pp. 1663
- POSCAR: pp. 1663

High-Temperature Cryolite (Na_3AlF_6) Structure: AB6C3_oI20_71_a_in_cj

http://aflow.org/prototype-encyclopedia/AB6C3_oI20_71_a_in_cj

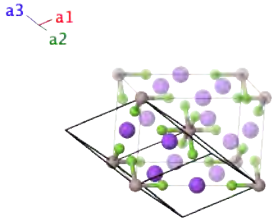


Prototype	:	AlF_6Na_3
AFLOW prototype label	:	AB6C3_oI20_71_a_in_cj
***Strukturbericht* designation**	:	None
Pearson symbol	:	oI20
Space group number	:	71
Space group symbol	:	*Immm*
AFLOW prototype command	:	`aflow --proto=AB6C3_oI20_71_a_in_cj --params=`$a, b/a, c/a, z_3, z_4, x_5, y_5$

- Cryolite undergoes a phase transition from the monoclinic $P2_1/c$ $J2_6$ phase to this orthorhombic phase at 890 K. We show structural data taken at 900 K.

Body-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = -\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}b\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}a\,\hat{\mathbf{x}} - \tfrac{1}{2}b\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}b\,\hat{\mathbf{y}} - \tfrac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($2a$)	Al
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	($2c$)	Na I
$\mathbf{B}_3$	=	$z_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_2$	=	$z_3 c\,\hat{\mathbf{z}}$	($4i$)	F I

$\mathbf{B}_4$	=	$-z_3\,\mathbf{a}_1 - z_3\,\mathbf{a}_2$	=	$-z_3 c\,\hat{\mathbf{z}}$	($4i$)	F I
$\mathbf{B}_5$	=	$z_4\,\mathbf{a}_1 + \left(\frac{1}{2} + z_4\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + z_4 c\,\hat{\mathbf{z}}$	($4j$)	Na II
$\mathbf{B}_6$	=	$-z_4\,\mathbf{a}_1 + \left(\frac{1}{2} - z_4\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - z_4 c\,\hat{\mathbf{z}}$	($4j$)	Na II
$\mathbf{B}_7$	=	$y_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + (x_5 + y_5)\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}}$	($8n$)	F II
$\mathbf{B}_8$	=	$-y_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + (-x_5 - y_5)\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}}$	($8n$)	F II
$\mathbf{B}_9$	=	$y_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + (-x_5 + y_5)\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + y_5 b\,\hat{\mathbf{y}}$	($8n$)	F II
$\mathbf{B}_{10}$	=	$-y_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + (x_5 - y_5)\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} - y_5 b\,\hat{\mathbf{y}}$	($8n$)	F II

References:

- H. Yang, S. Ghose, and D. M. Hatch, *Ferroelastic phase transition in cryolite, Na_3AlF_6, a mixed fluoride perovskite: High temperature single crystal X-ray diffraction study and symmetry analysis of the transition mechanism*, Phys. Chem. Miner. **19**, 528–544 (1993), doi:10.1007/BF00203053.

Found in:

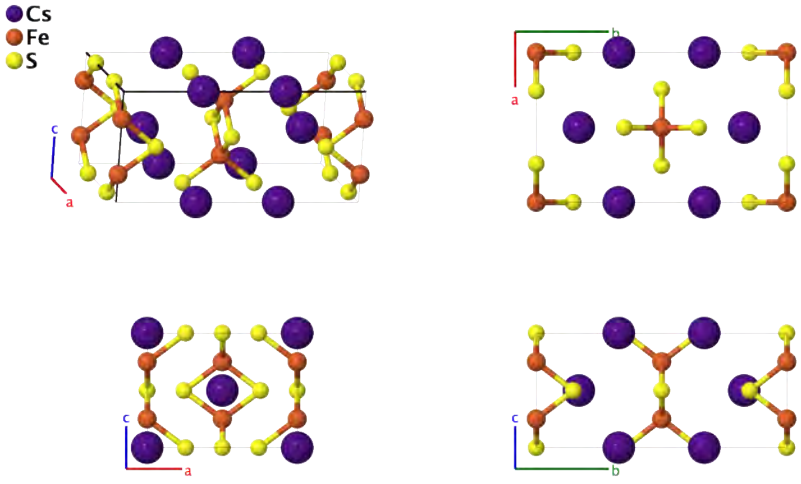
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1664
- POSCAR: pp. 1664

$CsFeS_2$ (100 K) Structure: ABC2_oI16_71_g_i_eh

http://aflow.org/prototype-encyclopedia/ABC2_oI16_71_g_i_eh



Prototype	:	$CsFeS_2$
AFLOW prototype label	:	ABC2_oI16_71_g_i_eh
***Strukturbericht* designation**	:	None
Pearson symbol	:	oI16
Space group number	:	71
Space group symbol	:	*Immm*
AFLOW prototype command	:	`aflow --proto=ABC2_oI16_71_g_i_eh --params=`$a, b/a, c/a, x_1, y_2, y_3, z_4$

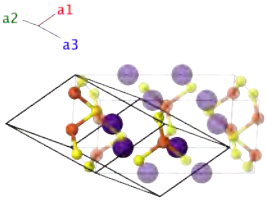
Other compounds with this structure

- $RbFeS_2$

- This structure is stable at 100 K and above. At 40 K the structure is triclinic, with as yet undetermined atomic positions (Ito, 1985).

Body-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = -\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} - \frac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_2 + x_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}}$	$(4e)$	S I
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_2 - x_1\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}}$	$(4e)$	S I

$\mathbf{B}_3$	=	$y_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_3$	=	$y_2 b\,\hat{\mathbf{y}}$	$(4g)$	Cs
$\mathbf{B}_4$	=	$-y_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_3$	=	$-y_2 b\,\hat{\mathbf{y}}$	$(4g)$	Cs
$\mathbf{B}_5$	=	$\left(\frac{1}{2} + y_3\right)\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + y_3\,\mathbf{a}_3$	=	$y_3 b\,\hat{\mathbf{y}} + \frac{1}{2} c\,\hat{\mathbf{z}}$	$(4h)$	S II
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - y_3\right)\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 - y_3\,\mathbf{a}_3$	=	$-y_3 b\,\hat{\mathbf{y}} + \frac{1}{2} c\,\hat{\mathbf{z}}$	$(4h)$	S II
$\mathbf{B}_7$	=	$z_4\,\mathbf{a}_1 + z_4\,\mathbf{a}_2$	=	$z_4 c\,\hat{\mathbf{z}}$	$(4i)$	Fe
$\mathbf{B}_8$	=	$-z_4\,\mathbf{a}_1 - z_4\,\mathbf{a}_2$	=	$-z_4 c\,\hat{\mathbf{z}}$	$(4i)$	Fe

References:

- Y. Ito, M. Nishi, C. F. Majkrzak, and L. Passell, *Low Temperature Powder Neutron Diffraction Studies of $CsFeS_2$*, J. Phys. Soc. Jpn. **54**, 348–357 (1985), doi:10.1143/JPSJ.54.348.

Found in:

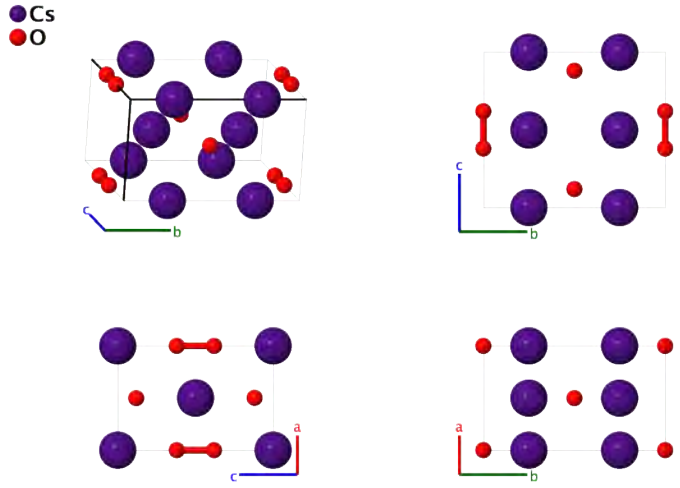
- P. Villars (Chief Editor), *CsFeS2 (100K) Crystal Structure*, http://materials.springer.com/isp/crystallographic/docs/sd_0382809 (2016). PAULING FILE in: Inorganic Solid Phases, SpringerMaterials (online database), Springer, Heidelberg (ed.) SpringerMaterials.

Geometry files:

- CIF: pp. 1664
- POSCAR: pp. 1665

CsO Structure: AB_oI8_71_g_i

http://aflow.org/prototype-encyclopedia/AB_oI8_71_g_i

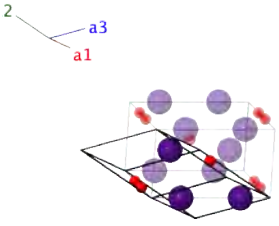


Prototype	:	CsO
AFLOW prototype label	:	AB_oI8_71_g_i
***Strukturbericht* designation**	:	None
Pearson symbol	:	oI8
Space group number	:	71
Space group symbol	:	*Immm*
AFLOW prototype command	:	`aflow --proto=AB_oI8_71_g_i --params=`$a, b/a, c/a, y_1, z_2$

- (Massalski, 1990) credits the discovery of this structure to (Rengade, 1909), but we have been unable to obtain a copy of this reference. (Downs, 2003) quotes (Wyckoff, 1963) giving an *Immm*-oI8 structure for CsO. Since this is the same space group and Pearson symbol as found in Massalski, we use Wyckoff's data.

Body-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = -\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} - \frac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$y_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_3$	=	$y_1 b\,\hat{\mathbf{y}}$	$(4g)$	Cs
$\mathbf{B}_2$	=	$-y_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_3$	=	$-y_1 b\,\hat{\mathbf{y}}$	$(4g)$	Cs
$\mathbf{B}_3$	=	$z_2\,\mathbf{a}_1 + z_2\,\mathbf{a}_2$	=	$z_2 c\,\hat{\mathbf{z}}$	$(4i)$	O

$\mathbf{B}_4$	=	$-z_2\,\mathbf{a}_1 - z_2\,\mathbf{a}_2$	=	$-z_2 c\,\hat{\mathbf{z}}$	$(4i)$	O

References:

- M. E. Rengade, *Sur les Sous-Oxydes de Caesium*, C. R. Acad. Sci. C **148**, 1199–1202 (1909).

Found in:

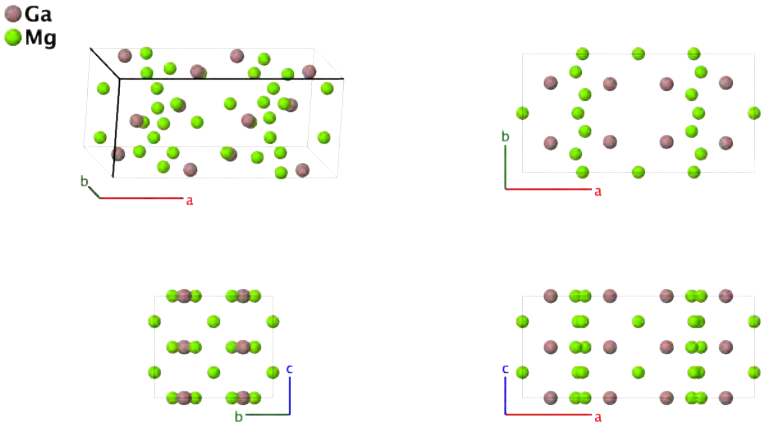
- T. B. Massalski, H. Okamoto, P. R. Subramanian, and L. Kacprzak, eds., *Binary Alloy Phase Diagrams*, vol. 2 (ASM International, Materials Park, Ohio, USA, 1990), 2^{nd} edn. Cd-Ce to Hf-Rb.
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).
- R. W. G. Wyckoff, *Crystal Structures*, vol. 1 (Interscience Publishers, New York, 1963), second edn.

Geometry files:

- CIF: pp. 1665
- POSCAR: pp. 1665

Ga_2Mg_5 ($D8_g$) Structure: A2B5_oI28_72_j_bfj

http://aflow.org/prototype-encyclopedia/A2B5_oI28_72_j_bfj



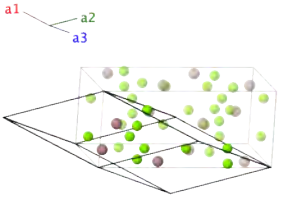
Prototype	:	Ga_2Mg_5
AFLOW prototype label	:	A2B5_oI28_72_j_bfj
***Strukturbericht* designation**	:	$D8_g$
Pearson symbol	:	oI28
Space group number	:	72
Space group symbol	:	*Ibam*
AFLOW prototype command	:	`aflow --proto=A2B5_oI28_72_j_bfj --params=`$a, b/a, c/a, x_2, x_3, y_3, x_4, y_4$

Other compounds with this structure

- As_2Cu_5, Tl_2Mg_5, and Mn_2Ge_5

Body-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = -\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}b\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}a\,\hat{\mathbf{x}} - \tfrac{1}{2}b\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}b\,\hat{\mathbf{y}} - \tfrac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\tfrac{1}{4}\mathbf{a}_1 + \tfrac{3}{4}\mathbf{a}_2 + \tfrac{1}{2}\mathbf{a}_3$	=	$\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{4}c\,\hat{\mathbf{z}}$	($4b$)	Mg I
$\mathbf{B}_2$	=	$\tfrac{3}{4}\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2 + \tfrac{1}{2}\mathbf{a}_3$	=	$\tfrac{1}{2}b\,\hat{\mathbf{y}} + \tfrac{1}{4}c\,\hat{\mathbf{z}}$	($4b$)	Mg I
$\mathbf{B}_3$	=	$\tfrac{1}{4}\mathbf{a}_1 + \left(\tfrac{1}{4} + x_2\right)\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + \tfrac{1}{4}c\,\hat{\mathbf{z}}$	($8f$)	Mg II
$\mathbf{B}_4$	=	$\tfrac{1}{4}\mathbf{a}_1 + \left(\tfrac{1}{4} - x_2\right)\mathbf{a}_2 - x_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \tfrac{1}{4}c\,\hat{\mathbf{z}}$	($8f$)	Mg II
$\mathbf{B}_5$	=	$\tfrac{3}{4}\mathbf{a}_1 + \left(\tfrac{3}{4} - x_2\right)\mathbf{a}_2 - x_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \tfrac{3}{4}c\,\hat{\mathbf{z}}$	($8f$)	Mg II
$\mathbf{B}_6$	=	$\tfrac{3}{4}\mathbf{a}_1 + \left(\tfrac{3}{4} + x_2\right)\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + \tfrac{3}{4}c\,\hat{\mathbf{z}}$	($8f$)	Mg II

$\mathbf{B}_7$	=	$y_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + (x_3 + y_3)\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}}$	$(8j)$	Ga
$\mathbf{B}_8$	=	$-y_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + (-x_3 - y_3)\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}}$	$(8j)$	Ga
$\mathbf{B}_9$	=	$\left(\frac{1}{2} + y_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 + (-x_3 + y_3)\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + y_3 b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(8j)$	Ga
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - y_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + (x_3 - y_3)\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} - y_3 b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(8j)$	Ga
$\mathbf{B}_{11}$	=	$y_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + (x_4 + y_4)\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}}$	$(8j)$	Mg III
$\mathbf{B}_{12}$	=	$-y_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + (-x_4 - y_4)\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}}$	$(8j)$	Mg III
$\mathbf{B}_{13}$	=	$\left(\frac{1}{2} + y_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_2 + (-x_4 + y_4)\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + y_4 b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(8j)$	Mg III
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - y_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_4\right)\mathbf{a}_2 + (x_4 - y_4)\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} - y_4 b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(8j)$	Mg III

References:

- K. Schubert, K. Frank, R. Gohle, A. Maldonado, H. G. Meissner, A. Raman, and W. Rossteutscher, *Einige Strukturdaten metallischer Phasen (8)*, Naturwissenschaften **50**, 41 (1963), doi:10.1007/BF00622812.

Found in:

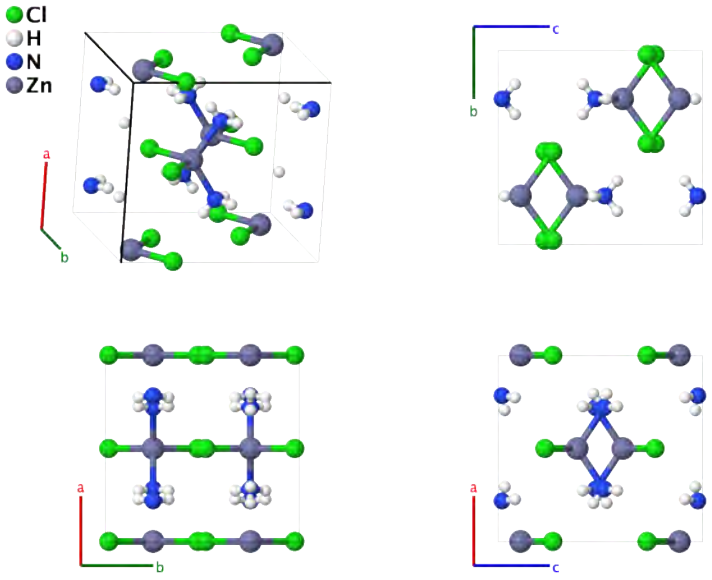
- V. Vreshch, *Crystal Structure of Ga_2Mg_5*, http://crystallography-online.com/structure/1522852 (2018). Crystallography online.com.

Geometry files:

- CIF: pp. 1665
- POSCAR: pp. 1666

$Zn(NH_3)_2Cl_2$ ($E1_2$) Structure: A2B6C2D_oI44_74_h_ij_i_e

http://aflow.org/prototype-encyclopedia/A2B6C2D_oI44_74_h_ij_i_e



Prototype	:	$Cl_2H_6N_2Zn$
AFLOW prototype label	:	A2B6C2D_oI44_74_h_ij_i_e
***Strukturbericht* designation**	:	$E1_2$
Pearson symbol	:	oI44
Space group number	:	74
Space group symbol	:	*Imma*
AFLOW prototype command	:	`aflow --proto=A2B6C2D_oI44_74_h_ij_i_e --params=`$a, b/a, c/a, z_1, y_2, z_2, x_3, z_3, x_4, z_4, x_5, y_5, z_5$

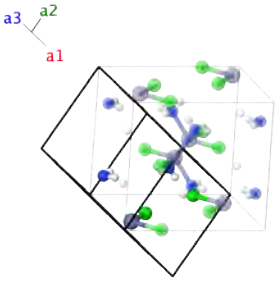
Other compounds with this structure

- $Zn(NH_3)_2Br_2$

- The recent work of (Ivšić, 2019) studied this system at 100 K and were able to locate the hydrogen atoms. The positions of the other atoms are similar to those in earlier works such as (Yamaguchi, 1981) and the space group is unchanged, so we use this as the prototype for the $E1_2$ label.

Body-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = -\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} - \frac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\left(\frac{1}{4}+z_1\right)\mathbf{a}_1+z_1\mathbf{a}_2+\frac{1}{4}\mathbf{a}_3$	=	$\frac{1}{4}b\,\hat{\mathbf{y}}+z_1c\,\hat{\mathbf{z}}$	(4*e*)	Zn
$\mathbf{B}_2$	=	$\left(\frac{3}{4}-z_1\right)\mathbf{a}_1-z_1\mathbf{a}_2+\frac{3}{4}\mathbf{a}_3$	=	$\frac{3}{4}b\,\hat{\mathbf{y}}-z_1c\,\hat{\mathbf{z}}$	(4*e*)	Zn
$\mathbf{B}_3$	=	$(y_2+z_2)\mathbf{a}_1+z_2\mathbf{a}_2+y_2\mathbf{a}_3$	=	$y_2b\,\hat{\mathbf{y}}+z_2c\,\hat{\mathbf{z}}$	(8*h*)	Cl
$\mathbf{B}_4$	=	$\left(\frac{1}{2}-y_2+z_2\right)\mathbf{a}_1+z_2\mathbf{a}_2+\left(\frac{1}{2}-y_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_2\right)b\,\hat{\mathbf{y}}+z_2c\,\hat{\mathbf{z}}$	(8*h*)	Cl
$\mathbf{B}_5$	=	$\left(\frac{1}{2}+y_2-z_2\right)\mathbf{a}_1-z_2\mathbf{a}_2+\left(\frac{1}{2}+y_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_2\right)b\,\hat{\mathbf{y}}-z_2c\,\hat{\mathbf{z}}$	(8*h*)	Cl
$\mathbf{B}_6$	=	$(-y_2-z_2)\mathbf{a}_1-z_2\mathbf{a}_2-y_2\mathbf{a}_3$	=	$-y_2b\,\hat{\mathbf{y}}-z_2c\,\hat{\mathbf{z}}$	(8*h*)	Cl
$\mathbf{B}_7$	=	$\left(\frac{1}{4}+z_3\right)\mathbf{a}_1+(x_3+z_3)\mathbf{a}_2+\left(\frac{1}{4}+x_3\right)\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+z_3c\,\hat{\mathbf{z}}$	(8*i*)	H I
$\mathbf{B}_8$	=	$\left(\frac{1}{4}+z_3\right)\mathbf{a}_1+(-x_3+z_3)\mathbf{a}_2+\left(\frac{1}{4}-x_3\right)\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+z_3c\,\hat{\mathbf{z}}$	(8*i*)	H I
$\mathbf{B}_9$	=	$\left(\frac{3}{4}-z_3\right)\mathbf{a}_1+(-x_3-z_3)\mathbf{a}_2+\left(\frac{3}{4}-x_3\right)\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}-z_3c\,\hat{\mathbf{z}}$	(8*i*)	H I
$\mathbf{B}_{10}$	=	$\left(\frac{3}{4}-z_3\right)\mathbf{a}_1+(x_3-z_3)\mathbf{a}_2+\left(\frac{3}{4}+x_3\right)\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}-z_3c\,\hat{\mathbf{z}}$	(8*i*)	H I
$\mathbf{B}_{11}$	=	$\left(\frac{1}{4}+z_4\right)\mathbf{a}_1+(x_4+z_4)\mathbf{a}_2+\left(\frac{1}{4}+x_4\right)\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+z_4c\,\hat{\mathbf{z}}$	(8*i*)	N
$\mathbf{B}_{12}$	=	$\left(\frac{1}{4}+z_4\right)\mathbf{a}_1+(-x_4+z_4)\mathbf{a}_2+\left(\frac{1}{4}-x_4\right)\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}}+\frac{1}{4}b\,\hat{\mathbf{y}}+z_4c\,\hat{\mathbf{z}}$	(8*i*)	N
$\mathbf{B}_{13}$	=	$\left(\frac{3}{4}-z_4\right)\mathbf{a}_1+(-x_4-z_4)\mathbf{a}_2+\left(\frac{3}{4}-x_4\right)\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}-z_4c\,\hat{\mathbf{z}}$	(8*i*)	N
$\mathbf{B}_{14}$	=	$\left(\frac{3}{4}-z_4\right)\mathbf{a}_1+(x_4-z_4)\mathbf{a}_2+\left(\frac{3}{4}+x_4\right)\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}}+\frac{3}{4}b\,\hat{\mathbf{y}}-z_4c\,\hat{\mathbf{z}}$	(8*i*)	N
$\mathbf{B}_{15}$	=	$(y_5+z_5)\mathbf{a}_1+(x_5+z_5)\mathbf{a}_2+(x_5+y_5)\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}}+y_5b\,\hat{\mathbf{y}}+z_5c\,\hat{\mathbf{z}}$	(16*j*)	H II
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2}-y_5+z_5\right)\mathbf{a}_1+(-x_5+z_5)\mathbf{a}_2+\left(\frac{1}{2}-x_5-y_5\right)\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_5\right)b\,\hat{\mathbf{y}}+z_5c\,\hat{\mathbf{z}}$	(16*j*)	H II
$\mathbf{B}_{17}$	=	$\left(\frac{1}{2}+y_5-z_5\right)\mathbf{a}_1+(-x_5-z_5)\mathbf{a}_2+\left(\frac{1}{2}-x_5+y_5\right)\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_5\right)b\,\hat{\mathbf{y}}-z_5c\,\hat{\mathbf{z}}$	(16*j*)	H II
$\mathbf{B}_{18}$	=	$(-y_5-z_5)\mathbf{a}_1+(x_5-z_5)\mathbf{a}_2+(x_5-y_5)\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}}-y_5b\,\hat{\mathbf{y}}-z_5c\,\hat{\mathbf{z}}$	(16*j*)	H II
$\mathbf{B}_{19}$	=	$(-y_5-z_5)\mathbf{a}_1+(-x_5-z_5)\mathbf{a}_2+(-x_5-y_5)\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}}-y_5b\,\hat{\mathbf{y}}-z_5c\,\hat{\mathbf{z}}$	(16*j*)	H II
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2}+y_5-z_5\right)\mathbf{a}_1+(x_5-z_5)\mathbf{a}_2+\left(\frac{1}{2}+x_5+y_5\right)\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_5\right)b\,\hat{\mathbf{y}}-z_5c\,\hat{\mathbf{z}}$	(16*j*)	H II
$\mathbf{B}_{21}$	=	$\left(\frac{1}{2}-y_5+z_5\right)\mathbf{a}_1+(x_5+z_5)\mathbf{a}_2+\left(\frac{1}{2}+x_5-y_5\right)\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_5\right)b\,\hat{\mathbf{y}}+z_5c\,\hat{\mathbf{z}}$	(16*j*)	H II
$\mathbf{B}_{22}$	=	$(y_5+z_5)\mathbf{a}_1+(-x_5+z_5)\mathbf{a}_2+(-x_5+y_5)\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}}+y_5b\,\hat{\mathbf{y}}+z_5c\,\hat{\mathbf{z}}$	(16*j*)	H II

References:

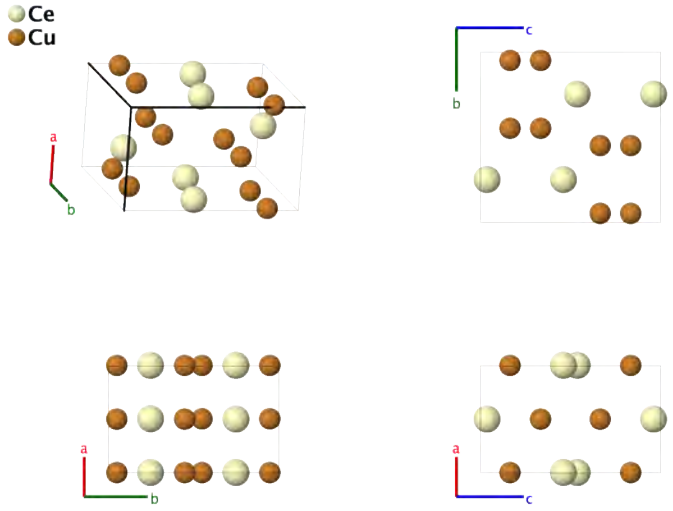
- T. Ivšić, D. W. Bi, and A. Magrez, *New refinement of the crystal structure of $Zn(NH_3)_2Cl_2$ at 100K*, Acta Crystallogr. E **75**, 1386–1388 (2019), doi:10.1107/S2056989019011757.
- T. Yamaguchi and O. Lindqvist, *The Crystal Structure of Diamminedichlorozinc(II), $ZnCl_2(NH_3)_2$. A New Refinement.*, Acta Chem. Scand. **35a**, 727–728 (1981), doi:10.3891/acta.chem.scand.35a-0727.

Geometry files:

- CIF: pp. 1666
- POSCAR: pp. 1666

$CeCu_2$ Structure: AB2_oI12_74_e_h

http://aflow.org/prototype-encyclopedia/AB2_oI12_74_e_h



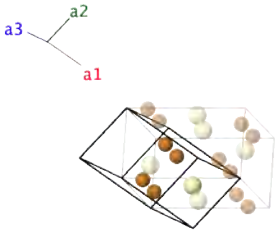
Prototype	:	$CeCu_2$
AFLOW prototype label	:	AB2_oI12_74_e_h
***Strukturbericht* designation**	:	None
Pearson symbol	:	oI12
Space group number	:	74
Space group symbol	:	*Imma*
AFLOW prototype command	:	`aflow --proto=AB2_oI12_74_e_h --params=`$a, b/a, c/a, z_1, y_2, z_2$

Other compounds with this structure

- $CeCu_2$, $PrCu_2$, $NdCu_2$, $SmCu_2$, $EuCu_2$, $GdCu_2$, $TbCu_2$, $DyCu_2$, $HoCu_2$, $ErCu_2$, $YbCu_2$, $LuCu_2$, $LaZn_2$, $CeZn_2$, $PrZn_2$, $NdZn_2$, $SmZn_2$, $EuZn_2$, $GdZn_2$, $ThZn_2$, $DyZn_2$, $HoZn_2$, $ErZn_2$, $TmZn_2$, $YbZn_2$, $LuZn_2$, $LaAg_2$, $CeAg_2$, $PrAg_2$, $NdAg_2$, $EuAg_2$, $YbAg_2$, $LaAu_2$, $CeAu_2$, and $EuAu_2$

Body-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = -\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} - \frac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$ =	$\left(\frac{1}{4} + z_1\right)\mathbf{a}_1 + z_1\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}b\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	($4e$)	Ce

$\mathbf{B}_2$	=	$\left(\frac{3}{4} - z_1\right)\mathbf{a}_1 - z_1\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}b\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	($4e$)	Ce
$\mathbf{B}_3$	=	$(y_2 + z_2)\,\mathbf{a}_1 + z_2\,\mathbf{a}_2 + y_2\,\mathbf{a}_3$	=	$y_2 b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	($8h$)	Cu
$\mathbf{B}_4$	=	$\left(\frac{1}{2} - y_2 + z_2\right)\mathbf{a}_1 + z_2\,\mathbf{a}_2 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_2\right)b\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	($8h$)	Cu
$\mathbf{B}_5$	=	$\left(\frac{1}{2} + y_2 - z_2\right)\mathbf{a}_1 - z_2\,\mathbf{a}_2 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_2\right)b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	($8h$)	Cu
$\mathbf{B}_6$	=	$(-y_2 - z_2)\,\mathbf{a}_1 - z_2\,\mathbf{a}_2 - y_2\,\mathbf{a}_3$	=	$-y_2 b\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	($8h$)	Cu

References:

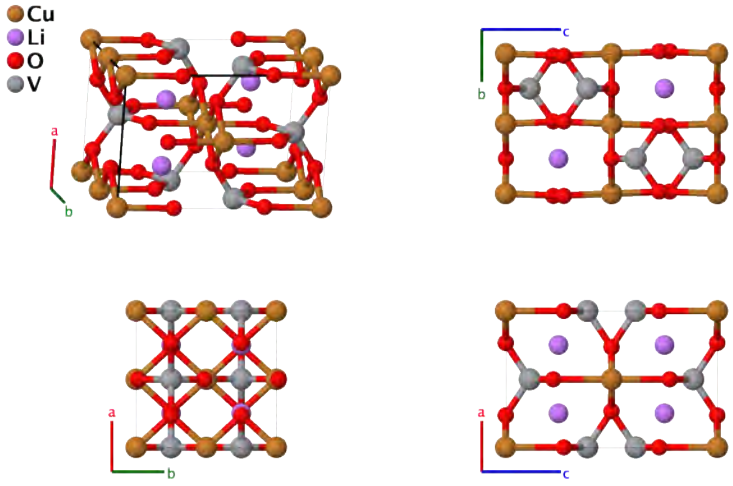
- A. C. Larson and D. T. Cromer, *The crystal structure of $CeCu_2$*, Acta Cryst. **14**, 73–74 (1961), doi:10.1107/S0365110X61000231.
- D. Debray, *Crystal Chemistry of the $CeCu_2$-type structure*, J. Less-Common Met. **30**, 237–248 (1973), doi:10.1016/0022-5088(73)90110-0.

Geometry files:

- CIF: pp. 1666
- POSCAR: pp. 1667

$LiCuVO_4$ Structure: ABC4D_oI28_74_a_d_hi_e

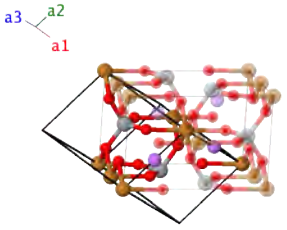
http://aflow.org/prototype-encyclopedia/ABC4D_oI28_74_a_d_hi_e



Prototype	:	$CuLiO_4V$
AFLOW prototype label	:	ABC4D_oI28_74_a_d_hi_e
***Strukturbericht* designation**	:	None
Pearson symbol	:	oI28
Space group number	:	74
Space group symbol	:	*Imma*
AFLOW prototype command	:	`aflow --proto=ABC4D_oI28_74_a_d_hi_e --params=`$a, b/a, c/a, z_3, y_4, z_4, x_5, z_5$

Body-centered Orthorhombic primitive vectors:

$$\mathbf{a}_1 = -\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} - \frac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(4a)$	Cu
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}b\,\hat{\mathbf{y}}$	$(4a)$	Cu
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}b\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(4d)$	Li
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} - \frac{1}{4}b\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(4d)$	Li
$\mathbf{B}_5$	=	$\left(\frac{1}{4} + z_3\right)\mathbf{a}_1 + z_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}b\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(4e)$	V
$\mathbf{B}_6$	=	$\left(\frac{3}{4} - z_3\right)\mathbf{a}_1 - z_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}b\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(4e)$	V
$\mathbf{B}_7$	=	$(y_4 + z_4)\,\mathbf{a}_1 + z_4\,\mathbf{a}_2 + y_4\,\mathbf{a}_3$	=	$y_4 b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(8h)$	O I

$\mathbf{B}_8$	=	$\left(\frac{1}{2} - y_4 + z_4\right)\mathbf{a}_1 + z_4\,\mathbf{a}_2 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_4\right) b\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*h*)	O I
$\mathbf{B}_9$	=	$\left(\frac{1}{2} + y_4 - z_4\right)\mathbf{a}_1 - z_4\,\mathbf{a}_2 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_4\right) b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*h*)	O I
$\mathbf{B}_{10}$	=	$(-y_4 - z_4)\,\mathbf{a}_1 - z_4\,\mathbf{a}_2 - y_4\,\mathbf{a}_3$	=	$-y_4 b\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*h*)	O I
$\mathbf{B}_{11}$	=	$\left(\frac{1}{4} + z_5\right)\mathbf{a}_1 + (x_5 + z_5)\,\mathbf{a}_2 + \left(\frac{1}{4} + x_5\right)\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(8*i*)	O II
$\mathbf{B}_{12}$	=	$\left(\frac{1}{4} + z_5\right)\mathbf{a}_1 + (-x_5 + z_5)\,\mathbf{a}_2 + \left(\frac{1}{4} - x_5\right)\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \frac{1}{4} b\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(8*i*)	O II
$\mathbf{B}_{13}$	=	$\left(\frac{3}{4} - z_5\right)\mathbf{a}_1 + (-x_5 - z_5)\,\mathbf{a}_2 + \left(\frac{3}{4} - x_5\right)\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(8*i*)	O II
$\mathbf{B}_{14}$	=	$\left(\frac{3}{4} - z_5\right)\mathbf{a}_1 + (x_5 - z_5)\,\mathbf{a}_2 + \left(\frac{3}{4} + x_5\right)\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + \frac{3}{4} b\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(8*i*)	O II

References:

- M. A. Lafontaine, M. Leblanc, and G. Ferey, *New refinement of the room-temperature structure of $LiCuVO_4$*, Acta Crystallogr. C **45**, 1205–1206 (1989), doi:10.1107/S0108270189001551.

Found in:

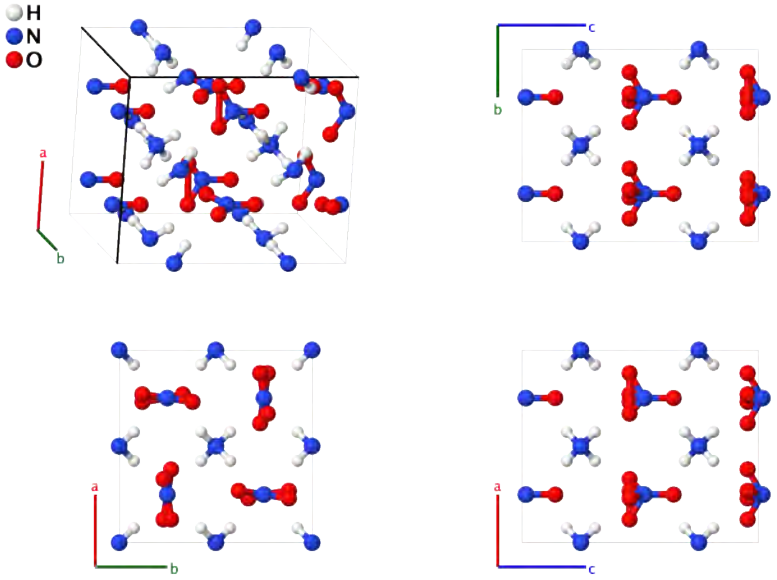
- A. V. Prokofiev, I. G. Vasilyeva, V. N. Ikorskii, V. V. Malakhov, I. P. Asanov, and W. Assmus, *Structure, stoichiometry and magnetic properties of the low-dimensional structure phase $LiCuVO_4$*, J. Solid State Chem. **177**, 3131–3139 (2004), doi:10.1016/j.jssc.2004.05.031.

Geometry files:

- CIF: pp. 1667
- POSCAR: pp. 1667

Gwihabaite [NH_4NO_3 (V)] Structure: A4B2C3_tP72_77_8d_ab2c2d_6d

http://aflow.org/prototype-encyclopedia/A4B2C3_tP72_77_8d_ab2c2d_6d



Prototype	:	$H_4N_2O_3$
AFLOW prototype label	:	A4B2C3_tP72_77_8d_ab2c2d_6d
***Strukturbericht* designation**	:	None
Pearson symbol	:	tP72
Space group number	:	77
Space group symbol	:	$P4_2$
AFLOW prototype command	:	`aflow --proto=A4B2C3_tP72_77_8d_ab2c2d_6d --params=`$a, c/a, z_1, z_2, z_3, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}, x_{17}, y_{17}, z_{17}, x_{18}, y_{18}, z_{18}, x_{19}, y_{19}, z_{19}, x_{20}, y_{20}, z_{20}$

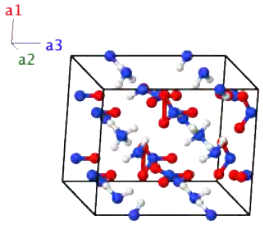
- Ammonium Nitrate exists in a variety of forms, (Hermann, 1937) depending on the temperature:

Phase	Temperature °C	*Strukturbericht*	Page
I	125 – 170	$G0_8$	AB_cP2_221_a_b.NH4.NO3
II	84 – 125	$G0_9$	ABC3_tP10_100_b_a_bc
III	32 – 84	$G0_{10}$	ABC3_oP20_62_c_c_cd.N.NH4.O
IV	-17 – 32	$G0_{11}$	A4B2C3_oP18_59_ef_ab_af
V	< −17	Gwihabaite	A4B2C3_tP72_77_8d_ab2c2d_6d2

- Data for this structure was taken at −150 °C.

Simple Tetragonal primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= a\,\hat{\mathbf{x}}\\ \mathbf{a}_2 &= a\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$z_1\,\mathbf{a}_3$	=	$z_1 c\,\hat{\mathbf{z}}$	$(2a)$	N I
$\mathbf{B}_2$	=	$\left(\frac{1}{2}+z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+z_1\right)c\,\hat{\mathbf{z}}$	$(2a)$	N I
$\mathbf{B}_3$	=	$\frac{1}{2}\mathbf{a}_1+\frac{1}{2}\mathbf{a}_2+z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}+\frac{1}{2}a\,\hat{\mathbf{y}}+z_2 c\,\hat{\mathbf{z}}$	$(2b)$	N II
$\mathbf{B}_4$	=	$\frac{1}{2}\mathbf{a}_1+\frac{1}{2}\mathbf{a}_2+\left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}+\frac{1}{2}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_2\right)c\,\hat{\mathbf{z}}$	$(2b)$	N II
$\mathbf{B}_5$	=	$\frac{1}{2}\mathbf{a}_2+z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}}+z_3 c\,\hat{\mathbf{z}}$	$(2c)$	N III
$\mathbf{B}_6$	=	$\frac{1}{2}\mathbf{a}_1+\left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+z_3\right)c\,\hat{\mathbf{z}}$	$(2c)$	N III
$\mathbf{B}_7$	=	$\frac{1}{2}\mathbf{a}_2+z_4\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}}+z_4 c\,\hat{\mathbf{z}}$	$(2c)$	N IV
$\mathbf{B}_8$	=	$\frac{1}{2}\mathbf{a}_1+\left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+z_4\right)c\,\hat{\mathbf{z}}$	$(2c)$	N IV
$\mathbf{B}_9$	=	$x_5\,\mathbf{a}_1+y_5\,\mathbf{a}_2+z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}}+y_5 a\,\hat{\mathbf{y}}+z_5 c\,\hat{\mathbf{z}}$	$(4d)$	H I
$\mathbf{B}_{10}$	=	$-x_5\,\mathbf{a}_1-y_5\,\mathbf{a}_2+z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}}-y_5 a\,\hat{\mathbf{y}}+z_5 c\,\hat{\mathbf{z}}$	$(4d)$	H I
$\mathbf{B}_{11}$	=	$-y_5\,\mathbf{a}_1+x_5\,\mathbf{a}_2+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$-y_5 a\,\hat{\mathbf{x}}+x_5 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	$(4d)$	H I
$\mathbf{B}_{12}$	=	$y_5\,\mathbf{a}_1-x_5\,\mathbf{a}_2+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$y_5 a\,\hat{\mathbf{x}}-x_5 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	$(4d)$	H I
$\mathbf{B}_{13}$	=	$x_6\,\mathbf{a}_1+y_6\,\mathbf{a}_2+z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}}+y_6 a\,\hat{\mathbf{y}}+z_6 c\,\hat{\mathbf{z}}$	$(4d)$	H II
$\mathbf{B}_{14}$	=	$-x_6\,\mathbf{a}_1-y_6\,\mathbf{a}_2+z_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}}-y_6 a\,\hat{\mathbf{y}}+z_6 c\,\hat{\mathbf{z}}$	$(4d)$	H II
$\mathbf{B}_{15}$	=	$-y_6\,\mathbf{a}_1+x_6\,\mathbf{a}_2+\left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	=	$-y_6 a\,\hat{\mathbf{x}}+x_6 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	$(4d)$	H II
$\mathbf{B}_{16}$	=	$y_6\,\mathbf{a}_1-x_6\,\mathbf{a}_2+\left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	=	$y_6 a\,\hat{\mathbf{x}}-x_6 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	$(4d)$	H II
$\mathbf{B}_{17}$	=	$x_7\,\mathbf{a}_1+y_7\,\mathbf{a}_2+z_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}}+y_7 a\,\hat{\mathbf{y}}+z_7 c\,\hat{\mathbf{z}}$	$(4d)$	H III
$\mathbf{B}_{18}$	=	$-x_7\,\mathbf{a}_1-y_7\,\mathbf{a}_2+z_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}}-y_7 a\,\hat{\mathbf{y}}+z_7 c\,\hat{\mathbf{z}}$	$(4d)$	H III
$\mathbf{B}_{19}$	=	$-y_7\,\mathbf{a}_1+x_7\,\mathbf{a}_2+\left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	=	$-y_7 a\,\hat{\mathbf{x}}+x_7 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	$(4d)$	H III
$\mathbf{B}_{20}$	=	$y_7\,\mathbf{a}_1-x_7\,\mathbf{a}_2+\left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	=	$y_7 a\,\hat{\mathbf{x}}-x_7 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	$(4d)$	H III
$\mathbf{B}_{21}$	=	$x_8\,\mathbf{a}_1+y_8\,\mathbf{a}_2+z_8\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}}+y_8 a\,\hat{\mathbf{y}}+z_8 c\,\hat{\mathbf{z}}$	$(4d)$	H IV
$\mathbf{B}_{22}$	=	$-x_8\,\mathbf{a}_1-y_8\,\mathbf{a}_2+z_8\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}}-y_8 a\,\hat{\mathbf{y}}+z_8 c\,\hat{\mathbf{z}}$	$(4d)$	H IV
$\mathbf{B}_{23}$	=	$-y_8\,\mathbf{a}_1+x_8\,\mathbf{a}_2+\left(\frac{1}{2}+z_8\right)\mathbf{a}_3$	=	$-y_8 a\,\hat{\mathbf{x}}+x_8 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_8\right)c\,\hat{\mathbf{z}}$	$(4d)$	H IV
$\mathbf{B}_{24}$	=	$y_8\,\mathbf{a}_1-x_8\,\mathbf{a}_2+\left(\frac{1}{2}+z_8\right)\mathbf{a}_3$	=	$y_8 a\,\hat{\mathbf{x}}-x_8 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_8\right)c\,\hat{\mathbf{z}}$	$(4d)$	H IV
$\mathbf{B}_{25}$	=	$x_9\,\mathbf{a}_1+y_9\,\mathbf{a}_2+z_9\,\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}}+y_9 a\,\hat{\mathbf{y}}+z_9 c\,\hat{\mathbf{z}}$	$(4d)$	H V
$\mathbf{B}_{26}$	=	$-x_9\,\mathbf{a}_1-y_9\,\mathbf{a}_2+z_9\,\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{x}}-y_9 a\,\hat{\mathbf{y}}+z_9 c\,\hat{\mathbf{z}}$	$(4d)$	H V
$\mathbf{B}_{27}$	=	$-y_9\,\mathbf{a}_1+x_9\,\mathbf{a}_2+\left(\frac{1}{2}+z_9\right)\mathbf{a}_3$	=	$-y_9 a\,\hat{\mathbf{x}}+x_9 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_9\right)c\,\hat{\mathbf{z}}$	$(4d)$	H V

$$
\begin{aligned}
\mathbf{B}_{28} &= y_9\,\mathbf{a}_1 - x_9\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_3 &&= y_9 a\,\hat{\mathbf{x}} - x_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_9\right) c\,\hat{\mathbf{z}} && (4d) && \text{H V}\\
\mathbf{B}_{29} &= x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 &&= x_{10} a\,\hat{\mathbf{x}} + y_{10} a\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}} && (4d) && \text{H VI}\\
\mathbf{B}_{30} &= -x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 &&= -x_{10} a\,\hat{\mathbf{x}} - y_{10} a\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}} && (4d) && \text{H VI}\\
\mathbf{B}_{31} &= -y_{10}\,\mathbf{a}_1 + x_{10}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{10}\right)\mathbf{a}_3 &&= -y_{10} a\,\hat{\mathbf{x}} + x_{10} a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{10}\right) c\,\hat{\mathbf{z}} && (4d) && \text{H VI}\\
\mathbf{B}_{32} &= y_{10}\,\mathbf{a}_1 - x_{10}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{10}\right)\mathbf{a}_3 &&= y_{10} a\,\hat{\mathbf{x}} - x_{10} a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{10}\right) c\,\hat{\mathbf{z}} && (4d) && \text{H VI}\\
\mathbf{B}_{33} &= x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 &&= x_{11} a\,\hat{\mathbf{x}} + y_{11} a\,\hat{\mathbf{y}} + z_{11} c\,\hat{\mathbf{z}} && (4d) && \text{H VII}\\
\mathbf{B}_{34} &= -x_{11}\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 &&= -x_{11} a\,\hat{\mathbf{x}} - y_{11} a\,\hat{\mathbf{y}} + z_{11} c\,\hat{\mathbf{z}} && (4d) && \text{H VII}\\
\mathbf{B}_{35} &= -y_{11}\,\mathbf{a}_1 + x_{11}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{11}\right)\mathbf{a}_3 &&= -y_{11} a\,\hat{\mathbf{x}} + x_{11} a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{11}\right) c\,\hat{\mathbf{z}} && (4d) && \text{H VII}\\
\mathbf{B}_{36} &= y_{11}\,\mathbf{a}_1 - x_{11}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{11}\right)\mathbf{a}_3 &&= y_{11} a\,\hat{\mathbf{x}} - x_{11} a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{11}\right) c\,\hat{\mathbf{z}} && (4d) && \text{H VII}\\
\mathbf{B}_{37} &= x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 &&= x_{12} a\,\hat{\mathbf{x}} + y_{12} a\,\hat{\mathbf{y}} + z_{12} c\,\hat{\mathbf{z}} && (4d) && \text{H VIII}\\
\mathbf{B}_{38} &= -x_{12}\,\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 &&= -x_{12} a\,\hat{\mathbf{x}} - y_{12} a\,\hat{\mathbf{y}} + z_{12} c\,\hat{\mathbf{z}} && (4d) && \text{H VIII}\\
\mathbf{B}_{39} &= -y_{12}\,\mathbf{a}_1 + x_{12}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{12}\right)\mathbf{a}_3 &&= -y_{12} a\,\hat{\mathbf{x}} + x_{12} a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{12}\right) c\,\hat{\mathbf{z}} && (4d) && \text{H VIII}\\
\mathbf{B}_{40} &= y_{12}\,\mathbf{a}_1 - x_{12}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{12}\right)\mathbf{a}_3 &&= y_{12} a\,\hat{\mathbf{x}} - x_{12} a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{12}\right) c\,\hat{\mathbf{z}} && (4d) && \text{H VIII}\\
\mathbf{B}_{41} &= x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3 &&= x_{13} a\,\hat{\mathbf{x}} + y_{13} a\,\hat{\mathbf{y}} + z_{13} c\,\hat{\mathbf{z}} && (4d) && \text{N V}\\
\mathbf{B}_{42} &= -x_{13}\,\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3 &&= -x_{13} a\,\hat{\mathbf{x}} - y_{13} a\,\hat{\mathbf{y}} + z_{13} c\,\hat{\mathbf{z}} && (4d) && \text{N V}\\
\mathbf{B}_{43} &= -y_{13}\,\mathbf{a}_1 + x_{13}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{13}\right)\mathbf{a}_3 &&= -y_{13} a\,\hat{\mathbf{x}} + x_{13} a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{13}\right) c\,\hat{\mathbf{z}} && (4d) && \text{N V}\\
\mathbf{B}_{44} &= y_{13}\,\mathbf{a}_1 - x_{13}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{13}\right)\mathbf{a}_3 &&= y_{13} a\,\hat{\mathbf{x}} - x_{13} a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{13}\right) c\,\hat{\mathbf{z}} && (4d) && \text{N V}\\
\mathbf{B}_{45} &= x_{14}\,\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3 &&= x_{14} a\,\hat{\mathbf{x}} + y_{14} a\,\hat{\mathbf{y}} + z_{14} c\,\hat{\mathbf{z}} && (4d) && \text{N VI}\\
\mathbf{B}_{46} &= -x_{14}\,\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3 &&= -x_{14} a\,\hat{\mathbf{x}} - y_{14} a\,\hat{\mathbf{y}} + z_{14} c\,\hat{\mathbf{z}} && (4d) && \text{N VI}\\
\mathbf{B}_{47} &= -y_{14}\,\mathbf{a}_1 + x_{14}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{14}\right)\mathbf{a}_3 &&= -y_{14} a\,\hat{\mathbf{x}} + x_{14} a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{14}\right) c\,\hat{\mathbf{z}} && (4d) && \text{N VI}\\
\mathbf{B}_{48} &= y_{14}\,\mathbf{a}_1 - x_{14}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{14}\right)\mathbf{a}_3 &&= y_{14} a\,\hat{\mathbf{x}} - x_{14} a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{14}\right) c\,\hat{\mathbf{z}} && (4d) && \text{N VI}\\
\mathbf{B}_{49} &= x_{15}\,\mathbf{a}_1 + y_{15}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 &&= x_{15} a\,\hat{\mathbf{x}} + y_{15} a\,\hat{\mathbf{y}} + z_{15} c\,\hat{\mathbf{z}} && (4d) && \text{O I}\\
\mathbf{B}_{50} &= -x_{15}\,\mathbf{a}_1 - y_{15}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 &&= -x_{15} a\,\hat{\mathbf{x}} - y_{15} a\,\hat{\mathbf{y}} + z_{15} c\,\hat{\mathbf{z}} && (4d) && \text{O I}\\
\mathbf{B}_{51} &= -y_{15}\,\mathbf{a}_1 + x_{15}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{15}\right)\mathbf{a}_3 &&= -y_{15} a\,\hat{\mathbf{x}} + x_{15} a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{15}\right) c\,\hat{\mathbf{z}} && (4d) && \text{O I}\\
\mathbf{B}_{52} &= y_{15}\,\mathbf{a}_1 - x_{15}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{15}\right)\mathbf{a}_3 &&= y_{15} a\,\hat{\mathbf{x}} - x_{15} a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{15}\right) c\,\hat{\mathbf{z}} && (4d) && \text{O I}\\
\mathbf{B}_{53} &= x_{16}\,\mathbf{a}_1 + y_{16}\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 &&= x_{16} a\,\hat{\mathbf{x}} + y_{16} a\,\hat{\mathbf{y}} + z_{16} c\,\hat{\mathbf{z}} && (4d) && \text{O II}\\
\mathbf{B}_{54} &= -x_{16}\,\mathbf{a}_1 - y_{16}\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 &&= -x_{16} a\,\hat{\mathbf{x}} - y_{16} a\,\hat{\mathbf{y}} + z_{16} c\,\hat{\mathbf{z}} && (4d) && \text{O II}\\
\mathbf{B}_{55} &= -y_{16}\,\mathbf{a}_1 + x_{16}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{16}\right)\mathbf{a}_3 &&= -y_{16} a\,\hat{\mathbf{x}} + x_{16} a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{16}\right) c\,\hat{\mathbf{z}} && (4d) && \text{O II}\\
\mathbf{B}_{56} &= y_{16}\,\mathbf{a}_1 - x_{16}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{16}\right)\mathbf{a}_3 &&= y_{16} a\,\hat{\mathbf{x}} - x_{16} a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{16}\right) c\,\hat{\mathbf{z}} && (4d) && \text{O II}\\
\mathbf{B}_{57} &= x_{17}\,\mathbf{a}_1 + y_{17}\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 &&= x_{17} a\,\hat{\mathbf{x}} + y_{17} a\,\hat{\mathbf{y}} + z_{17} c\,\hat{\mathbf{z}} && (4d) && \text{O III}\\
\mathbf{B}_{58} &= -x_{17}\,\mathbf{a}_1 - y_{17}\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 &&= -x_{17} a\,\hat{\mathbf{x}} - y_{17} a\,\hat{\mathbf{y}} + z_{17} c\,\hat{\mathbf{z}} && (4d) && \text{O III}\\
\mathbf{B}_{59} &= -y_{17}\,\mathbf{a}_1 + x_{17}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{17}\right)\mathbf{a}_3 &&= -y_{17} a\,\hat{\mathbf{x}} + x_{17} a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{17}\right) c\,\hat{\mathbf{z}} && (4d) && \text{O III}\\
\mathbf{B}_{60} &= y_{17}\,\mathbf{a}_1 - x_{17}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{17}\right)\mathbf{a}_3 &&= y_{17} a\,\hat{\mathbf{x}} - x_{17} a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{17}\right) c\,\hat{\mathbf{z}} && (4d) && \text{O III}\\
\mathbf{B}_{61} &= x_{18}\,\mathbf{a}_1 + y_{18}\,\mathbf{a}_2 + z_{18}\,\mathbf{a}_3 &&= x_{18} a\,\hat{\mathbf{x}} + y_{18} a\,\hat{\mathbf{y}} + z_{18} c\,\hat{\mathbf{z}} && (4d) && \text{O IV}\\
\mathbf{B}_{62} &= -x_{18}\,\mathbf{a}_1 - y_{18}\,\mathbf{a}_2 + z_{18}\,\mathbf{a}_3 &&= -x_{18} a\,\hat{\mathbf{x}} - y_{18} a\,\hat{\mathbf{y}} + z_{18} c\,\hat{\mathbf{z}} && (4d) && \text{O IV}\\
\mathbf{B}_{63} &= -y_{18}\,\mathbf{a}_1 + x_{18}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{18}\right)\mathbf{a}_3 &&= -y_{18} a\,\hat{\mathbf{x}} + x_{18} a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{18}\right) c\,\hat{\mathbf{z}} && (4d) && \text{O IV}
\end{aligned}
$$

$\mathbf{B}_{64}$	=	$y_{18}\,\mathbf{a}_1 - x_{18}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{18}\right)\mathbf{a}_3$	=	$y_{18}a\,\hat{\mathbf{x}} - x_{18}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{18}\right)c\,\hat{\mathbf{z}}$	(4*d*)	O IV
$\mathbf{B}_{65}$	=	$x_{19}\,\mathbf{a}_1 + y_{19}\,\mathbf{a}_2 + z_{19}\,\mathbf{a}_3$	=	$x_{19}a\,\hat{\mathbf{x}} + y_{19}a\,\hat{\mathbf{y}} + z_{19}c\,\hat{\mathbf{z}}$	(4*d*)	O V
$\mathbf{B}_{66}$	=	$-x_{19}\,\mathbf{a}_1 - y_{19}\,\mathbf{a}_2 + z_{19}\,\mathbf{a}_3$	=	$-x_{19}a\,\hat{\mathbf{x}} - y_{19}a\,\hat{\mathbf{y}} + z_{19}c\,\hat{\mathbf{z}}$	(4*d*)	O V
$\mathbf{B}_{67}$	=	$-y_{19}\,\mathbf{a}_1 + x_{19}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{19}\right)\mathbf{a}_3$	=	$-y_{19}a\,\hat{\mathbf{x}} + x_{19}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{19}\right)c\,\hat{\mathbf{z}}$	(4*d*)	O V
$\mathbf{B}_{68}$	=	$y_{19}\,\mathbf{a}_1 - x_{19}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{19}\right)\mathbf{a}_3$	=	$y_{19}a\,\hat{\mathbf{x}} - x_{19}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{19}\right)c\,\hat{\mathbf{z}}$	(4*d*)	O V
$\mathbf{B}_{69}$	=	$x_{20}\,\mathbf{a}_1 + y_{20}\,\mathbf{a}_2 + z_{20}\,\mathbf{a}_3$	=	$x_{20}a\,\hat{\mathbf{x}} + y_{20}a\,\hat{\mathbf{y}} + z_{20}c\,\hat{\mathbf{z}}$	(4*d*)	O VI
$\mathbf{B}_{70}$	=	$-x_{20}\,\mathbf{a}_1 - y_{20}\,\mathbf{a}_2 + z_{20}\,\mathbf{a}_3$	=	$-x_{20}a\,\hat{\mathbf{x}} - y_{20}a\,\hat{\mathbf{y}} + z_{20}c\,\hat{\mathbf{z}}$	(4*d*)	O VI
$\mathbf{B}_{71}$	=	$-y_{20}\,\mathbf{a}_1 + x_{20}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{20}\right)\mathbf{a}_3$	=	$-y_{20}a\,\hat{\mathbf{x}} + x_{20}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{20}\right)c\,\hat{\mathbf{z}}$	(4*d*)	O VI
$\mathbf{B}_{72}$	=	$y_{20}\,\mathbf{a}_1 - x_{20}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{20}\right)\mathbf{a}_3$	=	$y_{20}a\,\hat{\mathbf{x}} - x_{20}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{20}\right)c\,\hat{\mathbf{z}}$	(4*d*)	O VI

References:

- J. L. Amorós, F. Arrese, and M. Canut, *The crystal structure of the low-temperature phase of NH_4NO_3 (V) at – 150°C*, Zeitschrift für Kristallographie - Crystalline Materials **117**, 92–107 (1962), doi:10.1524/zkri.1962.117.2-3.92.
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

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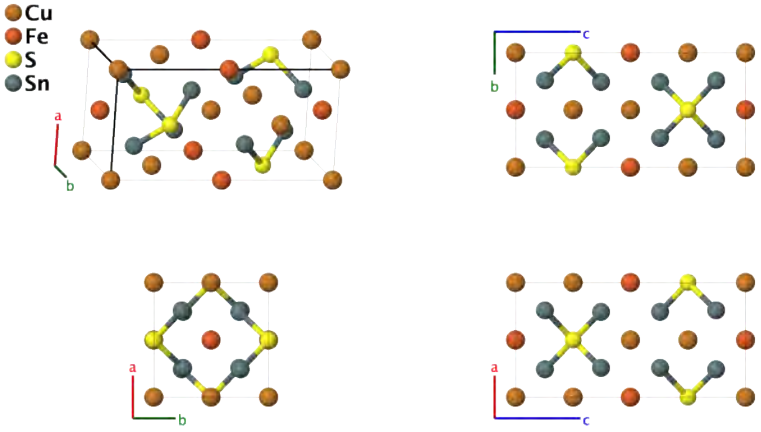
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1667
- POSCAR: pp. 1668

Kesterite [$Cu_2(Zn,Fe)SnS_4$] Structure: A2BCD4_tI16_82_ac_b_d_g

http://aflow.org/prototype-encyclopedia/A2BCD4_tI16_82_ac_b_d_g

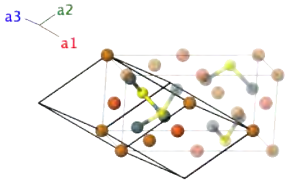


Prototype	:	$Cu_2(Fe,Zn)SnS_4$
AFLOW prototype label	:	A2BCD4_tI16_82_ac_b_d_g
***Strukturbericht* designation**	:	None
Pearson symbol	:	tI16
Space group number	:	82
Space group symbol	:	$I\bar{4}$
AFLOW prototype command	:	`aflow --proto=A2BCD4_tI16_82_ac_b_d_g --params=`$a, c/a, x_5, y_5, z_5$

- The kesterite structure is related to the stannite structure, with differences occurring due to the positioning of the Cu atoms (Hall, 1978). The (2*b*) Wyckoff position is partially occupied with Zn and Fe; however, we decorate it with Fe only.

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}a\,\hat{\mathbf{x}} - \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} - \tfrac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(2*a*)	Cu I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	(2*b*)	Fe
$\mathbf{B}_3$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(2*c*)	Cu II
$\mathbf{B}_4$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(2*d*)	S

$\mathbf{B}_5$	=	$(y_5 + z_5)\ \mathbf{a}_1 + (x_5 + z_5)\ \mathbf{a}_2 + (x_5 + y_5)\ \mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(8*g*)	Sn
$\mathbf{B}_6$	=	$(-y_5 + z_5)\ \mathbf{a}_1 + (-x_5 + z_5)\ \mathbf{a}_2 + (-x_5 - y_5)\ \mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(8*g*)	Sn
$\mathbf{B}_7$	=	$(-x_5 - z_5)\ \mathbf{a}_1 + (y_5 - z_5)\ \mathbf{a}_2 + (-x_5 + y_5)\ \mathbf{a}_3$	=	$y_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(8*g*)	Sn
$\mathbf{B}_8$	=	$(x_5 - z_5)\ \mathbf{a}_1 + (-y_5 - z_5)\ \mathbf{a}_2 + (x_5 - y_5)\ \mathbf{a}_3$	=	$-y_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(8*g*)	Sn

References:

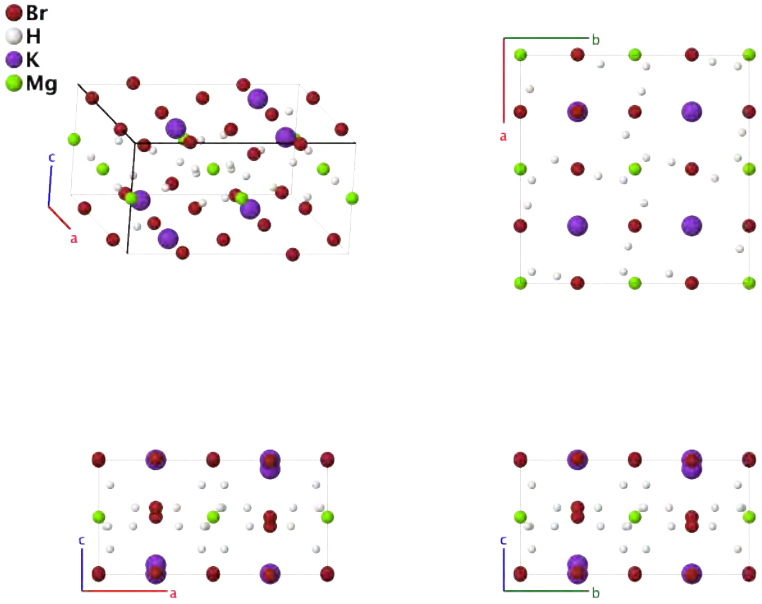
- S. R. Hall, J. T. Szymański, and J. M. Stewart, *Kesterite, $Cu_2(Zn,Fe)SnS_4$, and stannite, $Cu_2(Fe,Zn)SnS_4$, structurally similar but distinct minerals*, Can. Mineral. **16**, 131–137 (1978).
- S. Schorr, H.-J. Hoebler, and M. Tovar, *A neutron diffraction study of the stannite-kesterite solid solution series*, Eur. J. Mineral. **19**, 65–73 (2007), doi:10.1127/0935-1221/2007/0019-0065.
- S. Schorr, *The crystal structure of kesterite type compounds: A neutron and X-ray diffraction study*, Sol. Energy Mater. Sol. Cells **95**, 1482–1488 (2011), doi:10.1016/j.solmat.2011.01.002.

Geometry files:

- CIF: pp. 1668
- POSCAR: pp. 1668

Bromocarnallite ($KMg(H_2O)_6(Cl,Br)_3$, $E2_6$) Structure: A3B6CD_tP44_85_bcg_3g_ac_e

http://aflow.org/prototype-encyclopedia/A3B6CD_tP44_85_bcg_3g_ac_e

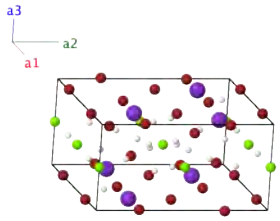


Prototype	:	$(Br,Cl)_3(H_2O)_6KMg$
AFLOW prototype label	:	A3B6CD_tP44_85_bcg_3g_ac_e
***Strukturbericht* designation**	:	$E2_6$
Pearson symbol	:	tP44
Space group number	:	85
Space group symbol	:	$P4/n$
AFLOW prototype command	:	`aflow --proto=A3B6CD_tP44_85_bcg_3g_ac_e --params=`$a, c/a, z_3, z_4, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9$

- (Andreß, 1939) first determined the structure of bromocarnallite, using a sample with 75% bromine on the halide site. They were unable to locate the hydrogen atoms, and placed the compound in space group $P4/n$ #85. The data was given by (Hermann, 1939) in setting 1 of this group, but we used FINDSYM to place it in the standard setting 2.
- (Hermann, 1939) assigned this compound the *Strukturbericht* symbol $E2_6$.
- (Schlemper, 1985) re-examined the structure for the pure chlorine version, carnallite. They located the hydrogen atoms and placed the system in space group *Pnna* #52. We present this structure in the carnallite structure page.

Simple Tetragonal primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= a\,\hat{\mathbf{x}}\\ \mathbf{a}_2 &= a\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2$	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}}$	(2*a*)	K I
$\mathbf{B}_2$	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2$	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}}$	(2*a*)	K I
$\mathbf{B}_3$	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(2*b*)	Br I
$\mathbf{B}_4$	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(2*b*)	Br I
$\mathbf{B}_5$	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	(2*c*)	Br II
$\mathbf{B}_6$	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	(2*c*)	Br II
$\mathbf{B}_7$	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	(2*c*)	K II
$\mathbf{B}_8$	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	(2*c*)	K II
$\mathbf{B}_9$	$\frac{1}{2}\,\mathbf{a}_3$	$\frac{1}{2}c\,\hat{\mathbf{z}}$	(4*e*)	Mg
$\mathbf{B}_{10}$	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*e*)	Mg
$\mathbf{B}_{11}$	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*e*)	Mg
$\mathbf{B}_{12}$	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*e*)	Mg
$\mathbf{B}_{13}$	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	$x_6a\,\hat{\mathbf{x}} + y_6a\,\hat{\mathbf{y}} + z_6c\,\hat{\mathbf{z}}$	(8*g*)	Br III
$\mathbf{B}_{14}$	$\left(\frac{1}{2} - x_6\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	$\left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right)a\,\hat{\mathbf{y}} + z_6c\,\hat{\mathbf{z}}$	(8*g*)	Br III
$\mathbf{B}_{15}$	$\left(\frac{1}{2} - y_6\right)\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	$\left(\frac{1}{2} - y_6\right)a\,\hat{\mathbf{x}} + x_6a\,\hat{\mathbf{y}} + z_6c\,\hat{\mathbf{z}}$	(8*g*)	Br III
$\mathbf{B}_{16}$	$y_6\,\mathbf{a}_1 + \left(\frac{1}{2} - x_6\right)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	$y_6a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{y}} + z_6c\,\hat{\mathbf{z}}$	(8*g*)	Br III
$\mathbf{B}_{17}$	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	$-x_6a\,\hat{\mathbf{x}} - y_6a\,\hat{\mathbf{y}} - z_6c\,\hat{\mathbf{z}}$	(8*g*)	Br III
$\mathbf{B}_{18}$	$\left(\frac{1}{2} + x_6\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right)a\,\hat{\mathbf{y}} - z_6c\,\hat{\mathbf{z}}$	(8*g*)	Br III
$\mathbf{B}_{19}$	$\left(\frac{1}{2} + y_6\right)\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	$\left(\frac{1}{2} + y_6\right)a\,\hat{\mathbf{x}} - x_6a\,\hat{\mathbf{y}} - z_6c\,\hat{\mathbf{z}}$	(8*g*)	Br III
$\mathbf{B}_{20}$	$-y_6\,\mathbf{a}_1 + \left(\frac{1}{2} + x_6\right)\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	$-y_6a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{y}} - z_6c\,\hat{\mathbf{z}}$	(8*g*)	Br III
$\mathbf{B}_{21}$	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	$x_7a\,\hat{\mathbf{x}} + y_7a\,\hat{\mathbf{y}} + z_7c\,\hat{\mathbf{z}}$	(8*g*)	H_2O I
$\mathbf{B}_{22}$	$\left(\frac{1}{2} - x_7\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - y_7\right)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	$\left(\frac{1}{2} - x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_7\right)a\,\hat{\mathbf{y}} + z_7c\,\hat{\mathbf{z}}$	(8*g*)	H_2O I
$\mathbf{B}_{23}$	$\left(\frac{1}{2} - y_7\right)\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	$\left(\frac{1}{2} - y_7\right)a\,\hat{\mathbf{x}} + x_7a\,\hat{\mathbf{y}} + z_7c\,\hat{\mathbf{z}}$	(8*g*)	H_2O I
$\mathbf{B}_{24}$	$y_7\,\mathbf{a}_1 + \left(\frac{1}{2} - x_7\right)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	$y_7a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_7\right)a\,\hat{\mathbf{y}} + z_7c\,\hat{\mathbf{z}}$	(8*g*)	H_2O I
$\mathbf{B}_{25}$	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	$-x_7a\,\hat{\mathbf{x}} - y_7a\,\hat{\mathbf{y}} - z_7c\,\hat{\mathbf{z}}$	(8*g*)	H_2O I
$\mathbf{B}_{26}$	$\left(\frac{1}{2} + x_7\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right)a\,\hat{\mathbf{y}} - z_7c\,\hat{\mathbf{z}}$	(8*g*)	H_2O I
$\mathbf{B}_{27}$	$\left(\frac{1}{2} + y_7\right)\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	$\left(\frac{1}{2} + y_7\right)a\,\hat{\mathbf{x}} - x_7a\,\hat{\mathbf{y}} - z_7c\,\hat{\mathbf{z}}$	(8*g*)	H_2O I

$\mathbf{B}_{28}$	=	$-y_7\,\mathbf{a}_1 + \left(\frac{1}{2} + x_7\right)\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$-y_7 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_7\right) a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	(8*g*)	H_2O I
$\mathbf{B}_{29}$	=	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + y_8 a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(8*g*)	H_2O II
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2} - x_8\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_8\right)\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_8\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_8\right) a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(8*g*)	H_2O II
$\mathbf{B}_{31}$	=	$\left(\frac{1}{2} - y_8\right)\mathbf{a}_1 + x_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_8\right) a\,\hat{\mathbf{x}} + x_8 a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(8*g*)	H_2O II
$\mathbf{B}_{32}$	=	$y_8\,\mathbf{a}_1 + \left(\frac{1}{2} - x_8\right)\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$y_8 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_8\right) a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(8*g*)	H_2O II
$\mathbf{B}_{33}$	=	$-x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} - y_8 a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(8*g*)	H_2O II
$\mathbf{B}_{34}$	=	$\left(\frac{1}{2} + x_8\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_8\right)\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_8\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_8\right) a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(8*g*)	H_2O II
$\mathbf{B}_{35}$	=	$\left(\frac{1}{2} + y_8\right)\mathbf{a}_1 - x_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_8\right) a\,\hat{\mathbf{x}} - x_8 a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(8*g*)	H_2O II
$\mathbf{B}_{36}$	=	$-y_8\,\mathbf{a}_1 + \left(\frac{1}{2} + x_8\right)\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$-y_8 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_8\right) a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(8*g*)	H_2O II
$\mathbf{B}_{37}$	=	$x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}} + y_9 a\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(8*g*)	H_2O III
$\mathbf{B}_{38}$	=	$\left(\frac{1}{2} - x_9\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_9\right)\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_9\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_9\right) a\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(8*g*)	H_2O III
$\mathbf{B}_{39}$	=	$\left(\frac{1}{2} - y_9\right)\mathbf{a}_1 + x_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_9\right) a\,\hat{\mathbf{x}} + x_9 a\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(8*g*)	H_2O III
$\mathbf{B}_{40}$	=	$y_9\,\mathbf{a}_1 + \left(\frac{1}{2} - x_9\right)\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$y_9 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_9\right) a\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(8*g*)	H_2O III
$\mathbf{B}_{41}$	=	$-x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{x}} - y_9 a\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}}$	(8*g*)	H_2O III
$\mathbf{B}_{42}$	=	$\left(\frac{1}{2} + x_9\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_9\right)\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_9\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_9\right) a\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}}$	(8*g*)	H_2O III
$\mathbf{B}_{43}$	=	$\left(\frac{1}{2} + y_9\right)\mathbf{a}_1 - x_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_9\right) a\,\hat{\mathbf{x}} - x_9 a\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}}$	(8*g*)	H_2O III
$\mathbf{B}_{44}$	=	$-y_9\,\mathbf{a}_1 + \left(\frac{1}{2} + x_9\right)\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$-y_9 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_9\right) a\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}}$	(8*g*)	H_2O III

References:

- K. R. Andreß and O. Saffe, *Röntgenographische Untersuchung der Mischkristallreihe Karnallit-Bromkarnallit*, Zeitschrift für Kristallographie - Crystalline Materials **101**, 451–469 (1939), doi:10.1524/zkri.1939.101.1.451.
- E. O. Schlemper, P. K. Sen Gupta, and T. Zoltai, *Refinement of the structure of carnallite, $Mg(H_2O)_6KCl_3$*, Am. Mineral. **70**, 1309–1313 (1985).

Found in:

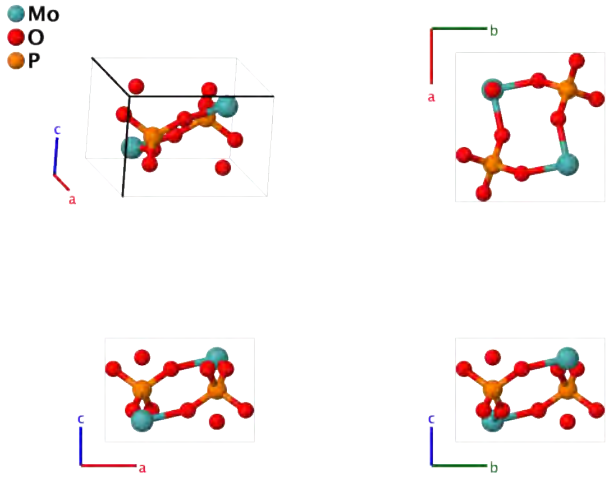
- K. Herrmann, ed., *Strukturbericht Band VII 1939* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1943).

Geometry files:

- CIF: pp. 1669
- POSCAR: pp. 1669

$MoPO_5$ Structure: AB5C_tP14_85_c_cg_b

http://aflow.org/prototype-encyclopedia/AB5C_tP14_85_c_cg_b

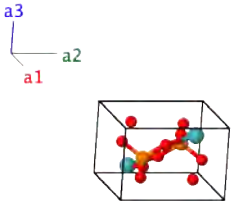


Prototype	:	MoO_5P
AFLOW prototype label	:	AB5C_tP14_85_c_cg_b
***Strukturbericht* designation**	:	None
Pearson symbol	:	tP14
Space group number	:	85
Space group symbol	:	$P4/n$
AFLOW prototype command	:	`aflow --proto=AB5C_tP14_85_c_cg_b --params=`$a, c/a, z_2, z_3, x_4, y_4, z_4$

- (Kierkegaard, 1964) gave this structure in setting 1 of space group $P4/n$ #85. We used FINDSYM to shift this to the standard setting 2.

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(2b)$	P
$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(2b)$	P
$\mathbf{B}_3$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + z_2c\,\hat{\mathbf{z}}$	$(2c)$	Mo

$\mathbf{B}_4$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(2*c*)	Mo
$\mathbf{B}_5$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(2*c*)	O I
$\mathbf{B}_6$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(2*c*)	O I
$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*g*)	O II
$\mathbf{B}_8$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*g*)	O II
$\mathbf{B}_9$	=	$\left(\frac{1}{2} - y_4\right)\mathbf{a}_1 + x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_4\right)a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*g*)	O II
$\mathbf{B}_{10}$	=	$y_4\,\mathbf{a}_1 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$y_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*g*)	O II
$\mathbf{B}_{11}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 a\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*g*)	O II
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)a\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*g*)	O II
$\mathbf{B}_{13}$	=	$\left(\frac{1}{2} + y_4\right)\mathbf{a}_1 - x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_4\right)a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*g*)	O II
$\mathbf{B}_{14}$	=	$-y_4\,\mathbf{a}_1 + \left(\frac{1}{2} + x_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-y_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*g*)	O II

References:

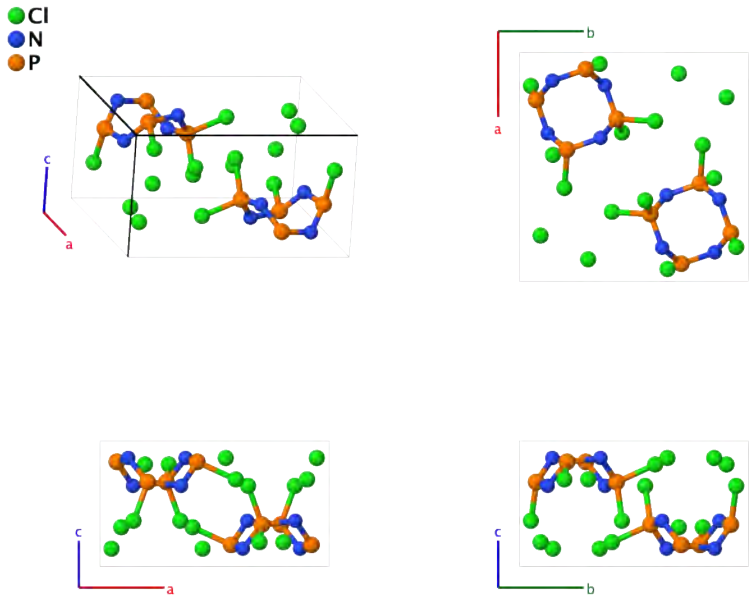
- P. Kierkegaard and M. Westerlund, *The Crystal Structure of $MoOPO_4$*, Acta Chem. Scand. **18**, 2217–2225 (1964), doi:10.3891/acta.chem.scand.18-2217.

Geometry files:

- CIF: pp. 1669
- POSCAR: pp. 1670

$PNCl_2$ ($E1_4$) Structure: A2BC_tP32_86_2g_g_g

http://aflow.org/prototype-encyclopedia/A2BC_tP32_86_2g_g_g

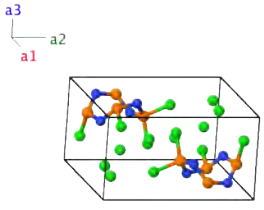


Prototype	:	Cl_2NP
AFLOW prototype label	:	A2BC_tP32_86_2g_g_g
***Strukturbericht* designation**	:	$E1_4$
Pearson symbol	:	tP32
Space group number	:	86
Space group symbol	:	$P4_2/n$
AFLOW prototype command	:	`aflow --proto=A2BC_tP32_86_2g_g_g --params=`$a, c/a, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4$

- (Ketellar, 1939) gave the atomic positions in setting 1 of space group #86. We have shifted the origin to the inversion site for this lattice and present the atomic positions in setting 2.

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + y_1 a\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(8*g*)	Cl I
$\mathbf{B}_2$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_1\right)\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_1\right)a\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(8*g*)	Cl I
$\mathbf{B}_3$	=	$-y_1\,\mathbf{a}_1 + \left(\frac{1}{2} + x_1\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$-y_1 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_1\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\,\hat{\mathbf{z}}$	(8*g*)	Cl I
$\mathbf{B}_4$	=	$\left(\frac{1}{2} + y_1\right)\mathbf{a}_1 - x_1\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_1\right)a\,\hat{\mathbf{x}} - x_1 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\,\hat{\mathbf{z}}$	(8*g*)	Cl I
$\mathbf{B}_5$	=	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} - y_1 a\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	(8*g*)	Cl I
$\mathbf{B}_6$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_1\right)\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_1\right)a\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	(8*g*)	Cl I
$\mathbf{B}_7$	=	$y_1\,\mathbf{a}_1 + \left(\frac{1}{2} - x_1\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	=	$y_1 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_1\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right)c\,\hat{\mathbf{z}}$	(8*g*)	Cl I
$\mathbf{B}_8$	=	$\left(\frac{1}{2} - y_1\right)\mathbf{a}_1 + x_1\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_1\right)a\,\hat{\mathbf{x}} + x_1 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right)c\,\hat{\mathbf{z}}$	(8*g*)	Cl I
$\mathbf{B}_9$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + y_2 a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(8*g*)	Cl II
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right)a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(8*g*)	Cl II
$\mathbf{B}_{11}$	=	$-y_2\,\mathbf{a}_1 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$-y_2 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	(8*g*)	Cl II
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + y_2\right)\mathbf{a}_1 - x_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_2\right)a\,\hat{\mathbf{x}} - x_2 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	(8*g*)	Cl II
$\mathbf{B}_{13}$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} - y_2 a\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(8*g*)	Cl II
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right)a\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(8*g*)	Cl II
$\mathbf{B}_{15}$	=	$y_2\,\mathbf{a}_1 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$y_2 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	(8*g*)	Cl II
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} - y_2\right)\mathbf{a}_1 + x_2\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_2\right)a\,\hat{\mathbf{x}} + x_2 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	(8*g*)	Cl II
$\mathbf{B}_{17}$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(8*g*)	N
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(8*g*)	N
$\mathbf{B}_{19}$	=	$-y_3\,\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$-y_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(8*g*)	N
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} + y_3\right)\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(8*g*)	N
$\mathbf{B}_{21}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(8*g*)	N
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(8*g*)	N
$\mathbf{B}_{23}$	=	$y_3\,\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$y_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(8*g*)	N
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} - y_3\right)\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(8*g*)	N
$\mathbf{B}_{25}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*g*)	P
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*g*)	P
$\mathbf{B}_{27}$	=	$-y_4\,\mathbf{a}_1 + \left(\frac{1}{2} + x_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$-y_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(8*g*)	P
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} + y_4\right)\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_4\right)a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(8*g*)	P
$\mathbf{B}_{29}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 a\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*g*)	P
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)a\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*g*)	P
$\mathbf{B}_{31}$	=	$y_4\,\mathbf{a}_1 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$y_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(8*g*)	P
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2} - y_4\right)\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_4\right)a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(8*g*)	P

References:

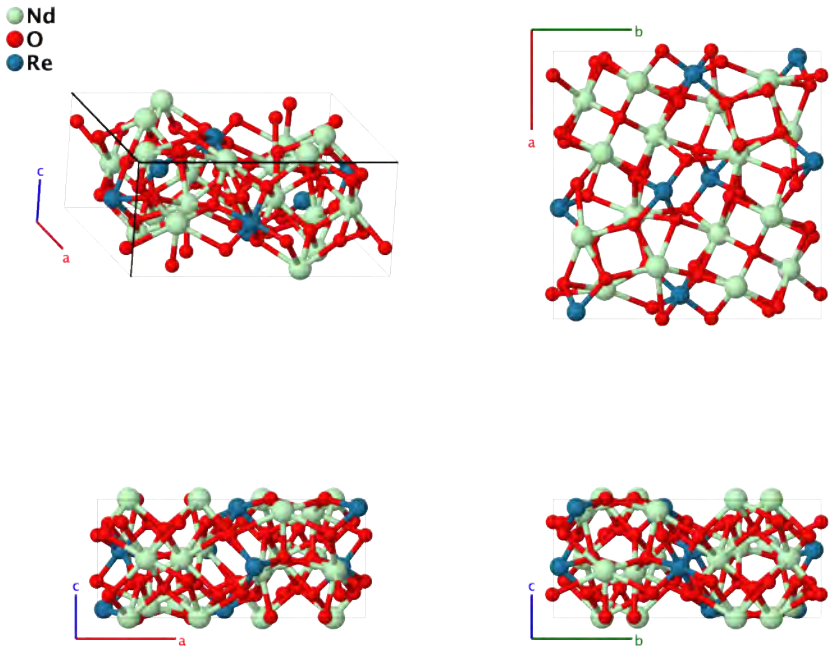
- J. A. A. Ketelaar and T. A. de Vries, *The crystal structure of tetra phosphonitrile chloride, $P_4N_4Cl_8$*, Rec. Trav. Chim. Pays-Bas **58**, 1081–1099 (1939), doi:10.1002/recl.19390581205.

Geometry files:

- CIF: pp. 1670
- POSCAR: pp. 1670

$Nd_4Re_2O_{11}$ Structure: A4B11C2_tP68_86_2g_ab5g_g

http://aflow.org/prototype-encyclopedia/A4B11C2_tP68_86_2g_ab5g_g

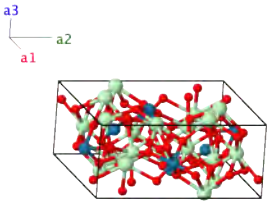


Prototype	:	$Nd_4O_{11}Re_2$
AFLOW prototype label	:	A4B11C2_tP68_86_2g_ab5g_g
***Strukturbericht* designation**	:	None
Pearson symbol	:	tP68
Space group number	:	86
Space group symbol	:	$P4_2/n$
AFLOW prototype command	:	`aflow --proto=A4B11C2_tP68_86_2g_ab5g_g --params=`$a, c/a, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}$

- (Wilhelmi, 1970) has a misprint for the Wyckoff position of the Nd-II atom, although the nearest-neighbor distances are correct. (Downs, 2003) corrects the position to be consistent with those distances, and we use their value.

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(2*a*)	O I
$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(2*a*)	O I
$\mathbf{B}_3$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(2*b*)	O II
$\mathbf{B}_4$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(2*b*)	O II
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + y_3a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	(8*g*)	Nd I
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	(8*g*)	Nd I
$\mathbf{B}_7$	=	$-y_3\,\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$-y_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(8*g*)	Nd I
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + y_3\right)\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(8*g*)	Nd I
$\mathbf{B}_9$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - y_3a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	(8*g*)	Nd I
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	(8*g*)	Nd I
$\mathbf{B}_{11}$	=	$y_3\,\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$y_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(8*g*)	Nd I
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} - y_3\right)\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(8*g*)	Nd I
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + y_4a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	(8*g*)	Nd II
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	(8*g*)	Nd II
$\mathbf{B}_{15}$	=	$-y_4\,\mathbf{a}_1 + \left(\frac{1}{2} + x_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$-y_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(8*g*)	Nd II
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + y_4\right)\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_4\right)a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(8*g*)	Nd II
$\mathbf{B}_{17}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} - y_4a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	(8*g*)	Nd II
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	(8*g*)	Nd II
$\mathbf{B}_{19}$	=	$y_4\,\mathbf{a}_1 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$y_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(8*g*)	Nd II
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} - y_4\right)\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_4\right)a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(8*g*)	Nd II
$\mathbf{B}_{21}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}} + y_5a\,\hat{\mathbf{y}} + z_5c\,\hat{\mathbf{z}}$	(8*g*)	O III
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right)a\,\hat{\mathbf{y}} + z_5c\,\hat{\mathbf{z}}$	(8*g*)	O III
$\mathbf{B}_{23}$	=	$-y_5\,\mathbf{a}_1 + \left(\frac{1}{2} + x_5\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$-y_5a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	(8*g*)	O III
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} + y_5\right)\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_5\right)a\,\hat{\mathbf{x}} - x_5a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	(8*g*)	O III
$\mathbf{B}_{25}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}} - y_5a\,\hat{\mathbf{y}} - z_5c\,\hat{\mathbf{z}}$	(8*g*)	O III
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right)a\,\hat{\mathbf{y}} - z_5c\,\hat{\mathbf{z}}$	(8*g*)	O III
$\mathbf{B}_{27}$	=	$y_5\,\mathbf{a}_1 + \left(\frac{1}{2} - x_5\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$y_5a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	(8*g*)	O III
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} - y_5\right)\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_5\right)a\,\hat{\mathbf{x}} + x_5a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	(8*g*)	O III
$\mathbf{B}_{29}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6a\,\hat{\mathbf{x}} + y_6a\,\hat{\mathbf{y}} + z_6c\,\hat{\mathbf{z}}$	(8*g*)	O IV
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right)a\,\hat{\mathbf{y}} + z_6c\,\hat{\mathbf{z}}$	(8*g*)	O IV
$\mathbf{B}_{31}$	=	$-y_6\,\mathbf{a}_1 + \left(\frac{1}{2} + x_6\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$-y_6a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\,\hat{\mathbf{z}}$	(8*g*)	O IV
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2} + y_6\right)\mathbf{a}_1 - x_6\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_6\right)a\,\hat{\mathbf{x}} - x_6a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\,\hat{\mathbf{z}}$	(8*g*)	O IV
$\mathbf{B}_{33}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$-x_6a\,\hat{\mathbf{x}} - y_6a\,\hat{\mathbf{y}} - z_6c\,\hat{\mathbf{z}}$	(8*g*)	O IV
$\mathbf{B}_{34}$	=	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right)a\,\hat{\mathbf{y}} - z_6c\,\hat{\mathbf{z}}$	(8*g*)	O IV

$$
\begin{array}{lcccccc}
\mathbf{B}_{35} & = & y_6\,\mathbf{a}_1 + \left(\frac{1}{2} - x_6\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3 & = & y_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_6\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right) c\,\hat{\mathbf{z}} & (8g) & \text{O IV} \\
\mathbf{B}_{36} & = & \left(\frac{1}{2} - y_6\right)\mathbf{a}_1 + x_6\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - y_6\right) a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right) c\,\hat{\mathbf{z}} & (8g) & \text{O IV} \\
\mathbf{B}_{37} & = & x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 & = & x_7 a\,\hat{\mathbf{x}} + y_7 a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}} & (8g) & \text{O V} \\
\mathbf{B}_{38} & = & \left(\frac{1}{2} - x_7\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_7\right)\mathbf{a}_2 + z_7\,\mathbf{a}_3 & = & \left(\frac{1}{2} - x_7\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_7\right) a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}} & (8g) & \text{O V} \\
\mathbf{B}_{39} & = & -y_7\,\mathbf{a}_1 + \left(\frac{1}{2} + x_7\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3 & = & -y_7 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_7\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right) c\,\hat{\mathbf{z}} & (8g) & \text{O V} \\
\mathbf{B}_{40} & = & \left(\frac{1}{2} + y_7\right)\mathbf{a}_1 - x_7\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3 & = & \left(\frac{1}{2} + y_7\right) a\,\hat{\mathbf{x}} - x_7 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right) c\,\hat{\mathbf{z}} & (8g) & \text{O V} \\
\mathbf{B}_{41} & = & -x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3 & = & -x_7 a\,\hat{\mathbf{x}} - y_7 a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}} & (8g) & \text{O V} \\
\mathbf{B}_{42} & = & \left(\frac{1}{2} + x_7\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\mathbf{a}_2 - z_7\,\mathbf{a}_3 & = & \left(\frac{1}{2} + x_7\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right) a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}} & (8g) & \text{O V} \\
\mathbf{B}_{43} & = & y_7\,\mathbf{a}_1 + \left(\frac{1}{2} - x_7\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\mathbf{a}_3 & = & y_7 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_7\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right) c\,\hat{\mathbf{z}} & (8g) & \text{O V} \\
\mathbf{B}_{44} & = & \left(\frac{1}{2} - y_7\right)\mathbf{a}_1 + x_7\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - y_7\right) a\,\hat{\mathbf{x}} + x_7 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right) c\,\hat{\mathbf{z}} & (8g) & \text{O V} \\
\mathbf{B}_{45} & = & x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 & = & x_8 a\,\hat{\mathbf{x}} + y_8 a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}} & (8g) & \text{O VI} \\
\mathbf{B}_{46} & = & \left(\frac{1}{2} - x_8\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_8\right)\mathbf{a}_2 + z_8\,\mathbf{a}_3 & = & \left(\frac{1}{2} - x_8\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_8\right) a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}} & (8g) & \text{O VI} \\
\mathbf{B}_{47} & = & -y_8\,\mathbf{a}_1 + \left(\frac{1}{2} + x_8\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3 & = & -y_8 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_8\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right) c\,\hat{\mathbf{z}} & (8g) & \text{O VI} \\
\mathbf{B}_{48} & = & \left(\frac{1}{2} + y_8\right)\mathbf{a}_1 - x_8\,\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3 & = & \left(\frac{1}{2} + y_8\right) a\,\hat{\mathbf{x}} - x_8 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right) c\,\hat{\mathbf{z}} & (8g) & \text{O VI} \\
\mathbf{B}_{49} & = & -x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3 & = & -x_8 a\,\hat{\mathbf{x}} - y_8 a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}} & (8g) & \text{O VI} \\
\mathbf{B}_{50} & = & \left(\frac{1}{2} + x_8\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_8\right)\mathbf{a}_2 - z_8\,\mathbf{a}_3 & = & \left(\frac{1}{2} + x_8\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_8\right) a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}} & (8g) & \text{O VI} \\
\mathbf{B}_{51} & = & y_8\,\mathbf{a}_1 + \left(\frac{1}{2} - x_8\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\mathbf{a}_3 & = & y_8 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_8\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right) c\,\hat{\mathbf{z}} & (8g) & \text{O VI} \\
\mathbf{B}_{52} & = & \left(\frac{1}{2} - y_8\right)\mathbf{a}_1 + x_8\,\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - y_8\right) a\,\hat{\mathbf{x}} + x_8 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right) c\,\hat{\mathbf{z}} & (8g) & \text{O VI} \\
\mathbf{B}_{53} & = & x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 & = & x_9 a\,\hat{\mathbf{x}} + y_9 a\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}} & (8g) & \text{O VII} \\
\mathbf{B}_{54} & = & \left(\frac{1}{2} - x_9\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_9\right)\mathbf{a}_2 + z_9\,\mathbf{a}_3 & = & \left(\frac{1}{2} - x_9\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_9\right) a\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}} & (8g) & \text{O VII} \\
\mathbf{B}_{55} & = & -y_9\,\mathbf{a}_1 + \left(\frac{1}{2} + x_9\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3 & = & -y_9 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_9\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right) c\,\hat{\mathbf{z}} & (8g) & \text{O VII} \\
\mathbf{B}_{56} & = & \left(\frac{1}{2} + y_9\right)\mathbf{a}_1 - x_9\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3 & = & \left(\frac{1}{2} + y_9\right) a\,\hat{\mathbf{x}} - x_9 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right) c\,\hat{\mathbf{z}} & (8g) & \text{O VII} \\
\mathbf{B}_{57} & = & -x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3 & = & -x_9 a\,\hat{\mathbf{x}} - y_9 a\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}} & (8g) & \text{O VII} \\
\mathbf{B}_{58} & = & \left(\frac{1}{2} + x_9\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_9\right)\mathbf{a}_2 - z_9\,\mathbf{a}_3 & = & \left(\frac{1}{2} + x_9\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_9\right) a\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}} & (8g) & \text{O VII} \\
\mathbf{B}_{59} & = & y_9\,\mathbf{a}_1 + \left(\frac{1}{2} - x_9\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_9\right)\mathbf{a}_3 & = & y_9 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_9\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_9\right) c\,\hat{\mathbf{z}} & (8g) & \text{O VII} \\
\mathbf{B}_{60} & = & \left(\frac{1}{2} - y_9\right)\mathbf{a}_1 + x_9\,\mathbf{a}_2 + \left(\frac{1}{2} - z_9\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - y_9\right) a\,\hat{\mathbf{x}} + x_9 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_9\right) c\,\hat{\mathbf{z}} & (8g) & \text{O VII} \\
\mathbf{B}_{61} & = & x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 & = & x_{10} a\,\hat{\mathbf{x}} + y_{10} a\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}} & (8g) & \text{Re} \\
\mathbf{B}_{62} & = & \left(\frac{1}{2} - x_{10}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{10}\right)\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 & = & \left(\frac{1}{2} - x_{10}\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{10}\right) a\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}} & (8g) & \text{Re} \\
\mathbf{B}_{63} & = & -y_{10}\,\mathbf{a}_1 + \left(\frac{1}{2} + x_{10}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{10}\right)\mathbf{a}_3 & = & -y_{10} a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_{10}\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{10}\right) c\,\hat{\mathbf{z}} & (8g) & \text{Re} \\
\mathbf{B}_{64} & = & \left(\frac{1}{2} + y_{10}\right)\mathbf{a}_1 - x_{10}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{10}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} + y_{10}\right) a\,\hat{\mathbf{x}} - x_{10} a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{10}\right) c\,\hat{\mathbf{z}} & (8g) & \text{Re} \\
\mathbf{B}_{65} & = & -x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 & = & -x_{10} a\,\hat{\mathbf{x}} - y_{10} a\,\hat{\mathbf{y}} - z_{10} c\,\hat{\mathbf{z}} & (8g) & \text{Re} \\
\mathbf{B}_{66} & = & \left(\frac{1}{2} + x_{10}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{10}\right)\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 & = & \left(\frac{1}{2} + x_{10}\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{10}\right) a\,\hat{\mathbf{y}} - z_{10} c\,\hat{\mathbf{z}} & (8g) & \text{Re} \\
\mathbf{B}_{67} & = & y_{10}\,\mathbf{a}_1 + \left(\frac{1}{2} - x_{10}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{10}\right)\mathbf{a}_3 & = & y_{10} a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_{10}\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{10}\right) c\,\hat{\mathbf{z}} & (8g) & \text{Re} \\
\mathbf{B}_{68} & = & \left(\frac{1}{2} - y_{10}\right)\mathbf{a}_1 + x_{10}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{10}\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - y_{10}\right) a\,\hat{\mathbf{x}} + x_{10} a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{10}\right) c\,\hat{\mathbf{z}} & (8g) & \text{Re}
\end{array}
$$

References:

- K.-A. Wilhelmi, E. Lagervall, and O. Muller, *On the Crystal Structure of* $Nd_4Re_2O_{11}$, Acta Chem. Scand. **24**, 3406–3408 (1970), doi:10.3891/acta.chem.scand.24-3406.

Found in:

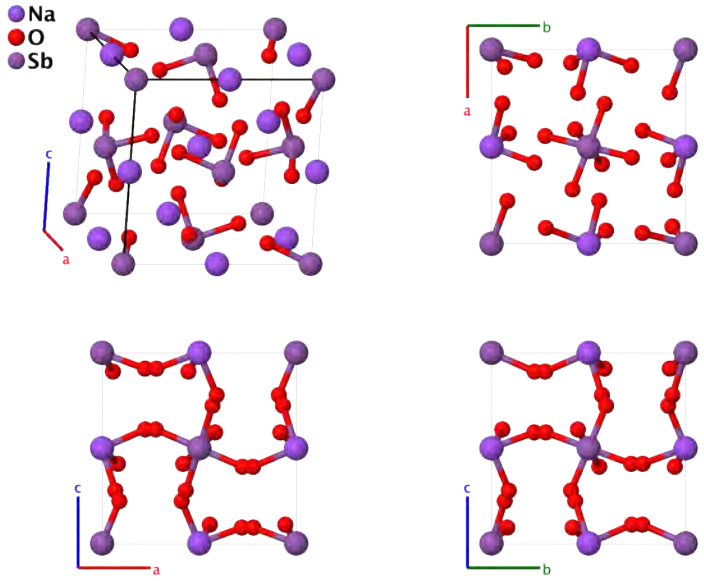
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1670
- POSCAR: pp. 1671

$NaSb(OH)_6$ ($J1_{11}$) Structure: AB6C_tP32_86_d_3g_c

http://aflow.org/prototype-encyclopedia/AB6C_tP32_86_d_3g_c



Prototype	:	$Na(OH)_6Sb$
AFLOW prototype label	:	AB6C_tP32_86_d_3g_c
***Strukturbericht* designation**	:	$J1_{11}$
Pearson symbol	:	tP32
Space group number	:	86
Space group symbol	:	$P4_2/n$
AFLOW prototype command	:	`aflow --proto=AB6C_tP32_86_d_3g_c --params=`$a, c/a, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5$

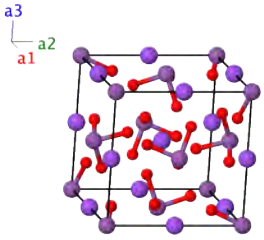
Other compounds with this structure

- $AgSb(OH)_6$

- The atomic positions were originally determined using setting 1 of space group #86. We used FINDSYM to change the origin to setting 2.
- The positions of the hydrogen atoms in the OH ion were not determined, so we only provide the positions of the oxygen atoms (labeled as OH).
- Although the replacement of fluorine by OH does not affect the shape of the Sb-$(F,OH)_6$ ions, it has a profound effect on the structure, as can be seen by looking at $NaSbF_6$ and $NaSbF_4(OH)_2$ ($J1_{12}$).

Simple Tetragonal primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= a\,\hat{\mathbf{x}}\\ \mathbf{a}_2 &= a\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(4*c*)	Sb
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$	(4*c*)	Sb
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*c*)	Sb
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*c*)	Sb
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	(4*d*)	Na
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*d*)	Na
$\mathbf{B}_7$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{y}}$	(4*d*)	Na
$\mathbf{B}_8$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	(4*d*)	Na
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + y_3a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	(8*g*)	O I
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	(8*g*)	O I
$\mathbf{B}_{11}$	=	$-y_3\,\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$-y_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(8*g*)	O I
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + y_3\right)\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(8*g*)	O I
$\mathbf{B}_{13}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - y_3a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	(8*g*)	O I
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	(8*g*)	O I
$\mathbf{B}_{15}$	=	$y_3\,\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$y_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(8*g*)	O I
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} - y_3\right)\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(8*g*)	O I
$\mathbf{B}_{17}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + y_4a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	(8*g*)	O II
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	(8*g*)	O II
$\mathbf{B}_{19}$	=	$-y_4\,\mathbf{a}_1 + \left(\frac{1}{2} + x_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$-y_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(8*g*)	O II
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} + y_4\right)\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_4\right)a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(8*g*)	O II
$\mathbf{B}_{21}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} - y_4a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	(8*g*)	O II
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	(8*g*)	O II
$\mathbf{B}_{23}$	=	$y_4\,\mathbf{a}_1 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$y_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(8*g*)	O II
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} - y_4\right)\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_4\right)a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(8*g*)	O II
$\mathbf{B}_{25}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}} + y_5a\,\hat{\mathbf{y}} + z_5c\,\hat{\mathbf{z}}$	(8*g*)	O III
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right)a\,\hat{\mathbf{y}} + z_5c\,\hat{\mathbf{z}}$	(8*g*)	O III
$\mathbf{B}_{27}$	=	$-y_5\,\mathbf{a}_1 + \left(\frac{1}{2} + x_5\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$-y_5a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	(8*g*)	O III

$$\mathbf{B}_{28} = \left(\tfrac{1}{2}+y_5\right)\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+y_5\right)a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_5\right)c\,\hat{\mathbf{z}} \quad (8g) \quad \text{O III}$$

$$\mathbf{B}_{29} = -x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3 = -x_5 a\,\hat{\mathbf{x}} - y_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}} \quad (8g) \quad \text{O III}$$

$$\mathbf{B}_{30} = \left(\tfrac{1}{2}+x_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3 = \left(\tfrac{1}{2}+x_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_5\right)a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}} \quad (8g) \quad \text{O III}$$

$$\mathbf{B}_{31} = y_5\,\mathbf{a}_1 + \left(\tfrac{1}{2}-x_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_5\right)\mathbf{a}_3 = y_5 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_5\right)c\,\hat{\mathbf{z}} \quad (8g) \quad \text{O III}$$

$$\mathbf{B}_{32} = \left(\tfrac{1}{2}-y_5\right)\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-y_5\right)a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_5\right)c\,\hat{\mathbf{z}} \quad (8g) \quad \text{O III}$$

References:

- T. Asai, *Refinement of the Crystal Structure of Sodium Hexahydroxoantimonate(V), $NaSb(OH)_6$*, Bull. Chem. Soc. Jpn. **48**, 2677–2679 (1975), doi:10.1246/bcsj.48.2677.

Found in:

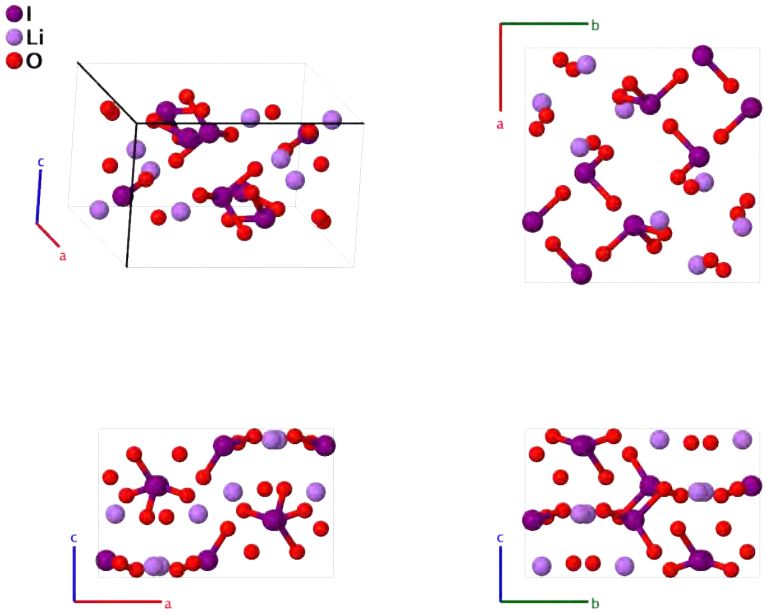
- F. Hoffmann, M. Sartor, and M. Fröba, *The Fascination of Crystals and Symmetry*, http://crystalsymmetry.wordpress.com/tag/nasboh6/ (2014). $NaSb(OH)_6$.

Geometry files:

- CIF: pp. 1671
- POSCAR: pp. 1671

β-$LiIO_3$ Structure: ABC3_tP40_86_g_g_3g

http://aflow.org/prototype-encyclopedia/ABC3_tP40_86_g_g_3g

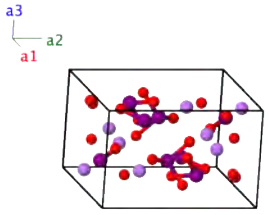


Prototype	:	$ILiO_3$
AFLOW prototype label	:	ABC3_tP40_86_g_g_3g
***Strukturbericht* designation**	:	None
Pearson symbol	:	tP40
Space group number	:	86
Space group symbol	:	$P4_2/n$
AFLOW prototype command	:	`aflow --proto=ABC3_tP40_86_g_g_3g --params=`$a, c/a, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5$

- $LiIO_3$ is known to exist in three forms:
  - α-$LiIO_3$, stable below 473 K: (Zachariasen, 1931) originally determined that the structure of α-$LiIO_3$ was in space group $P6_322$ #182, which (Hermann, 1937) designated *Strukturbericht* $E2_3$. (Rosenzweig, 1966) subsequently determined that the true structure was in space group $P6_3$ #173.
  - β-$LiIO_3$, stable from 573 K up to the melting point at 708 K (this structure).
  - γ-$LiIO_3$, stable between the α- and β-phases, with an orthorhombic structure in space group $Pna2_1$ #33.

Simple Tetragonal primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= a\,\hat{\mathbf{x}}\\ \mathbf{a}_2 &= a\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	$x_1 a\,\hat{\mathbf{x}} + y_1 a\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	$(8g)$	I
$\mathbf{B}_2$	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_1\right)\mathbf{a}_2 + z_1\,\mathbf{a}_3$	$\left(\frac{1}{2} - x_1\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_1\right)a\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	$(8g)$	I
$\mathbf{B}_3$	$-y_1\,\mathbf{a}_1 + \left(\frac{1}{2} + x_1\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	$-y_1 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_1\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\,\hat{\mathbf{z}}$	$(8g)$	I
$\mathbf{B}_4$	$\left(\frac{1}{2} + y_1\right)\mathbf{a}_1 - x_1\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	$\left(\frac{1}{2} + y_1\right)a\,\hat{\mathbf{x}} - x_1 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\,\hat{\mathbf{z}}$	$(8g)$	I
$\mathbf{B}_5$	$-x_1\,\mathbf{a}_1 - y_1\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	$-x_1 a\,\hat{\mathbf{x}} - y_1 a\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	$(8g)$	I
$\mathbf{B}_6$	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_1\right)\mathbf{a}_2 - z_1\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_1\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_1\right)a\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	$(8g)$	I
$\mathbf{B}_7$	$y_1\,\mathbf{a}_1 + \left(\frac{1}{2} - x_1\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	$y_1 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_1\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right)c\,\hat{\mathbf{z}}$	$(8g)$	I
$\mathbf{B}_8$	$\left(\frac{1}{2} - y_1\right)\mathbf{a}_1 + x_1\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	$\left(\frac{1}{2} - y_1\right)a\,\hat{\mathbf{x}} + x_1 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right)c\,\hat{\mathbf{z}}$	$(8g)$	I
$\mathbf{B}_9$	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	$x_2 a\,\hat{\mathbf{x}} + y_2 a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(8g)$	Li
$\mathbf{B}_{10}$	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 + z_2\,\mathbf{a}_3$	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right)a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(8g)$	Li
$\mathbf{B}_{11}$	$-y_2\,\mathbf{a}_1 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	$-y_2 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	$(8g)$	Li
$\mathbf{B}_{12}$	$\left(\frac{1}{2} + y_2\right)\mathbf{a}_1 - x_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	$\left(\frac{1}{2} + y_2\right)a\,\hat{\mathbf{x}} - x_2 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	$(8g)$	Li
$\mathbf{B}_{13}$	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	$-x_2 a\,\hat{\mathbf{x}} - y_2 a\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	$(8g)$	Li
$\mathbf{B}_{14}$	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 - z_2\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right)a\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	$(8g)$	Li
$\mathbf{B}_{15}$	$y_2\,\mathbf{a}_1 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	$y_2 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	$(8g)$	Li
$\mathbf{B}_{16}$	$\left(\frac{1}{2} - y_2\right)\mathbf{a}_1 + x_2\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	$\left(\frac{1}{2} - y_2\right)a\,\hat{\mathbf{x}} + x_2 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	$(8g)$	Li
$\mathbf{B}_{17}$	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$x_3 a\,\hat{\mathbf{x}} + y_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(8g)$	O I
$\mathbf{B}_{18}$	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(8g)$	O I
$\mathbf{B}_{19}$	$-y_3\,\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	$-y_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	$(8g)$	O I
$\mathbf{B}_{20}$	$\left(\frac{1}{2} + y_3\right)\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	$\left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	$(8g)$	O I
$\mathbf{B}_{21}$	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	$-x_3 a\,\hat{\mathbf{x}} - y_3 a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(8g)$	O I
$\mathbf{B}_{22}$	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 - z_3\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(8g)$	O I
$\mathbf{B}_{23}$	$y_3\,\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	$y_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	$(8g)$	O I
$\mathbf{B}_{24}$	$\left(\frac{1}{2} - y_3\right)\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	$\left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	$(8g)$	O I
$\mathbf{B}_{25}$	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$x_4 a\,\hat{\mathbf{x}} + y_4 a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(8g)$	O II
$\mathbf{B}_{26}$	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(8g)$	O II
$\mathbf{B}_{27}$	$-y_4\,\mathbf{a}_1 + \left(\frac{1}{2} + x_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	$-y_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	$(8g)$	O II

$\mathbf{B}_{28}$	=	$\left(\frac{1}{2}+y_4\right)\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_4\right)a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_4\right)c\,\hat{\mathbf{z}}$	(8*g*)	O II
$\mathbf{B}_{29}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 a\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*g*)	O II
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2}+x_4\right)\mathbf{a}_1 + \left(\frac{1}{2}+y_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_4\right)a\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*g*)	O II
$\mathbf{B}_{31}$	=	$y_4\,\mathbf{a}_1 + \left(\frac{1}{2}-x_4\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_4\right)\mathbf{a}_3$	=	$y_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_4\right)c\,\hat{\mathbf{z}}$	(8*g*)	O II
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2}-y_4\right)\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \left(\frac{1}{2}-z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_4\right)a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_4\right)c\,\hat{\mathbf{z}}$	(8*g*)	O II
$\mathbf{B}_{33}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(8*g*)	O III
$\mathbf{B}_{34}$	=	$\left(\frac{1}{2}-x_5\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_5\right)\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_5\right)a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(8*g*)	O III
$\mathbf{B}_{35}$	=	$-y_5\,\mathbf{a}_1 + \left(\frac{1}{2}+x_5\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$-y_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	(8*g*)	O III
$\mathbf{B}_{36}$	=	$\left(\frac{1}{2}+y_5\right)\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_5\right)a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	(8*g*)	O III
$\mathbf{B}_{37}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(8*g*)	O III
$\mathbf{B}_{38}$	=	$\left(\frac{1}{2}+x_5\right)\mathbf{a}_1 + \left(\frac{1}{2}+y_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_5\right)a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(8*g*)	O III
$\mathbf{B}_{39}$	=	$y_5\,\mathbf{a}_1 + \left(\frac{1}{2}-x_5\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_5\right)\mathbf{a}_3$	=	$y_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_5\right)c\,\hat{\mathbf{z}}$	(8*g*)	O III
$\mathbf{B}_{40}$	=	$\left(\frac{1}{2}-y_5\right)\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \left(\frac{1}{2}-z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_5\right)a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_5\right)c\,\hat{\mathbf{z}}$	(8*g*)	O III

References:

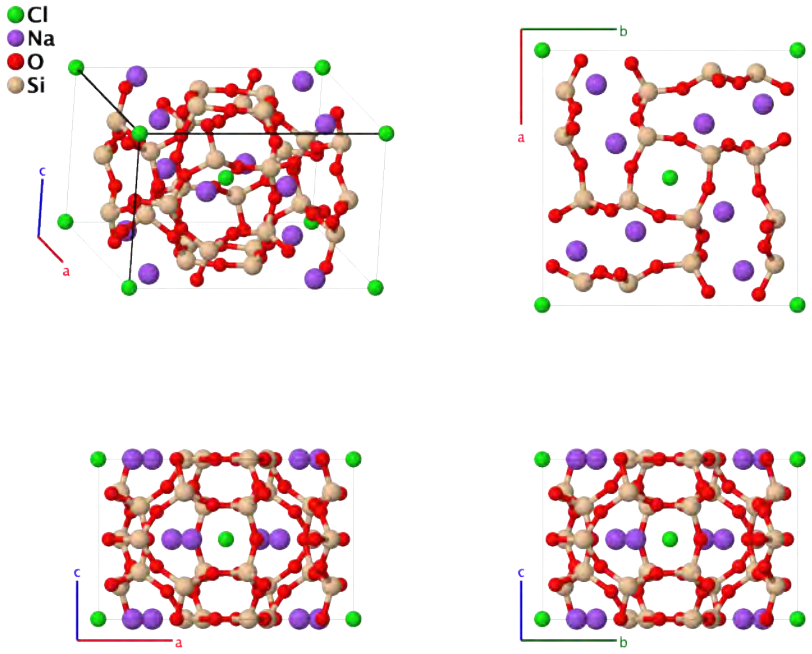
- H. Schulz, *The structure of β-$LiIO_3$*, Acta Crystallogr. Sect. B Struct. Sci. **29**, 2285–2289 (1973), doi:10.1107/S0567740873006485.
- W. H. Zachariasen and F. A. Barta, *Crystal Structure of Lithium Iodate*, Phys. Rev. **37**, 1626–1630 (1931), doi:10.1103/PhysRev.37.1626.
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).
- A. Rosenzweig and B. Morosin, *A reinvestigation of the crystal structure of $LiIO_3$*, Acta Cryst. **20**, 758–761 (1966), doi:10.1107/S0365110X66001804.

Geometry files:

- CIF: pp. 1672
- POSCAR: pp. 1672

Marialite Scapolite [$Na_4Cl(AlSi_3)_3O_{24}$, $S6_4$] Structure: AB4C24D12_tI82_87_a_h_2h2i_hi

http://aflow.org/prototype-encyclopedia/AB4C24D12_tI82_87_a_h_2h2i_hi

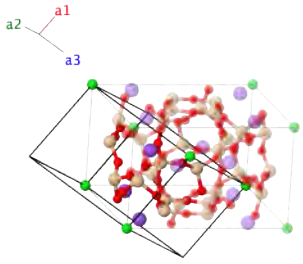


Prototype	:	$ClNa_4O_{24}(Al_3Si_9)$
AFLOW prototype label	:	AB4C24D12_tI82_87_a_h_2h2i_hi
***Strukturbericht* designation**	:	$S6_4$
Pearson symbol	:	tI82
Space group number	:	87
Space group symbol	:	$I4/m$
AFLOW prototype command	:	`aflow --proto=AB4C24D12_tI82_87_a_h_2h2i_hi --params=`$a, c/a, x_2, y_2, x_3, y_3, x_4, y_4, x_5, y_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8$

- (Papike, 1965) found that the composition of the Si-II (16*i*) site was actually $Al_{0.458}Si_{0.542}$. This is richer in aluminum than assumed by Pauling, who gave the total aluminum/silicon composition as $AlSi_3$ (Pauling, 1930). If we assume that the Si-I (8*h*) site is only filled by silicon atoms, then Pauling's composition for the Si-II site is $Al_{0.375}Si_{0.625}$. The name "marialite" scapolite distinguishes this from meionite scapolite, which replaces the sodium atoms by calcium but also includes SiO_4 and CO_3, which replace the chlorine atoms. According to Pauling, "The minerals of the scapolite group can be considered as solid solution of two end-members, marialite, $Na_4Al_3Si_9O_{24}Cl$, and meionite, $Ca_4Al_6Si_6O_{24}(SO_4,CO_3)$, in various proportions."
- Since the Al-Si sites are partially occupied with a higher concentration of Si, the positions are labeled as Si.

Body-centered Tetragonal primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= -\tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{y}} + \tfrac{1}{2}\,c\,\hat{\mathbf{z}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2}\,a\,\hat{\mathbf{y}} + \tfrac{1}{2}\,c\,\hat{\mathbf{z}}\\ \mathbf{a}_3 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{y}} - \tfrac{1}{2}\,c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(2*a*)	Cl
$\mathbf{B}_2$	=	$y_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + (x_2 + y_2)\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + y_2 a\,\hat{\mathbf{y}}$	(8*h*)	Na
$\mathbf{B}_3$	=	$-y_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 + (-x_2 - y_2)\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} - y_2 a\,\hat{\mathbf{y}}$	(8*h*)	Na
$\mathbf{B}_4$	=	$x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + (x_2 - y_2)\,\mathbf{a}_3$	=	$-y_2 a\,\hat{\mathbf{x}} + x_2 a\,\hat{\mathbf{y}}$	(8*h*)	Na
$\mathbf{B}_5$	=	$-x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + (-x_2 + y_2)\,\mathbf{a}_3$	=	$y_2 a\,\hat{\mathbf{x}} - x_2 a\,\hat{\mathbf{y}}$	(8*h*)	Na
$\mathbf{B}_6$	=	$y_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + (x_3 + y_3)\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 a\,\hat{\mathbf{y}}$	(8*h*)	O I
$\mathbf{B}_7$	=	$-y_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + (-x_3 - y_3)\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 a\,\hat{\mathbf{y}}$	(8*h*)	O I
$\mathbf{B}_8$	=	$x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + (x_3 - y_3)\,\mathbf{a}_3$	=	$-y_3 a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}}$	(8*h*)	O I
$\mathbf{B}_9$	=	$-x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + (-x_3 + y_3)\,\mathbf{a}_3$	=	$y_3 a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}}$	(8*h*)	O I
$\mathbf{B}_{10}$	=	$y_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + (x_4 + y_4)\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 a\,\hat{\mathbf{y}}$	(8*h*)	O II
$\mathbf{B}_{11}$	=	$-y_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + (-x_4 - y_4)\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 a\,\hat{\mathbf{y}}$	(8*h*)	O II
$\mathbf{B}_{12}$	=	$x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + (x_4 - y_4)\,\mathbf{a}_3$	=	$-y_4 a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}}$	(8*h*)	O II
$\mathbf{B}_{13}$	=	$-x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + (-x_4 + y_4)\,\mathbf{a}_3$	=	$y_4 a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}}$	(8*h*)	O II
$\mathbf{B}_{14}$	=	$y_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + (x_5 + y_5)\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 a\,\hat{\mathbf{y}}$	(8*h*)	Si I
$\mathbf{B}_{15}$	=	$-y_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + (-x_5 - y_5)\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 a\,\hat{\mathbf{y}}$	(8*h*)	Si I
$\mathbf{B}_{16}$	=	$x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + (x_5 - y_5)\,\mathbf{a}_3$	=	$-y_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}}$	(8*h*)	Si I
$\mathbf{B}_{17}$	=	$-x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + (-x_5 + y_5)\,\mathbf{a}_3$	=	$y_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}}$	(8*h*)	Si I
$\mathbf{B}_{18}$	=	$(y_6 + z_6)\,\mathbf{a}_1 + (x_6 + z_6)\,\mathbf{a}_2 + (x_6 + y_6)\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + y_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(16*i*)	O III
$\mathbf{B}_{19}$	=	$(-y_6 + z_6)\,\mathbf{a}_1 + (-x_6 + z_6)\,\mathbf{a}_2 + (-x_6 - y_6)\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - y_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(16*i*)	O III
$\mathbf{B}_{20}$	=	$(x_6 + z_6)\,\mathbf{a}_1 + (-y_6 + z_6)\,\mathbf{a}_2 + (x_6 - y_6)\,\mathbf{a}_3$	=	$-y_6 a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(16*i*)	O III
$\mathbf{B}_{21}$	=	$(-x_6 + z_6)\,\mathbf{a}_1 + (y_6 + z_6)\,\mathbf{a}_2 + (-x_6 + y_6)\,\mathbf{a}_3$	=	$y_6 a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(16*i*)	O III
$\mathbf{B}_{22}$	=	$(-y_6 - z_6)\,\mathbf{a}_1 + (-x_6 - z_6)\,\mathbf{a}_2 + (-x_6 - y_6)\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - y_6 a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(16*i*)	O III
$\mathbf{B}_{23}$	=	$(y_6 - z_6)\,\mathbf{a}_1 + (x_6 - z_6)\,\mathbf{a}_2 + (x_6 + y_6)\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + y_6 a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(16*i*)	O III
$\mathbf{B}_{24}$	=	$(-x_6 - z_6)\,\mathbf{a}_1 + (y_6 - z_6)\,\mathbf{a}_2 + (-x_6 + y_6)\,\mathbf{a}_3$	=	$y_6 a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(16*i*)	O III
$\mathbf{B}_{25}$	=	$(x_6 - z_6)\,\mathbf{a}_1 + (-y_6 - z_6)\,\mathbf{a}_2 + (x_6 - y_6)\,\mathbf{a}_3$	=	$-y_6 a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(16*i*)	O III
$\mathbf{B}_{26}$	=	$(y_7 + z_7)\,\mathbf{a}_1 + (x_7 + z_7)\,\mathbf{a}_2 + (x_7 + y_7)\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + y_7 a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(16*i*)	O IV
$\mathbf{B}_{27}$	=	$(-y_7 + z_7)\,\mathbf{a}_1 + (-x_7 + z_7)\,\mathbf{a}_2 + (-x_7 - y_7)\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} - y_7 a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(16*i*)	O IV

$\mathbf{B}_{28}$	=	$(x_7 + z_7)\,\mathbf{a}_1 + (-y_7 + z_7)\,\mathbf{a}_2 + (x_7 - y_7)\,\mathbf{a}_3$	=	$-y_7 a\,\hat{\mathbf{x}} + x_7 a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(16*i*)	O IV
$\mathbf{B}_{29}$	=	$(-x_7 + z_7)\,\mathbf{a}_1 + (y_7 + z_7)\,\mathbf{a}_2 + (-x_7 + y_7)\,\mathbf{a}_3$	=	$y_7 a\,\hat{\mathbf{x}} - x_7 a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(16*i*)	O IV
$\mathbf{B}_{30}$	=	$(-y_7 - z_7)\,\mathbf{a}_1 + (-x_7 - z_7)\,\mathbf{a}_2 + (-x_7 - y_7)\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} - y_7 a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	(16*i*)	O IV
$\mathbf{B}_{31}$	=	$(y_7 - z_7)\,\mathbf{a}_1 + (x_7 - z_7)\,\mathbf{a}_2 + (x_7 + y_7)\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + y_7 a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	(16*i*)	O IV
$\mathbf{B}_{32}$	=	$(-x_7 - z_7)\,\mathbf{a}_1 + (y_7 - z_7)\,\mathbf{a}_2 + (-x_7 + y_7)\,\mathbf{a}_3$	=	$y_7 a\,\hat{\mathbf{x}} - x_7 a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	(16*i*)	O IV
$\mathbf{B}_{33}$	=	$(x_7 - z_7)\,\mathbf{a}_1 + (-y_7 - z_7)\,\mathbf{a}_2 + (x_7 - y_7)\,\mathbf{a}_3$	=	$-y_7 a\,\hat{\mathbf{x}} + x_7 a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	(16*i*)	O IV
$\mathbf{B}_{34}$	=	$(y_8 + z_8)\,\mathbf{a}_1 + (x_8 + z_8)\,\mathbf{a}_2 + (x_8 + y_8)\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + y_8 a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(16*i*)	Si II
$\mathbf{B}_{35}$	=	$(-y_8 + z_8)\,\mathbf{a}_1 + (-x_8 + z_8)\,\mathbf{a}_2 + (-x_8 - y_8)\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} - y_8 a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(16*i*)	Si II
$\mathbf{B}_{36}$	=	$(x_8 + z_8)\,\mathbf{a}_1 + (-y_8 + z_8)\,\mathbf{a}_2 + (x_8 - y_8)\,\mathbf{a}_3$	=	$-y_8 a\,\hat{\mathbf{x}} + x_8 a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(16*i*)	Si II
$\mathbf{B}_{37}$	=	$(-x_8 + z_8)\,\mathbf{a}_1 + (y_8 + z_8)\,\mathbf{a}_2 + (-x_8 + y_8)\,\mathbf{a}_3$	=	$y_8 a\,\hat{\mathbf{x}} - x_8 a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(16*i*)	Si II
$\mathbf{B}_{38}$	=	$(-y_8 - z_8)\,\mathbf{a}_1 + (-x_8 - z_8)\,\mathbf{a}_2 + (-x_8 - y_8)\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} - y_8 a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(16*i*)	Si II
$\mathbf{B}_{39}$	=	$(y_8 - z_8)\,\mathbf{a}_1 + (x_8 - z_8)\,\mathbf{a}_2 + (x_8 + y_8)\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + y_8 a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(16*i*)	Si II
$\mathbf{B}_{40}$	=	$(-x_8 - z_8)\,\mathbf{a}_1 + (y_8 - z_8)\,\mathbf{a}_2 + (-x_8 + y_8)\,\mathbf{a}_3$	=	$y_8 a\,\hat{\mathbf{x}} - x_8 a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(16*i*)	Si II
$\mathbf{B}_{41}$	=	$(x_8 - z_8)\,\mathbf{a}_1 + (-y_8 - z_8)\,\mathbf{a}_2 + (x_8 - y_8)\,\mathbf{a}_3$	=	$-y_8 a\,\hat{\mathbf{x}} + x_8 a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(16*i*)	Si II

References:

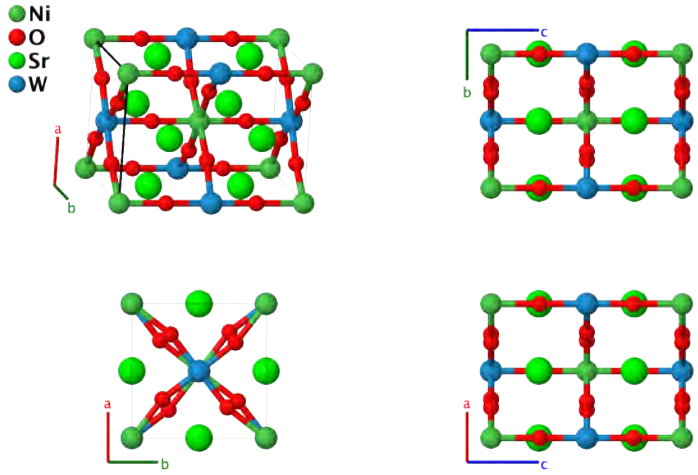
- J. J. Papike and T. Zoltai, *The crystal structure of a marialite scapolite*, Am. Mineral. **50**, 641–655 (1965).
- L. Pauling, *The Structure of Some Sodium and Calcium Aluminosilicates*, Proc. Natl. Acad. Sci. **16**, 453–459 (1930), doi:10.1073/pnas.16.7.453.

Geometry files:

- CIF: pp. 1672
- POSCAR: pp. 1673

Sr_2NiWO_6 Structure: AB6C2D_tI20_87_a_eh_d_b

http://aflow.org/prototype-encyclopedia/AB6C2D_tI20_87_a_eh_d_b



Prototype	:	NiO_6Sr_2W
AFLOW prototype label	:	AB6C2D_tI20_87_a_eh_d_b
***Strukturbericht* designation**	:	None
Pearson symbol	:	tI20
Space group number	:	87
Space group symbol	:	$I4/m$
AFLOW prototype command	:	`aflow --proto=AB6C2D_tI20_87_a_eh_d_b --params=`$a, c/a, z_4, x_5, y_5$

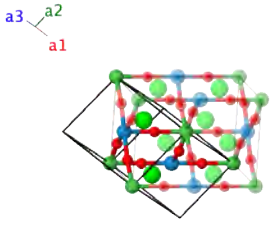
Other compounds with this structure

- Sr_2MgWO_6

- This double perovskite crystal is the ground state structure of Sr_2NiWO_6. Above 300 °C, it transforms into the cubic perovskite $E2_1$ structure. (Iwanga, 2000) places the strontium atoms on the (4*c*) Wyckoff position, but gives the coordinates for the (4*d*) site. The interatomic distances they give with this structure are consistent with the (4*d*) site.

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} - \frac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(2*a*)	Ni

$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	($2b$)	W
$\mathbf{B}_3$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($4d$)	Sr
$\mathbf{B}_4$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($4d$)	Sr
$\mathbf{B}_5$	=	$z_4\,\mathbf{a}_1 + z_4\,\mathbf{a}_2$	=	$z_4 c\,\hat{\mathbf{z}}$	($4e$)	O I
$\mathbf{B}_6$	=	$-z_4\,\mathbf{a}_1 - z_4\,\mathbf{a}_2$	=	$-z_4 c\,\hat{\mathbf{z}}$	($4e$)	O I
$\mathbf{B}_7$	=	$y_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + (x_5 + y_5)\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 a\,\hat{\mathbf{y}}$	($8h$)	O II
$\mathbf{B}_8$	=	$-y_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + (-x_5 - y_5)\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 a\,\hat{\mathbf{y}}$	($8h$)	O II
$\mathbf{B}_9$	=	$x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + (x_5 - y_5)\,\mathbf{a}_3$	=	$-y_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}}$	($8h$)	O II
$\mathbf{B}_{10}$	=	$-x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + (-x_5 + y_5)\,\mathbf{a}_3$	=	$y_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}}$	($8h$)	O II

References:

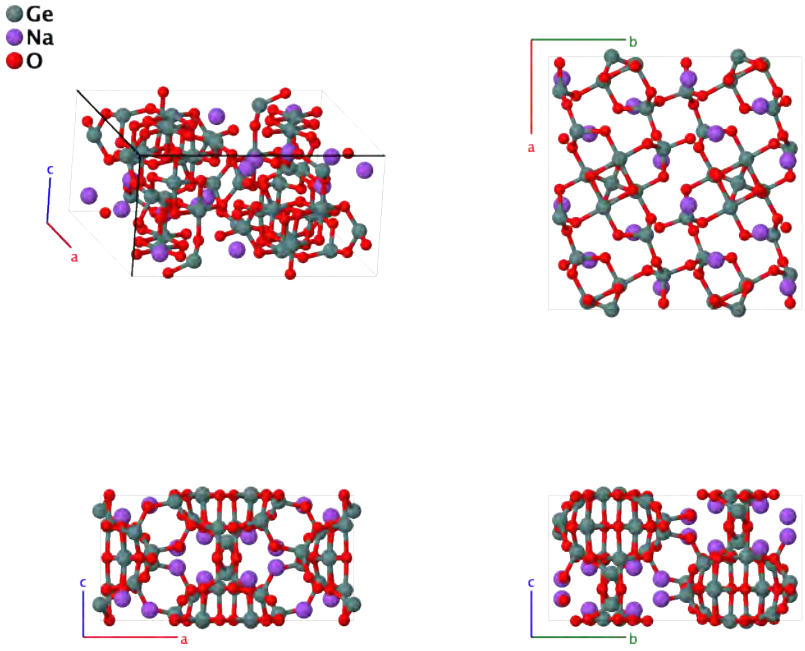
- D. Iwanaga, Y. Inaguma, and M. Itoh, *Structure and Magnetic Properties of Sr_2NiAO_6 (A = W, Te)*, Mater. Res. Bull. **35**, 449–457 (2000), doi:10.1016/S0025-5408(00)00222-1.

Geometry files:

- CIF: pp. 1673
- POSCAR: pp. 1673

$Na_4Ge_9O_{20}$ Structure: A9B4C20_tI132_88_a2f_f_5f

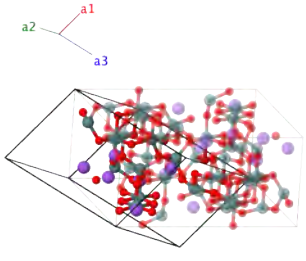
http://aflow.org/prototype-encyclopedia/A9B4C20_tI132_88_a2f_f_5f



Prototype	:	$Ge_9Na_4O_{20}$
AFLOW prototype label	:	A9B4C20_tI132_88_a2f_f_5f
***Strukturbericht* designation**	:	None
Pearson symbol	:	tI132
Space group number	:	88
Space group symbol	:	$I4_1/a$
AFLOW prototype command	:	`aflow --proto=A9B4C20_tI132_88_a2f_f_5f --params=`$a, c/a, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9$

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} - \frac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$\mathbf{B}_1$	=	$\frac{3}{8}\,\mathbf{a}_1 + \frac{1}{8}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{8}c\,\hat{\mathbf{z}}$	(4*a*)	Ge I
$\mathbf{B}_2$	=	$\frac{5}{8}\,\mathbf{a}_1 + \frac{7}{8}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{3}{8}c\,\hat{\mathbf{z}}$	(4*a*)	Ge I
$\mathbf{B}_3$	=	$(y_2 + z_2)\,\mathbf{a}_1 + (x_2 + z_2)\,\mathbf{a}_2 + (x_2 + y_2)\,\mathbf{a}_3$	=	$x_2a\,\hat{\mathbf{x}} + y_2a\,\hat{\mathbf{y}} + z_2c\,\hat{\mathbf{z}}$	(16*f*)	Ge II
$\mathbf{B}_4$	=	$\left(\frac{1}{2} - y_2 + z_2\right)\mathbf{a}_1 + (-x_2 + z_2)\,\mathbf{a}_2 + \left(\frac{1}{2} - x_2 - y_2\right)\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right)a\,\hat{\mathbf{y}} + z_2c\,\hat{\mathbf{z}}$	(16*f*)	Ge II
$\mathbf{B}_5$	=	$\left(\frac{1}{2} + x_2 + z_2\right)\mathbf{a}_1 + (-y_2 + z_2)\,\mathbf{a}_2 + (x_2 - y_2)\,\mathbf{a}_3$	=	$\left(\frac{3}{4} - y_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_2\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_2\right)c\,\hat{\mathbf{z}}$	(16*f*)	Ge II
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_2 + z_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_2 + z_2\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_2 + y_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + y_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_2\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_2\right)c\,\hat{\mathbf{z}}$	(16*f*)	Ge II
$\mathbf{B}_7$	=	$(-y_2 - z_2)\,\mathbf{a}_1 + (-x_2 - z_2)\,\mathbf{a}_2 + (-x_2 - y_2)\,\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}} - y_2a\,\hat{\mathbf{y}} - z_2c\,\hat{\mathbf{z}}$	(16*f*)	Ge II
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + y_2 - z_2\right)\mathbf{a}_1 + (x_2 - z_2)\,\mathbf{a}_2 + \left(\frac{1}{2} + x_2 + y_2\right)\mathbf{a}_3$	=	$x_2a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right)a\,\hat{\mathbf{y}} - z_2c\,\hat{\mathbf{z}}$	(16*f*)	Ge II
$\mathbf{B}_9$	=	$\left(\frac{1}{2} - x_2 - z_2\right)\mathbf{a}_1 + (y_2 - z_2)\,\mathbf{a}_2 + (-x_2 + y_2)\,\mathbf{a}_3$	=	$\left(-\frac{1}{4} + y_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_2\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_2\right)c\,\hat{\mathbf{z}}$	(16*f*)	Ge II
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} + x_2 - z_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_2 - z_2\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_2 - y_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - y_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_2\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_2\right)c\,\hat{\mathbf{z}}$	(16*f*)	Ge II
$\mathbf{B}_{11}$	=	$(y_3 + z_3)\,\mathbf{a}_1 + (x_3 + z_3)\,\mathbf{a}_2 + (x_3 + y_3)\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + y_3a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	(16*f*)	Ge III
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} - y_3 + z_3\right)\mathbf{a}_1 + (-x_3 + z_3)\,\mathbf{a}_2 + \left(\frac{1}{2} - x_3 - y_3\right)\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	(16*f*)	Ge III
$\mathbf{B}_{13}$	=	$\left(\frac{1}{2} + x_3 + z_3\right)\mathbf{a}_1 + (-y_3 + z_3)\,\mathbf{a}_2 + (x_3 - y_3)\,\mathbf{a}_3$	=	$\left(\frac{3}{4} - y_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_3\right)c\,\hat{\mathbf{z}}$	(16*f*)	Ge III
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - x_3 + z_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3 + z_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_3 + y_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + y_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_3\right)c\,\hat{\mathbf{z}}$	(16*f*)	Ge III
$\mathbf{B}_{15}$	=	$(-y_3 - z_3)\,\mathbf{a}_1 + (-x_3 - z_3)\,\mathbf{a}_2 + (-x_3 - y_3)\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - y_3a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	(16*f*)	Ge III
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + y_3 - z_3\right)\mathbf{a}_1 + (x_3 - z_3)\,\mathbf{a}_2 + \left(\frac{1}{2} + x_3 + y_3\right)\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	(16*f*)	Ge III
$\mathbf{B}_{17}$	=	$\left(\frac{1}{2} - x_3 - z_3\right)\mathbf{a}_1 + (y_3 - z_3)\,\mathbf{a}_2 + (-x_3 + y_3)\,\mathbf{a}_3$	=	$\left(-\frac{1}{4} + y_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_3\right)c\,\hat{\mathbf{z}}$	(16*f*)	Ge III
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} + x_3 - z_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3 - z_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_3 - y_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - y_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_3\right)c\,\hat{\mathbf{z}}$	(16*f*)	Ge III
$\mathbf{B}_{19}$	=	$(y_4 + z_4)\,\mathbf{a}_1 + (x_4 + z_4)\,\mathbf{a}_2 + (x_4 + y_4)\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + y_4a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	(16*f*)	Na
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} - y_4 + z_4\right)\mathbf{a}_1 + (-x_4 + z_4)\,\mathbf{a}_2 + \left(\frac{1}{2} - x_4 - y_4\right)\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	(16*f*)	Na
$\mathbf{B}_{21}$	=	$\left(\frac{1}{2} + x_4 + z_4\right)\mathbf{a}_1 + (-y_4 + z_4)\,\mathbf{a}_2 + (x_4 - y_4)\,\mathbf{a}_3$	=	$\left(\frac{3}{4} - y_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_4\right)c\,\hat{\mathbf{z}}$	(16*f*)	Na
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} - x_4 + z_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4 + z_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_4 + y_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + y_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_4\right)c\,\hat{\mathbf{z}}$	(16*f*)	Na

$\mathbf{B}_{23}$	=	$(-y_4 - z_4)\,\mathbf{a}_1 + (-x_4 - z_4)\,\mathbf{a}_2 + (-x_4 - y_4)\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 a\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	$(16f)$	Na
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} + y_4 - z_4\right)\mathbf{a}_1 + (x_4 - z_4)\,\mathbf{a}_2 + \left(\frac{1}{2} + x_4 + y_4\right)\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right) a\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	$(16f)$	Na
$\mathbf{B}_{25}$	=	$\left(\frac{1}{2} - x_4 - z_4\right)\mathbf{a}_1 + (y_4 - z_4)\,\mathbf{a}_2 + (-x_4 + y_4)\,\mathbf{a}_3$	=	$\left(-\frac{1}{4} + y_4\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_4\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_4\right) c\,\hat{\mathbf{z}}$	$(16f)$	Na
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} + x_4 - z_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4 - z_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_4 - y_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - y_4\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_4\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_4\right) c\,\hat{\mathbf{z}}$	$(16f)$	Na
$\mathbf{B}_{27}$	=	$(y_5 + z_5)\,\mathbf{a}_1 + (x_5 + z_5)\,\mathbf{a}_2 + (x_5 + y_5)\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(16f)$	O I
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} - y_5 + z_5\right)\mathbf{a}_1 + (-x_5 + z_5)\,\mathbf{a}_2 + \left(\frac{1}{2} - x_5 - y_5\right)\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right) a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(16f)$	O I
$\mathbf{B}_{29}$	=	$\left(\frac{1}{2} + x_5 + z_5\right)\mathbf{a}_1 + (-y_5 + z_5)\,\mathbf{a}_2 + (x_5 - y_5)\,\mathbf{a}_3$	=	$\left(\frac{3}{4} - y_5\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_5\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_5\right) c\,\hat{\mathbf{z}}$	$(16f)$	O I
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2} - x_5 + z_5\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_5 + z_5\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_5 + y_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + y_5\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_5\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_5\right) c\,\hat{\mathbf{z}}$	$(16f)$	O I
$\mathbf{B}_{31}$	=	$(-y_5 - z_5)\,\mathbf{a}_1 + (-x_5 - z_5)\,\mathbf{a}_2 + (-x_5 - y_5)\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	$(16f)$	O I
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2} + y_5 - z_5\right)\mathbf{a}_1 + (x_5 - z_5)\,\mathbf{a}_2 + \left(\frac{1}{2} + x_5 + y_5\right)\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right) a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	$(16f)$	O I
$\mathbf{B}_{33}$	=	$\left(\frac{1}{2} - x_5 - z_5\right)\mathbf{a}_1 + (y_5 - z_5)\,\mathbf{a}_2 + (-x_5 + y_5)\,\mathbf{a}_3$	=	$\left(-\frac{1}{4} + y_5\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_5\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_5\right) c\,\hat{\mathbf{z}}$	$(16f)$	O I
$\mathbf{B}_{34}$	=	$\left(\frac{1}{2} + x_5 - z_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_5 - z_5\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_5 - y_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - y_5\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_5\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_5\right) c\,\hat{\mathbf{z}}$	$(16f)$	O I
$\mathbf{B}_{35}$	=	$(y_6 + z_6)\,\mathbf{a}_1 + (x_6 + z_6)\,\mathbf{a}_2 + (x_6 + y_6)\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + y_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(16f)$	O II
$\mathbf{B}_{36}$	=	$\left(\frac{1}{2} - y_6 + z_6\right)\mathbf{a}_1 + (-x_6 + z_6)\,\mathbf{a}_2 + \left(\frac{1}{2} - x_6 - y_6\right)\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right) a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(16f)$	O II
$\mathbf{B}_{37}$	=	$\left(\frac{1}{2} + x_6 + z_6\right)\mathbf{a}_1 + (-y_6 + z_6)\,\mathbf{a}_2 + (x_6 - y_6)\,\mathbf{a}_3$	=	$\left(\frac{3}{4} - y_6\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_6\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_6\right) c\,\hat{\mathbf{z}}$	$(16f)$	O II
$\mathbf{B}_{38}$	=	$\left(\frac{1}{2} - x_6 + z_6\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_6 + z_6\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_6 + y_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + y_6\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_6\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_6\right) c\,\hat{\mathbf{z}}$	$(16f)$	O II
$\mathbf{B}_{39}$	=	$(-y_6 - z_6)\,\mathbf{a}_1 + (-x_6 - z_6)\,\mathbf{a}_2 + (-x_6 - y_6)\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - y_6 a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	$(16f)$	O II
$\mathbf{B}_{40}$	=	$\left(\frac{1}{2} + y_6 - z_6\right)\mathbf{a}_1 + (x_6 - z_6)\,\mathbf{a}_2 + \left(\frac{1}{2} + x_6 + y_6\right)\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right) a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	$(16f)$	O II
$\mathbf{B}_{41}$	=	$\left(\frac{1}{2} - x_6 - z_6\right)\mathbf{a}_1 + (y_6 - z_6)\,\mathbf{a}_2 + (-x_6 + y_6)\,\mathbf{a}_3$	=	$\left(-\frac{1}{4} + y_6\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_6\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_6\right) c\,\hat{\mathbf{z}}$	$(16f)$	O II
$\mathbf{B}_{42}$	=	$\left(\frac{1}{2} + x_6 - z_6\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_6 - z_6\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_6 - y_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - y_6\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_6\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_6\right) c\,\hat{\mathbf{z}}$	$(16f)$	O II
$\mathbf{B}_{43}$	=	$(y_7 + z_7)\,\mathbf{a}_1 + (x_7 + z_7)\,\mathbf{a}_2 + (x_7 + y_7)\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + y_7 a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	$(16f)$	O III

$$\mathbf{B}_{44} = \left(\tfrac{1}{2} - y_7 + z_7\right)\mathbf{a}_1 + (-x_7 + z_7)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_7 - y_7\right)\mathbf{a}_3 = -x_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_7\right)a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}} \quad (16f) \quad \text{O III}$$

$$\mathbf{B}_{45} = \left(\tfrac{1}{2} + x_7 + z_7\right)\mathbf{a}_1 + (-y_7 + z_7)\,\mathbf{a}_2 + (x_7 - y_7)\,\mathbf{a}_3 = \left(\tfrac{3}{4} - y_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_7\right)c\,\hat{\mathbf{z}} \quad (16f) \quad \text{O III}$$

$$\mathbf{B}_{46} = \left(\tfrac{1}{2} - x_7 + z_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_7 + z_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_7 + y_7\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + y_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_7\right)c\,\hat{\mathbf{z}} \quad (16f) \quad \text{O III}$$

$$\mathbf{B}_{47} = (-y_7 - z_7)\,\mathbf{a}_1 + (-x_7 - z_7)\,\mathbf{a}_2 + (-x_7 - y_7)\,\mathbf{a}_3 = -x_7 a\,\hat{\mathbf{x}} - y_7 a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}} \quad (16f) \quad \text{O III}$$

$$\mathbf{B}_{48} = \left(\tfrac{1}{2} + y_7 - z_7\right)\mathbf{a}_1 + (x_7 - z_7)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_7 + y_7\right)\mathbf{a}_3 = x_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_7\right)a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}} \quad (16f) \quad \text{O III}$$

$$\mathbf{B}_{49} = \left(\tfrac{1}{2} - x_7 - z_7\right)\mathbf{a}_1 + (y_7 - z_7)\,\mathbf{a}_2 + (-x_7 + y_7)\,\mathbf{a}_3 = \left(-\tfrac{1}{4} + y_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - z_7\right)c\,\hat{\mathbf{z}} \quad (16f) \quad \text{O III}$$

$$\mathbf{B}_{50} = \left(\tfrac{1}{2} + x_7 - z_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_7 - z_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_7 - y_7\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - y_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - z_7\right)c\,\hat{\mathbf{z}} \quad (16f) \quad \text{O III}$$

$$\mathbf{B}_{51} = (y_8 + z_8)\,\mathbf{a}_1 + (x_8 + z_8)\,\mathbf{a}_2 + (x_8 + y_8)\,\mathbf{a}_3 = x_8 a\,\hat{\mathbf{x}} + y_8 a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}} \quad (16f) \quad \text{O IV}$$

$$\mathbf{B}_{52} = \left(\tfrac{1}{2} - y_8 + z_8\right)\mathbf{a}_1 + (-x_8 + z_8)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_8 - y_8\right)\mathbf{a}_3 = -x_8 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_8\right)a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}} \quad (16f) \quad \text{O IV}$$

$$\mathbf{B}_{53} = \left(\tfrac{1}{2} + x_8 + z_8\right)\mathbf{a}_1 + (-y_8 + z_8)\,\mathbf{a}_2 + (x_8 - y_8)\,\mathbf{a}_3 = \left(\tfrac{3}{4} - y_8\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_8\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_8\right)c\,\hat{\mathbf{z}} \quad (16f) \quad \text{O IV}$$

$$\mathbf{B}_{54} = \left(\tfrac{1}{2} - x_8 + z_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_8 + z_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_8 + y_8\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + y_8\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_8\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_8\right)c\,\hat{\mathbf{z}} \quad (16f) \quad \text{O IV}$$

$$\mathbf{B}_{55} = (-y_8 - z_8)\,\mathbf{a}_1 + (-x_8 - z_8)\,\mathbf{a}_2 + (-x_8 - y_8)\,\mathbf{a}_3 = -x_8 a\,\hat{\mathbf{x}} - y_8 a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}} \quad (16f) \quad \text{O IV}$$

$$\mathbf{B}_{56} = \left(\tfrac{1}{2} + y_8 - z_8\right)\mathbf{a}_1 + (x_8 - z_8)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_8 + y_8\right)\mathbf{a}_3 = x_8 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_8\right)a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}} \quad (16f) \quad \text{O IV}$$

$$\mathbf{B}_{57} = \left(\tfrac{1}{2} - x_8 - z_8\right)\mathbf{a}_1 + (y_8 - z_8)\,\mathbf{a}_2 + (-x_8 + y_8)\,\mathbf{a}_3 = \left(-\tfrac{1}{4} + y_8\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_8\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - z_8\right)c\,\hat{\mathbf{z}} \quad (16f) \quad \text{O IV}$$

$$\mathbf{B}_{58} = \left(\tfrac{1}{2} + x_8 - z_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_8 - z_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_8 - y_8\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - y_8\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_8\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - z_8\right)c\,\hat{\mathbf{z}} \quad (16f) \quad \text{O IV}$$

$$\mathbf{B}_{59} = (y_9 + z_9)\,\mathbf{a}_1 + (x_9 + z_9)\,\mathbf{a}_2 + (x_9 + y_9)\,\mathbf{a}_3 = x_9 a\,\hat{\mathbf{x}} + y_9 a\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}} \quad (16f) \quad \text{O V}$$

$$\mathbf{B}_{60} = \left(\tfrac{1}{2} - y_9 + z_9\right)\mathbf{a}_1 + (-x_9 + z_9)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_9 - y_9\right)\mathbf{a}_3 = -x_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_9\right)a\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}} \quad (16f) \quad \text{O V}$$

$$\mathbf{B}_{61} = \left(\tfrac{1}{2} + x_9 + z_9\right)\mathbf{a}_1 + (-y_9 + z_9)\,\mathbf{a}_2 + (x_9 - y_9)\,\mathbf{a}_3 = \left(\tfrac{3}{4} - y_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_9\right)c\,\hat{\mathbf{z}} \quad (16f) \quad \text{O V}$$

$$\mathbf{B}_{62} = \left(\tfrac{1}{2} - x_9 + z_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_9 + z_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_9 + y_9\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + y_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_9\right)c\,\hat{\mathbf{z}} \quad (16f) \quad \text{O V}$$

$$\mathbf{B}_{63} = (-y_9 - z_9)\,\mathbf{a}_1 + (-x_9 - z_9)\,\mathbf{a}_2 + (-x_9 - y_9)\,\mathbf{a}_3 = -x_9 a\,\hat{\mathbf{x}} - y_9 a\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}} \quad (16f) \quad \text{O V}$$

$$\mathbf{B}_{64} = \left(\tfrac{1}{2} + y_9 - z_9\right)\mathbf{a}_1 + (x_9 - z_9)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_9 + y_9\right)\mathbf{a}_3 = x_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_9\right)a\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}} \quad (16f) \quad \text{O V}$$

$\mathbf{B}_{65}$	=	$\left(\frac{1}{2} - x_9 - z_9\right)\mathbf{a}_1 + (y_9 - z_9)\,\mathbf{a}_2 + (-x_9 + y_9)\,\mathbf{a}_3$	=	$\left(-\frac{1}{4} + y_9\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_9\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_9\right)c\,\hat{\mathbf{z}}$	$(16f)$	O V
$\mathbf{B}_{66}$	=	$\left(\frac{1}{2} + x_9 - z_9\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_9 - z_9\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_9 - y_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - y_9\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_9\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_9\right)c\,\hat{\mathbf{z}}$	$(16f)$	O V

References:

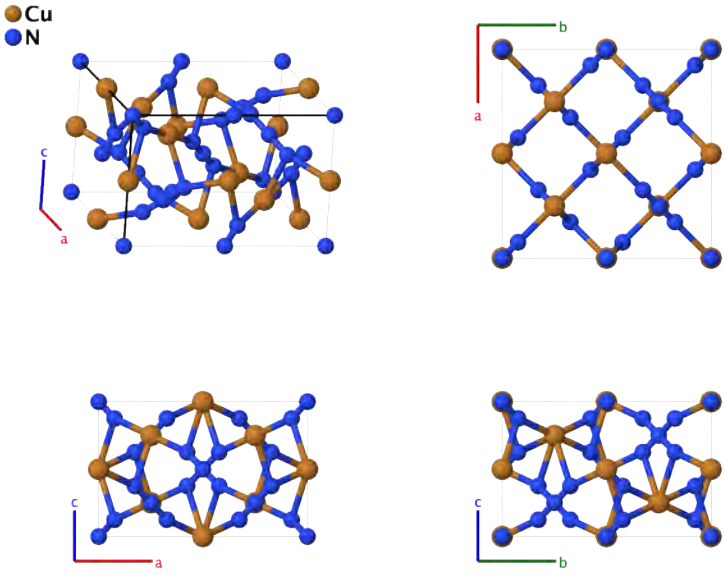
- N. Ingri and G. Lundgren, *The Crystal Structure of* $Na_4Ge_9O_{20}$, Acta Chem. Scand. **17**, 617–633 (1963), doi:10.3891/acta.chem.scand.17-0617.

Geometry files:

- CIF: pp. 1673
- POSCAR: pp. 1674

Copper (I) Azide (CuN_3) Structure: AB3_tI32_88_d_cf

http://aflow.org/prototype-encyclopedia/AB3_tI32_88_d_cf



Prototype	:	CuN_3
AFLOW prototype label	:	AB3_tI32_88_d_cf
***Strukturbericht* designation**	:	None
Pearson symbol	:	tI32
Space group number	:	88
Space group symbol	:	$I4_1/a$
AFLOW prototype command	:	`aflow --proto=AB3_tI32_88_d_cf --params=`$a, c/a, x_3, y_3, z_3$

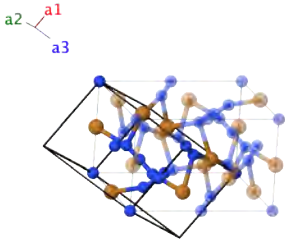
Other compounds with this structure

- AgN_3 and TiN_3

- Not to be confused with Copper (II) Azide, $Cu(N_3)_2$, an explosive.
- (Wilsdorf, 1948) gave the Wyckoff positions in setting 1 of space group #88. We used FINDSYM to translate this to the standard setting 2.

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{y}} + \tfrac{1}{2}\,c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2}\,a\,\hat{\mathbf{y}} + \tfrac{1}{2}\,c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{y}} - \tfrac{1}{2}\,c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($8c$)	N I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}}$	($8c$)	N I
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($8c$)	N I
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($8c$)	N I
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	($8d$)	Cu
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	($8d$)	Cu
$\mathbf{B}_7$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} - \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($8d$)	Cu
$\mathbf{B}_8$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} - \frac{1}{4}c\,\hat{\mathbf{z}}$	($8d$)	Cu
$\mathbf{B}_9$	=	$(y_3 + z_3)\,\mathbf{a}_1 + (x_3 + z_3)\,\mathbf{a}_2 + (x_3 + y_3)\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + y_3a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	($16f$)	N II
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - y_3 + z_3\right)\mathbf{a}_1 + (-x_3 + z_3)\,\mathbf{a}_2 + \left(\frac{1}{2} - x_3 - y_3\right)\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	($16f$)	N II
$\mathbf{B}_{11}$	=	$\left(\frac{1}{2} + x_3 + z_3\right)\mathbf{a}_1 + (-y_3 + z_3)\,\mathbf{a}_2 + (x_3 - y_3)\,\mathbf{a}_3$	=	$\left(\frac{3}{4} - y_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_3\right)c\,\hat{\mathbf{z}}$	($16f$)	N II
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} - x_3 + z_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3 + z_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_3 + y_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + y_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_3\right)c\,\hat{\mathbf{z}}$	($16f$)	N II
$\mathbf{B}_{13}$	=	$(-y_3 - z_3)\,\mathbf{a}_1 + (-x_3 - z_3)\,\mathbf{a}_2 + (-x_3 - y_3)\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - y_3a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	($16f$)	N II
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} + y_3 - z_3\right)\mathbf{a}_1 + (x_3 - z_3)\,\mathbf{a}_2 + \left(\frac{1}{2} + x_3 + y_3\right)\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	($16f$)	N II
$\mathbf{B}_{15}$	=	$\left(\frac{1}{2} - x_3 - z_3\right)\mathbf{a}_1 + (y_3 - z_3)\,\mathbf{a}_2 + (-x_3 + y_3)\,\mathbf{a}_3$	=	$\left(-\frac{1}{4} + y_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_3\right)c\,\hat{\mathbf{z}}$	($16f$)	N II
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_3 - z_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3 - z_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_3 - y_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - y_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_3\right)c\,\hat{\mathbf{z}}$	($16f$)	N II

References:

- H. Wilsdorf, *Die Kristallstruktur des einwertigen Kupferazids, CuN_3*, Acta Cryst. **1**, 115–118 (1948), doi:10.1107/S0365110X48000314.

Found in:

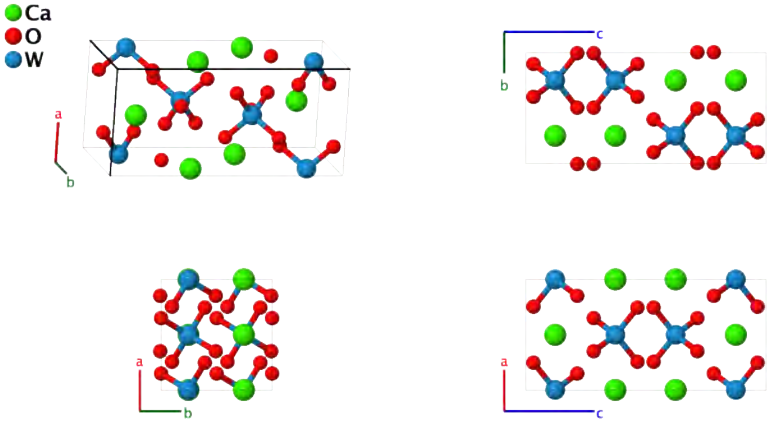
- W. Zhu and H. Xiao, Ab initio *study of electronic structure and optical properties of heavy-metal azides: TlN_3, AgN_3, and CuN_3*, J. Comput. Chem. **29**, 176–184 (2008), doi:10.1002/jcc.20682.

Geometry files:

- CIF: pp. 1674
- POSCAR: pp. 1675

Scheelite ($CaWO_4$, $H0_4$) Structure: AB4C_tI24_88_b_f_a

http://aflow.org/prototype-encyclopedia/AB4C_tI24_88_b_f_a



Prototype	:	CaO_4W
AFLOW prototype label	:	AB4C_tI24_88_b_f_a
***Strukturbericht* designation**	:	$H0_4$
Pearson symbol	:	tI24
Space group number	:	88
Space group symbol	:	$I4_1/a$
AFLOW prototype command	:	`aflow --proto=AB4C_tI24_88_b_f_a --params=`$a, c/a, x_3, y_3, z_3$

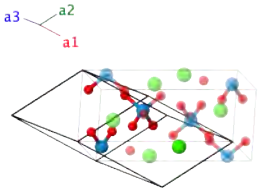
Other compounds with this structure

- $ZrSiO_4$, $LaNbO_4$, $YNbO_4$, (Y,RE)NbO_4 (fergusonite), YVO_4, $BiVO_4$, $BaWO_4$, $PbWO_4$ (wulfenite), $SrWO_4$, $EuWO_4$, $PbMoO_4$ (stolzite), $SrMoO_4$, $CaMoO_4$ (powellite), $CdMoO_4$, $KReO_4$, $TlReO_4$, $AgReO_4$, and $NaAlH_4$

- (Ewald, 1931) originally gave this structure the *Strukturbericht* designation $H4$, but this was changed to $H0_4$ in (Gottfried, 1937).

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}a\,\hat{\mathbf{x}} - \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} - \tfrac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\tfrac{3}{8}\mathbf{a}_1 + \tfrac{1}{8}\mathbf{a}_2 + \tfrac{1}{4}\mathbf{a}_3$	=	$\tfrac{1}{4}a\,\hat{\mathbf{y}} + \tfrac{1}{8}c\,\hat{\mathbf{z}}$	$(4a)$	W
$\mathbf{B}_2$	=	$\tfrac{5}{8}\mathbf{a}_1 + \tfrac{7}{8}\mathbf{a}_2 + \tfrac{3}{4}\mathbf{a}_3$	=	$\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{4}a\,\hat{\mathbf{y}} + \tfrac{3}{8}c\,\hat{\mathbf{z}}$	$(4a)$	W

$\mathbf{B}_3$	=	$\frac{7}{8}\,\mathbf{a}_1 + \frac{5}{8}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{y}} + \frac{5}{8}c\,\hat{\mathbf{z}}$	(4*b*)	Ca
$\mathbf{B}_4$	=	$\frac{1}{8}\,\mathbf{a}_1 + \frac{3}{8}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{7}{8}c\,\hat{\mathbf{z}}$	(4*b*)	Ca
$\mathbf{B}_5$	=	$(y_3 + z_3)\,\mathbf{a}_1 + (x_3 + z_3)\,\mathbf{a}_2 + (x_3 + y_3)\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + y_3a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	(16*f*)	O
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - y_3 + z_3\right)\mathbf{a}_1 + (-x_3 + z_3)\,\mathbf{a}_2 + \left(\frac{1}{2} - x_3 - y_3\right)\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	(16*f*)	O
$\mathbf{B}_7$	=	$\left(\frac{1}{2} + x_3 + z_3\right)\mathbf{a}_1 + (-y_3 + z_3)\,\mathbf{a}_2 + (x_3 - y_3)\,\mathbf{a}_3$	=	$\left(\frac{3}{4} - y_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_3\right)c\,\hat{\mathbf{z}}$	(16*f*)	O
$\mathbf{B}_8$	=	$\left(\frac{1}{2} - x_3 + z_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3 + z_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_3 + y_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + y_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_3\right)c\,\hat{\mathbf{z}}$	(16*f*)	O
$\mathbf{B}_9$	=	$(-y_3 - z_3)\,\mathbf{a}_1 + (-x_3 - z_3)\,\mathbf{a}_2 + (-x_3 - y_3)\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - y_3a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	(16*f*)	O
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} + y_3 - z_3\right)\mathbf{a}_1 + (x_3 - z_3)\,\mathbf{a}_2 + \left(\frac{1}{2} + x_3 + y_3\right)\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	(16*f*)	O
$\mathbf{B}_{11}$	=	$\left(\frac{1}{2} - x_3 - z_3\right)\mathbf{a}_1 + (y_3 - z_3)\,\mathbf{a}_2 + (-x_3 + y_3)\,\mathbf{a}_3$	=	$\left(-\frac{1}{4} + y_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_3\right)c\,\hat{\mathbf{z}}$	(16*f*)	O
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3 - z_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3 - z_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_3 - y_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - y_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_3\right)c\,\hat{\mathbf{z}}$	(16*f*)	O

References:

- R. M. Hazen, L. W. Finger, and J. W. E. Mariathasan, *High-pressure crystal chemistry of scheelite-type tungstates and molybdates*, J. Phys. Chem. Solids **46**, 253–263 (1985), doi:10.1016/0022-3697(85)90039-3.
- P. P. Ewald and C. Hermann, eds., *Strukturbericht 1913-1928* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1931).
- C. Gottfried and F. Schossberger, eds., *Strukturbericht Band III 1933-1935* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Found in:

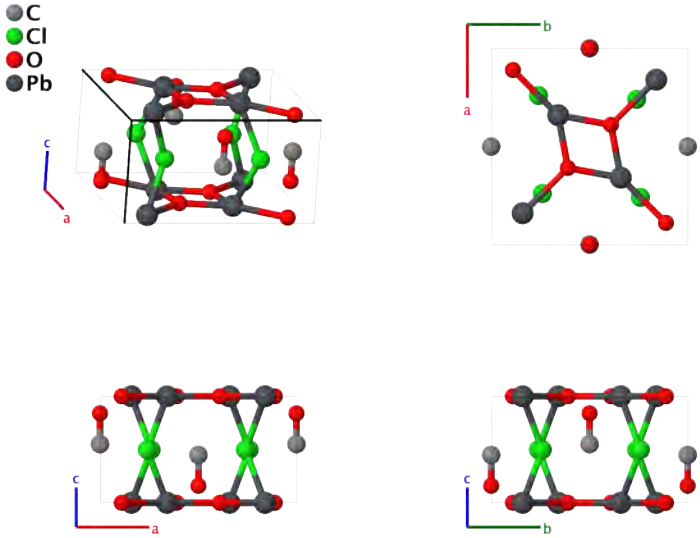
- Y. Zhang, N. A. W. Holzwarth, and R. T. Williams, *Electronic band structures of the scheelite materials $CaMoO_4$, $CaWO_4$, $PbMoO_4$, and $PbWO_4$*, Phys. Rev. B **57**, 12738–12750 (1988), doi:10.1103/PhysRevB.57.12738.

Geometry files:

- CIF: pp. 1675
- POSCAR: pp. 1675

$G7_5$ ($PbCO_3 \cdot PbCl_2$, Phosgenite) (*obsolete*) Structure: AB2C3D2_tP16_90_c_f_ce_e

http://aflow.org/prototype-encyclopedia/AB2C3D2_tP16_90_c_f_ce_e

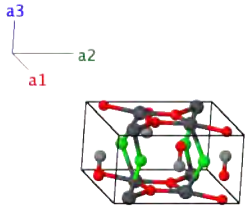


Prototype	:	$CCl_2O_3Pb_2$
AFLOW prototype label	:	AB2C3D2_tP16_90_c_f_ce_e
***Strukturbericht* designation**	:	$G7_5$
Pearson symbol	:	tP16
Space group number	:	90
Space group symbol	:	$P42_12$
AFLOW prototype command	:	`aflow --proto=AB2C3D2_tP16_90_c_f_ce_e --params=`$a, c/a, z_1, z_2, x_3, x_4, x_5$

- (Onorato, 1934) made an early determination of the structure of phosgenite, and (Gottfried, 1937) assigned it the *Strukturbericht* designation $G7_5$. However, the "proposed structure [was] based on photographic data and partly on steric considerations" (Giuseppetti, 1974). Subsequent investigations showed that the true phosgenite structure was in space group $P4/mbm$ #127. We list the original structure here as part of the historical record.

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{2}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	($2c$)	C
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 - z_1\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - z_1 c\,\hat{\mathbf{z}}$	($2c$)	C
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	($2c$)	O I
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 - z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - z_2 c\,\hat{\mathbf{z}}$	($2c$)	O I
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2$	=	$x_3 a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}}$	($4e$)	O II
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2$	=	$-x_3 a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}}$	($4e$)	O II
$\mathbf{B}_7$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}}$	($4e$)	O II
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}}$	($4e$)	O II
$\mathbf{B}_9$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2$	=	$x_4 a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}}$	($4e$)	Pb
$\mathbf{B}_{10}$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2$	=	$-x_4 a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}}$	($4e$)	Pb
$\mathbf{B}_{11}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_4\right)\mathbf{a}_2$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{y}}$	($4e$)	Pb
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_2$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{y}}$	($4e$)	Pb
$\mathbf{B}_{13}$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4f$)	Cl
$\mathbf{B}_{14}$	=	$-x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4f$)	Cl
$\mathbf{B}_{15}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_5\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4f$)	Cl
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_5\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4f$)	Cl

References:

- E. Onorato, *La struttura della Fosgenite*, Period. Mineral. **5**, 37–61 (1934).
- C. Gottfried and F. Schossberger, eds., *Strukturbericht Band III 1933-1935* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Found in:

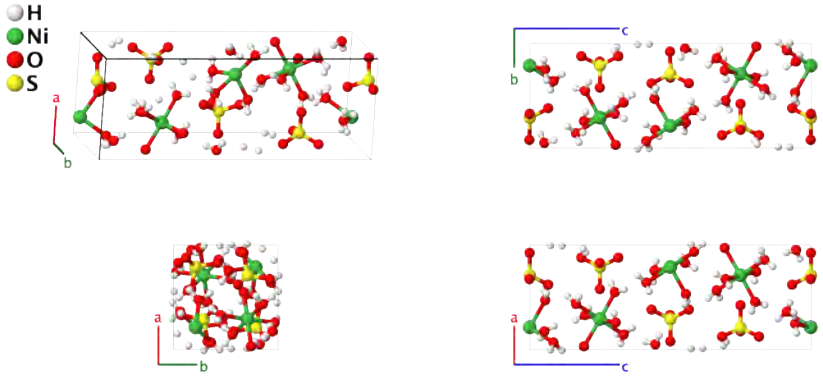
- G. Giuseppetti and C. Tadini, *Reexamination of the crystal structure of phosgenite, $Pb_2Cl_2(CO_3)$*, Tschermaks Min. Petr. Mitt. **21**, 101–109 (1974), doi:10.1007/BF01081262.

Geometry files:

- CIF: pp. 1675
- POSCAR: pp. 1676

Retgersite (α-$NiSO_4 \cdot 6H_2O$, $H4_5$) Structure: A12BC10D_tP96_92_6b_a_5b_a

http://aflow.org/prototype-encyclopedia/A12BC10D_tP96_92_6b_a_5b_a

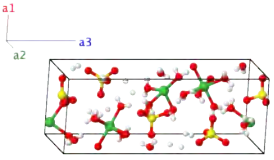


Prototype	:	$H_{12}NiO_{10}S$
AFLOW prototype label	:	A12BC10D_tP96_92_6b_a_5b_a
***Strukturbericht* designation**	:	$H4_5$
Pearson symbol	:	tP96
Space group number	:	92
Space group symbol	:	$P4_12_12$
AFLOW prototype command	:	`aflow --proto=A12BC10D_tP96_92_6b_a_5b_a --params=`$a, c/a, x_1, x_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}$

- This structure is nearly identical to the one presented in (Hermann, 1937) as $H4_5$, but now includes the positions of the hydrogen atoms.
- This compound can also be found in the enantiomorphic space group $P4_32_12$ #98.

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + x_1\,\mathbf{a}_2$	=	$x_1 a\,\hat{\mathbf{x}} + x_1 a\,\hat{\mathbf{y}}$	($4a$)	Ni
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 - x_1\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} - x_1 a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4a$)	Ni
$\mathbf{B}_3$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_1\right)\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_1\right)a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($4a$)	Ni
$\mathbf{B}_4$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_1\right)\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_1\right)a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	($4a$)	Ni
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2$	=	$x_2 a\,\hat{\mathbf{x}} + x_2 a\,\hat{\mathbf{y}}$	($4a$)	S

$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} - x_2 a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*a*)	S
$\mathbf{B}_7$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*a*)	S
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(4*a*)	S
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(8*b*)	H I
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(8*b*)	H I
$\mathbf{B}_{11}$	=	$\left(\frac{1}{2} - y_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + \left(\frac{1}{4} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_3\right)c\,\hat{\mathbf{z}}$	(8*b*)	H I
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + y_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 + \left(\frac{3}{4} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{3}{4} + z_3\right)c\,\hat{\mathbf{z}}$	(8*b*)	H I
$\mathbf{B}_{13}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\frac{1}{4} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_3\right)c\,\hat{\mathbf{z}}$	(8*b*)	H I
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + \left(\frac{3}{4} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{y}} + \left(\frac{3}{4} - z_3\right)c\,\hat{\mathbf{z}}$	(8*b*)	H I
$\mathbf{B}_{15}$	=	$y_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$y_3 a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(8*b*)	H I
$\mathbf{B}_{16}$	=	$-y_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$-y_3 a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(8*b*)	H I
$\mathbf{B}_{17}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*b*)	H II
$\mathbf{B}_{18}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(8*b*)	H II
$\mathbf{B}_{19}$	=	$\left(\frac{1}{2} - y_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_4\right)\mathbf{a}_2 + \left(\frac{1}{4} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_4\right)c\,\hat{\mathbf{z}}$	(8*b*)	H II
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} + y_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_2 + \left(\frac{3}{4} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{3}{4} + z_4\right)c\,\hat{\mathbf{z}}$	(8*b*)	H II
$\mathbf{B}_{21}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\frac{1}{4} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_4\right)c\,\hat{\mathbf{z}}$	(8*b*)	H II
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + \left(\frac{3}{4} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)a\,\hat{\mathbf{y}} + \left(\frac{3}{4} - z_4\right)c\,\hat{\mathbf{z}}$	(8*b*)	H II
$\mathbf{B}_{23}$	=	$y_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$y_4 a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*b*)	H II
$\mathbf{B}_{24}$	=	$-y_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$-y_4 a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(8*b*)	H II
$\mathbf{B}_{25}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(8*b*)	H III
$\mathbf{B}_{26}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	(8*b*)	H III
$\mathbf{B}_{27}$	=	$\left(\frac{1}{2} - y_5\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_5\right)\mathbf{a}_2 + \left(\frac{1}{4} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_5\right)c\,\hat{\mathbf{z}}$	(8*b*)	H III
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} + y_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_5\right)\mathbf{a}_2 + \left(\frac{3}{4} + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{y}} + \left(\frac{3}{4} + z_5\right)c\,\hat{\mathbf{z}}$	(8*b*)	H III
$\mathbf{B}_{29}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 + \left(\frac{1}{4} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_5\right)c\,\hat{\mathbf{z}}$	(8*b*)	H III
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + \left(\frac{3}{4} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right)a\,\hat{\mathbf{y}} + \left(\frac{3}{4} - z_5\right)c\,\hat{\mathbf{z}}$	(8*b*)	H III
$\mathbf{B}_{31}$	=	$y_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$y_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(8*b*)	H III
$\mathbf{B}_{32}$	=	$-y_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$-y_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	(8*b*)	H III

$$\mathbf{B}_{33} = x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 = x_6 a\,\hat{\mathbf{x}} + y_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}} \quad (8b) \quad \text{H IV}$$

$$\mathbf{B}_{34} = -x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 = -x_6 a\,\hat{\mathbf{x}} - y_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right) c\,\hat{\mathbf{z}} \quad (8b) \quad \text{H IV}$$

$$\mathbf{B}_{35} = \left(\tfrac{1}{2} - y_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_6\right)\mathbf{a}_2 + \left(\tfrac{1}{4} + z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - y_6\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_6\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_6\right) c\,\hat{\mathbf{z}} \quad (8b) \quad \text{H IV}$$

$$\mathbf{B}_{36} = \left(\tfrac{1}{2} + y_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_2 + \left(\tfrac{3}{4} + z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + y_6\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_6\right) a\,\hat{\mathbf{y}} + \left(\tfrac{3}{4} + z_6\right) c\,\hat{\mathbf{z}} \quad (8b) \quad \text{H IV}$$

$$\mathbf{B}_{37} = \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{4} - z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_6\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - z_6\right) c\,\hat{\mathbf{z}} \quad (8b) \quad \text{H IV}$$

$$\mathbf{B}_{38} = \left(\tfrac{1}{2} + x_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_6\right)\mathbf{a}_2 + \left(\tfrac{3}{4} - z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_6\right) a\,\hat{\mathbf{y}} + \left(\tfrac{3}{4} - z_6\right) c\,\hat{\mathbf{z}} \quad (8b) \quad \text{H IV}$$

$$\mathbf{B}_{39} = y_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3 = y_6 a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}} \quad (8b) \quad \text{H IV}$$

$$\mathbf{B}_{40} = -y_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_3 = -y_6 a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_6\right) c\,\hat{\mathbf{z}} \quad (8b) \quad \text{H IV}$$

$$\mathbf{B}_{41} = x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 = x_7 a\,\hat{\mathbf{x}} + y_7 a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}} \quad (8b) \quad \text{H V}$$

$$\mathbf{B}_{42} = -x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_7\right)\mathbf{a}_3 = -x_7 a\,\hat{\mathbf{x}} - y_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_7\right) c\,\hat{\mathbf{z}} \quad (8b) \quad \text{H V}$$

$$\mathbf{B}_{43} = \left(\tfrac{1}{2} - y_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_7\right)\mathbf{a}_2 + \left(\tfrac{1}{4} + z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - y_7\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_7\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_7\right) c\,\hat{\mathbf{z}} \quad (8b) \quad \text{H V}$$

$$\mathbf{B}_{44} = \left(\tfrac{1}{2} + y_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_7\right)\mathbf{a}_2 + \left(\tfrac{3}{4} + z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + y_7\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_7\right) a\,\hat{\mathbf{y}} + \left(\tfrac{3}{4} + z_7\right) c\,\hat{\mathbf{z}} \quad (8b) \quad \text{H V}$$

$$\mathbf{B}_{45} = \left(\tfrac{1}{2} - x_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{4} - z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_7\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_7\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - z_7\right) c\,\hat{\mathbf{z}} \quad (8b) \quad \text{H V}$$

$$\mathbf{B}_{46} = \left(\tfrac{1}{2} + x_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_7\right)\mathbf{a}_2 + \left(\tfrac{3}{4} - z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_7\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_7\right) a\,\hat{\mathbf{y}} + \left(\tfrac{3}{4} - z_7\right) c\,\hat{\mathbf{z}} \quad (8b) \quad \text{H V}$$

$$\mathbf{B}_{47} = y_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3 = y_7 a\,\hat{\mathbf{x}} + x_7 a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}} \quad (8b) \quad \text{H V}$$

$$\mathbf{B}_{48} = -y_7\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_7\right)\mathbf{a}_3 = -y_7 a\,\hat{\mathbf{x}} - x_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_7\right) c\,\hat{\mathbf{z}} \quad (8b) \quad \text{H V}$$

$$\mathbf{B}_{49} = x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 = x_8 a\,\hat{\mathbf{x}} + y_8 a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}} \quad (8b) \quad \text{H VI}$$

$$\mathbf{B}_{50} = -x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_8\right)\mathbf{a}_3 = -x_8 a\,\hat{\mathbf{x}} - y_8 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_8\right) c\,\hat{\mathbf{z}} \quad (8b) \quad \text{H VI}$$

$$\mathbf{B}_{51} = \left(\tfrac{1}{2} - y_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_8\right)\mathbf{a}_2 + \left(\tfrac{1}{4} + z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - y_8\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_8\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_8\right) c\,\hat{\mathbf{z}} \quad (8b) \quad \text{H VI}$$

$$\mathbf{B}_{52} = \left(\tfrac{1}{2} + y_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_8\right)\mathbf{a}_2 + \left(\tfrac{3}{4} + z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + y_8\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_8\right) a\,\hat{\mathbf{y}} + \left(\tfrac{3}{4} + z_8\right) c\,\hat{\mathbf{z}} \quad (8b) \quad \text{H VI}$$

$$\mathbf{B}_{53} = \left(\tfrac{1}{2} - x_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_8\right)\mathbf{a}_2 + \left(\tfrac{1}{4} - z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_8\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_8\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - z_8\right) c\,\hat{\mathbf{z}} \quad (8b) \quad \text{H VI}$$

$$\mathbf{B}_{54} = \left(\tfrac{1}{2} + x_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_8\right)\mathbf{a}_2 + \left(\tfrac{3}{4} - z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_8\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_8\right) a\,\hat{\mathbf{y}} + \left(\tfrac{3}{4} - z_8\right) c\,\hat{\mathbf{z}} \quad (8b) \quad \text{H VI}$$

$$\mathbf{B}_{55} = y_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3 = y_8 a\,\hat{\mathbf{x}} + x_8 a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}} \quad (8b) \quad \text{H VI}$$

$$\mathbf{B}_{56} = -y_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_8\right)\mathbf{a}_3 = -y_8 a\,\hat{\mathbf{x}} - x_8 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_8\right) c\,\hat{\mathbf{z}} \quad (8b) \quad \text{H VI}$$

$$\mathbf{B}_{57} = x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 = x_9 a\,\hat{\mathbf{x}} + y_9 a\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O I}$$

$$\mathbf{B}_{58} = -x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_3 = -x_9 a\,\hat{\mathbf{x}} - y_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_9\right) c\,\hat{\mathbf{z}} \quad (8b) \quad \text{O I}$$

$\mathbf{B}_{59}$	=	$\left(\frac{1}{2}-y_9\right)\mathbf{a}_1+\left(\frac{1}{2}+x_9\right)\mathbf{a}_2+\left(\frac{1}{4}+z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_9\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_9\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+z_9\right)c\,\hat{\mathbf{z}}$	(8*b*)	O I
$\mathbf{B}_{60}$	=	$\left(\frac{1}{2}+y_9\right)\mathbf{a}_1+\left(\frac{1}{2}-x_9\right)\mathbf{a}_2+\left(\frac{3}{4}+z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_9\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_9\right)a\,\hat{\mathbf{y}}+\left(\frac{3}{4}+z_9\right)c\,\hat{\mathbf{z}}$	(8*b*)	O I
$\mathbf{B}_{61}$	=	$\left(\frac{1}{2}-x_9\right)\mathbf{a}_1+\left(\frac{1}{2}+y_9\right)\mathbf{a}_2+\left(\frac{1}{4}-z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_9\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_9\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-z_9\right)c\,\hat{\mathbf{z}}$	(8*b*)	O I
$\mathbf{B}_{62}$	=	$\left(\frac{1}{2}+x_9\right)\mathbf{a}_1+\left(\frac{1}{2}-y_9\right)\mathbf{a}_2+\left(\frac{3}{4}-z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_9\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_9\right)a\,\hat{\mathbf{y}}+\left(\frac{3}{4}-z_9\right)c\,\hat{\mathbf{z}}$	(8*b*)	O I
$\mathbf{B}_{63}$	=	$y_9\,\mathbf{a}_1+x_9\,\mathbf{a}_2-z_9\,\mathbf{a}_3$	=	$y_9a\,\hat{\mathbf{x}}+x_9a\,\hat{\mathbf{y}}-z_9c\,\hat{\mathbf{z}}$	(8*b*)	O I
$\mathbf{B}_{64}$	=	$-y_9\,\mathbf{a}_1-x_9\,\mathbf{a}_2+\left(\frac{1}{2}-z_9\right)\mathbf{a}_3$	=	$-y_9a\,\hat{\mathbf{x}}-x_9a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_9\right)c\,\hat{\mathbf{z}}$	(8*b*)	O I
$\mathbf{B}_{65}$	=	$x_{10}\,\mathbf{a}_1+y_{10}\,\mathbf{a}_2+z_{10}\,\mathbf{a}_3$	=	$x_{10}a\,\hat{\mathbf{x}}+y_{10}a\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}}$	(8*b*)	O II
$\mathbf{B}_{66}$	=	$-x_{10}\,\mathbf{a}_1-y_{10}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{10}\right)\mathbf{a}_3$	=	$-x_{10}a\,\hat{\mathbf{x}}-y_{10}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{10}\right)c\,\hat{\mathbf{z}}$	(8*b*)	O II
$\mathbf{B}_{67}$	=	$\left(\frac{1}{2}-y_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}+x_{10}\right)\mathbf{a}_2+\left(\frac{1}{4}+z_{10}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+z_{10}\right)c\,\hat{\mathbf{z}}$	(8*b*)	O II
$\mathbf{B}_{68}$	=	$\left(\frac{1}{2}+y_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}-x_{10}\right)\mathbf{a}_2+\left(\frac{3}{4}+z_{10}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{y}}+\left(\frac{3}{4}+z_{10}\right)c\,\hat{\mathbf{z}}$	(8*b*)	O II
$\mathbf{B}_{69}$	=	$\left(\frac{1}{2}-x_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{10}\right)\mathbf{a}_2+\left(\frac{1}{4}-z_{10}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{10}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-z_{10}\right)c\,\hat{\mathbf{z}}$	(8*b*)	O II
$\mathbf{B}_{70}$	=	$\left(\frac{1}{2}+x_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{10}\right)\mathbf{a}_2+\left(\frac{3}{4}-z_{10}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{10}\right)a\,\hat{\mathbf{y}}+\left(\frac{3}{4}-z_{10}\right)c\,\hat{\mathbf{z}}$	(8*b*)	O II
$\mathbf{B}_{71}$	=	$y_{10}\,\mathbf{a}_1+x_{10}\,\mathbf{a}_2-z_{10}\,\mathbf{a}_3$	=	$y_{10}a\,\hat{\mathbf{x}}+x_{10}a\,\hat{\mathbf{y}}-z_{10}c\,\hat{\mathbf{z}}$	(8*b*)	O II
$\mathbf{B}_{72}$	=	$-y_{10}\,\mathbf{a}_1-x_{10}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{10}\right)\mathbf{a}_3$	=	$-y_{10}a\,\hat{\mathbf{x}}-x_{10}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{10}\right)c\,\hat{\mathbf{z}}$	(8*b*)	O II
$\mathbf{B}_{73}$	=	$x_{11}\,\mathbf{a}_1+y_{11}\,\mathbf{a}_2+z_{11}\,\mathbf{a}_3$	=	$x_{11}a\,\hat{\mathbf{x}}+y_{11}a\,\hat{\mathbf{y}}+z_{11}c\,\hat{\mathbf{z}}$	(8*b*)	O III
$\mathbf{B}_{74}$	=	$-x_{11}\,\mathbf{a}_1-y_{11}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{11}\right)\mathbf{a}_3$	=	$-x_{11}a\,\hat{\mathbf{x}}-y_{11}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{11}\right)c\,\hat{\mathbf{z}}$	(8*b*)	O III
$\mathbf{B}_{75}$	=	$\left(\frac{1}{2}-y_{11}\right)\mathbf{a}_1+\left(\frac{1}{2}+x_{11}\right)\mathbf{a}_2+\left(\frac{1}{4}+z_{11}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_{11}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+z_{11}\right)c\,\hat{\mathbf{z}}$	(8*b*)	O III
$\mathbf{B}_{76}$	=	$\left(\frac{1}{2}+y_{11}\right)\mathbf{a}_1+\left(\frac{1}{2}-x_{11}\right)\mathbf{a}_2+\left(\frac{3}{4}+z_{11}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_{11}\right)a\,\hat{\mathbf{y}}+\left(\frac{3}{4}+z_{11}\right)c\,\hat{\mathbf{z}}$	(8*b*)	O III
$\mathbf{B}_{77}$	=	$\left(\frac{1}{2}-x_{11}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{11}\right)\mathbf{a}_2+\left(\frac{1}{4}-z_{11}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{11}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-z_{11}\right)c\,\hat{\mathbf{z}}$	(8*b*)	O III
$\mathbf{B}_{78}$	=	$\left(\frac{1}{2}+x_{11}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{11}\right)\mathbf{a}_2+\left(\frac{3}{4}-z_{11}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{11}\right)a\,\hat{\mathbf{y}}+\left(\frac{3}{4}-z_{11}\right)c\,\hat{\mathbf{z}}$	(8*b*)	O III
$\mathbf{B}_{79}$	=	$y_{11}\,\mathbf{a}_1+x_{11}\,\mathbf{a}_2-z_{11}\,\mathbf{a}_3$	=	$y_{11}a\,\hat{\mathbf{x}}+x_{11}a\,\hat{\mathbf{y}}-z_{11}c\,\hat{\mathbf{z}}$	(8*b*)	O III
$\mathbf{B}_{80}$	=	$-y_{11}\,\mathbf{a}_1-x_{11}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{11}\right)\mathbf{a}_3$	=	$-y_{11}a\,\hat{\mathbf{x}}-x_{11}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{11}\right)c\,\hat{\mathbf{z}}$	(8*b*)	O III
$\mathbf{B}_{81}$	=	$x_{12}\,\mathbf{a}_1+y_{12}\,\mathbf{a}_2+z_{12}\,\mathbf{a}_3$	=	$x_{12}a\,\hat{\mathbf{x}}+y_{12}a\,\hat{\mathbf{y}}+z_{12}c\,\hat{\mathbf{z}}$	(8*b*)	O IV
$\mathbf{B}_{82}$	=	$-x_{12}\,\mathbf{a}_1-y_{12}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{12}\right)\mathbf{a}_3$	=	$-x_{12}a\,\hat{\mathbf{x}}-y_{12}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{12}\right)c\,\hat{\mathbf{z}}$	(8*b*)	O IV
$\mathbf{B}_{83}$	=	$\left(\frac{1}{2}-y_{12}\right)\mathbf{a}_1+\left(\frac{1}{2}+x_{12}\right)\mathbf{a}_2+\left(\frac{1}{4}+z_{12}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_{12}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_{12}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+z_{12}\right)c\,\hat{\mathbf{z}}$	(8*b*)	O IV

$\mathbf{B}_{84}$	=	$\left(\frac{1}{2}+y_{12}\right)\mathbf{a}_1+\left(\frac{1}{2}-x_{12}\right)\mathbf{a}_2+\left(\frac{3}{4}+z_{12}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_{12}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_{12}\right)a\,\hat{\mathbf{y}}+\left(\frac{3}{4}+z_{12}\right)c\,\hat{\mathbf{z}}$	(8*b*)	O IV
$\mathbf{B}_{85}$	=	$\left(\frac{1}{2}-x_{12}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{12}\right)\mathbf{a}_2+\left(\frac{1}{4}-z_{12}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{12}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{12}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-z_{12}\right)c\,\hat{\mathbf{z}}$	(8*b*)	O IV
$\mathbf{B}_{86}$	=	$\left(\frac{1}{2}+x_{12}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{12}\right)\mathbf{a}_2+\left(\frac{3}{4}-z_{12}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{12}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{12}\right)a\,\hat{\mathbf{y}}+\left(\frac{3}{4}-z_{12}\right)c\,\hat{\mathbf{z}}$	(8*b*)	O IV
$\mathbf{B}_{87}$	=	$y_{12}\,\mathbf{a}_1+x_{12}\,\mathbf{a}_2-z_{12}\,\mathbf{a}_3$	=	$y_{12}a\,\hat{\mathbf{x}}+x_{12}a\,\hat{\mathbf{y}}-z_{12}c\,\hat{\mathbf{z}}$	(8*b*)	O IV
$\mathbf{B}_{88}$	=	$-y_{12}\,\mathbf{a}_1-x_{12}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{12}\right)\mathbf{a}_3$	=	$-y_{12}a\,\hat{\mathbf{x}}-x_{12}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{12}\right)c\,\hat{\mathbf{z}}$	(8*b*)	O IV
$\mathbf{B}_{89}$	=	$x_{13}\,\mathbf{a}_1+y_{13}\,\mathbf{a}_2+z_{13}\,\mathbf{a}_3$	=	$x_{13}a\,\hat{\mathbf{x}}+y_{13}a\,\hat{\mathbf{y}}+z_{13}c\,\hat{\mathbf{z}}$	(8*b*)	O V
$\mathbf{B}_{90}$	=	$-x_{13}\,\mathbf{a}_1-y_{13}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{13}\right)\mathbf{a}_3$	=	$-x_{13}a\,\hat{\mathbf{x}}-y_{13}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{13}\right)c\,\hat{\mathbf{z}}$	(8*b*)	O V
$\mathbf{B}_{91}$	=	$\left(\frac{1}{2}-y_{13}\right)\mathbf{a}_1+\left(\frac{1}{2}+x_{13}\right)\mathbf{a}_2+\left(\frac{1}{4}+z_{13}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_{13}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_{13}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+z_{13}\right)c\,\hat{\mathbf{z}}$	(8*b*)	O V
$\mathbf{B}_{92}$	=	$\left(\frac{1}{2}+y_{13}\right)\mathbf{a}_1+\left(\frac{1}{2}-x_{13}\right)\mathbf{a}_2+\left(\frac{3}{4}+z_{13}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_{13}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_{13}\right)a\,\hat{\mathbf{y}}+\left(\frac{3}{4}+z_{13}\right)c\,\hat{\mathbf{z}}$	(8*b*)	O V
$\mathbf{B}_{93}$	=	$\left(\frac{1}{2}-x_{13}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{13}\right)\mathbf{a}_2+\left(\frac{1}{4}-z_{13}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{13}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{13}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-z_{13}\right)c\,\hat{\mathbf{z}}$	(8*b*)	O V
$\mathbf{B}_{94}$	=	$\left(\frac{1}{2}+x_{13}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{13}\right)\mathbf{a}_2+\left(\frac{3}{4}-z_{13}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{13}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{13}\right)a\,\hat{\mathbf{y}}+\left(\frac{3}{4}-z_{13}\right)c\,\hat{\mathbf{z}}$	(8*b*)	O V
$\mathbf{B}_{95}$	=	$y_{13}\,\mathbf{a}_1+x_{13}\,\mathbf{a}_2-z_{13}\,\mathbf{a}_3$	=	$y_{13}a\,\hat{\mathbf{x}}+x_{13}a\,\hat{\mathbf{y}}-z_{13}c\,\hat{\mathbf{z}}$	(8*b*)	O V
$\mathbf{B}_{96}$	=	$-y_{13}\,\mathbf{a}_1-x_{13}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{13}\right)\mathbf{a}_3$	=	$-y_{13}a\,\hat{\mathbf{x}}-x_{13}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{13}\right)c\,\hat{\mathbf{z}}$	(8*b*)	O V

References:

- K. Stadnicka, A. M. Glazer, and M. Koralewski, *Structure, absolute configuration and optical activity of α-nickel sulfate hexahydrate*, Acta Crystallogr. Sect. B Struct. Sci. **43**, 319–325 (1987), doi:10.1107/S0108768187097787.
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Found in:

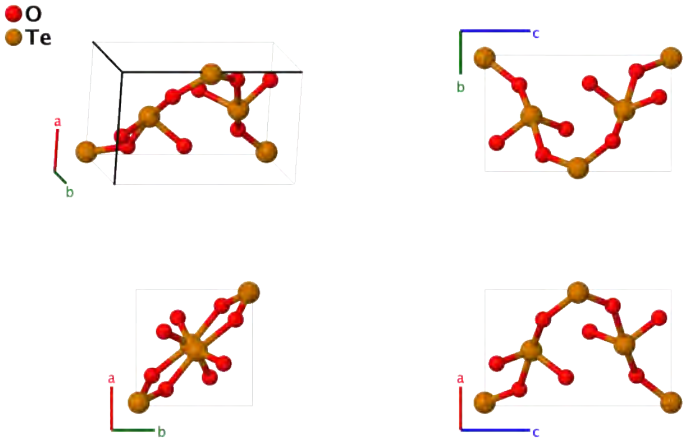
- J. W. Anthony, R. A. Bideaux, K. W. Bladh, and M. C. Nichols, eds., *Handbook of Mineralogy* (Mineralogical Society of America, 2004), chap. Retgersite, $NiSO_4{\cdot}6H_2O$.

Geometry files:

- CIF: pp. 1676
- POSCAR: pp. 1676

Paratellurite (α-TeO_2) Structure: A2B_tP12_92_b_a

http://aflow.org/prototype-encyclopedia/A2B_tP12_92_b_a.TeO2

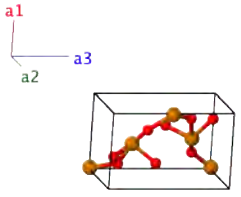


Prototype	:	O_2Te
AFLOW prototype label	:	A2B_tP12_92_b_a
***Strukturbericht* designation**	:	None
Pearson symbol	:	tP12
Space group number	:	92
Space group symbol	:	$P4_12_12$
AFLOW prototype command	:	`aflow --proto=A2B_tP12_92_b_a --params=`$a, c/a, x_1, x_2, y_2, z_2$

- Although this as the same space group and Wyckoff positions as α-cristobalite ($C30$), the actual positions of the atoms are substantially different, so we have given the two compounds separate entries in the encyclopedia.

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + x_1\,\mathbf{a}_2$	=	$x_1 a\,\hat{\mathbf{x}} + x_1 a\,\hat{\mathbf{y}}$	($4a$)	Te
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 - x_1\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} - x_1 a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4a$)	Te
$\mathbf{B}_3$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_1\right)\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_1\right)a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($4a$)	Te
$\mathbf{B}_4$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_1\right)\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_1\right)a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	($4a$)	Te
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + y_2 a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	($8b$)	O

$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} - y_2 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right) c\,\hat{\mathbf{z}}$	(8*b*)	O
$\mathbf{B}_7$	=	$\left(\frac{1}{2} - y_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_2 + \left(\frac{1}{4} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_2\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_2\right) c\,\hat{\mathbf{z}}$	(8*b*)	O
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + y_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_2 + \left(\frac{3}{4} + z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_2\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{y}} + \left(\frac{3}{4} + z_2\right) c\,\hat{\mathbf{z}}$	(8*b*)	O
$\mathbf{B}_9$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_2\right)\mathbf{a}_2 + \left(\frac{1}{4} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_2\right) c\,\hat{\mathbf{z}}$	(8*b*)	O
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_2\right)\mathbf{a}_2 + \left(\frac{3}{4} - z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right) a\,\hat{\mathbf{y}} + \left(\frac{3}{4} - z_2\right) c\,\hat{\mathbf{z}}$	(8*b*)	O
$\mathbf{B}_{11}$	=	$y_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$y_2 a\,\hat{\mathbf{x}} + x_2 a\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(8*b*)	O
$\mathbf{B}_{12}$	=	$-y_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$-y_2 a\,\hat{\mathbf{x}} - x_2 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right) c\,\hat{\mathbf{z}}$	(8*b*)	O

References:

- P. A. Thomas, *The crystal structure and absolute optical chirality of paratellurite, α-TeO_2*, J. Phys. C: Solid State Phys. **21**, 4611–4627 (1988), doi:10.1088/0022-3719/21/25/009.

Found in:

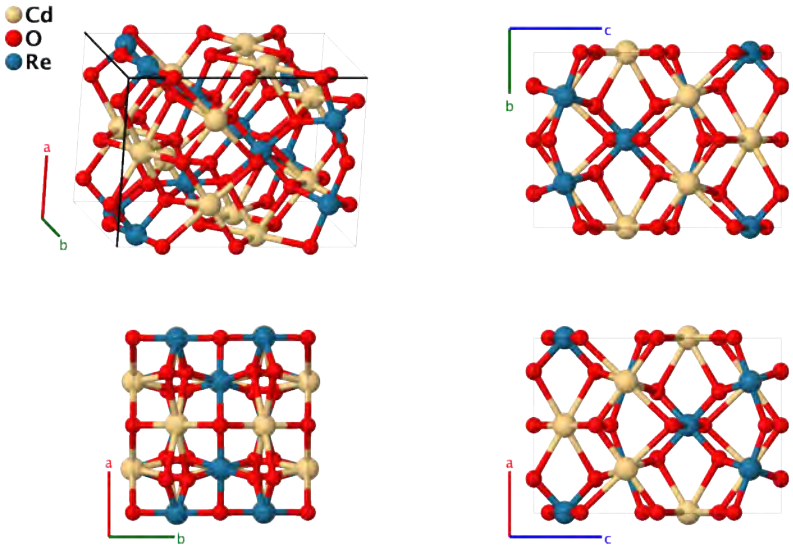
- M. Ceriotti, F. Pietrucci, and M. Bernasconi, Ab initio *study of the vibrational properties of crystalline TeO_2: The α, β, and γ phases*, Phys. Rev. B **73**, 104304 (2006), doi:10.1103/PhysRevB.73.104304.

Geometry files:

- CIF: pp. 1677
- POSCAR: pp. 1677

Phase III $Cd_2Re_2O_7$ Structure: A2B7C2_tI44_98_f_bcde_f

http://aflow.org/prototype-encyclopedia/A2B7C2_tI44_98_f_bcde_f

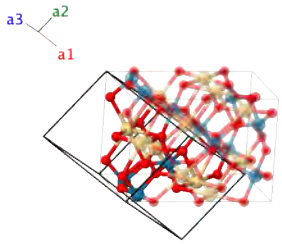


Prototype	:	$Cd_2O_7Re_2$
AFLOW prototype label	:	A2B7C2_tI44_98_f_bcde_f
***Strukturbericht* designation**	:	None
Pearson symbol	:	tI44
Space group number	:	98
Space group symbol	:	$I4_122$
AFLOW prototype command	:	`aflow --proto=A2B7C2_tI44_98_f_bcde_f --params=`$a, c/a, z_2, x_3, x_4, x_5, x_6$

- $Cd_2Re_2O_7$ exhibits a number of phases. We will use the notation of (Kapcia, 2019) to describe them:
- – Phase I: above 200 K, the system takes on the cubic pyrochlore ($E8_1$) structure.
  – Phase II: in the range 120-200 K the system is in the tetragonal $I\bar{4}m2$ #119 structure.
  – Phase III: in the range 80-120 K the system is in the tetragonal $I4_122$ #98 structure (this structure).
  – Phase IV: (Kapcia, 2019) did a first-principles study of this system and found that below 80 K Phase III develops a soft phonon mode which transforms the system into an orthorhombic $F222$ #22 structure.
  – (Norman, 2019) points out that both Phase III and Phase IV structures have issues.
- Both Phase II and Phase III are extremely close to Phase I. If AFLOW-SYM and FINDSYM allow a 0.2 Å uncertainty in lattice vectors and atomic positions both of the tetragonal phases become cubic.
- Phase IV is extremely close to Phase II. If AFLOW-SYM and FINDSYM allow a 0.002 Å uncertainty in the lattice vectors and atomic positions the orthorhombic phase becomes tetragonal.
- Data for the Phase III structure was taken at 90 K.

Body-centered Tetragonal primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= -\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}\\ \mathbf{a}_2 &= \tfrac{1}{2}a\,\hat{\mathbf{x}} - \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}\\ \mathbf{a}_3 &= \tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} - \tfrac{1}{2}c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{2}\mathbf{a}_1 + \frac{1}{2}\mathbf{a}_2$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	(4*b*)	O I
$\mathbf{B}_2$	=	$\frac{1}{4}\mathbf{a}_1 + \frac{3}{4}\mathbf{a}_2 + \frac{1}{2}\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*b*)	O I
$\mathbf{B}_3$	=	$z_2\,\mathbf{a}_1 + z_2\,\mathbf{a}_2$	=	$z_2 c\,\hat{\mathbf{z}}$	(8*c*)	O II
$\mathbf{B}_4$	=	$\left(\frac{3}{4}+z_2\right)\mathbf{a}_1 + \left(\frac{1}{4}+z_2\right)\mathbf{a}_2 + \frac{1}{2}\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \left(\frac{1}{4}+z_2\right)c\,\hat{\mathbf{z}}$	(8*c*)	O II
$\mathbf{B}_5$	=	$\left(\frac{3}{4}-z_2\right)\mathbf{a}_1 + \left(\frac{1}{4}-z_2\right)\mathbf{a}_2 + \frac{1}{2}\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \left(\frac{1}{4}-z_2\right)c\,\hat{\mathbf{z}}$	(8*c*)	O II
$\mathbf{B}_6$	=	$-z_2\,\mathbf{a}_1 - z_2\,\mathbf{a}_2$	=	$-z_2 c\,\hat{\mathbf{z}}$	(8*c*)	O II
$\mathbf{B}_7$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + 2x_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}}$	(8*d*)	O III
$\mathbf{B}_8$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - 2x_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}}$	(8*d*)	O III
$\mathbf{B}_9$	=	$\left(\frac{3}{4}+x_3\right)\mathbf{a}_1 + \left(\frac{1}{4}-x_3\right)\mathbf{a}_2 + \frac{1}{2}\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{10}$	=	$\left(\frac{3}{4}-x_3\right)\mathbf{a}_1 + \left(\frac{1}{4}+x_3\right)\mathbf{a}_2 + \frac{1}{2}\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(8*d*)	O III
$\mathbf{B}_{11}$	=	$x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2$	=	$-x_4 a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}}$	(8*e*)	O IV
$\mathbf{B}_{12}$	=	$-x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2$	=	$x_4 a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}}$	(8*e*)	O IV
$\mathbf{B}_{13}$	=	$\left(\frac{3}{4}-x_4\right)\mathbf{a}_1 + \left(\frac{1}{4}-x_4\right)\mathbf{a}_2 + \left(\frac{1}{2}-2x_4\right)\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(8*e*)	O IV
$\mathbf{B}_{14}$	=	$\left(\frac{3}{4}+x_4\right)\mathbf{a}_1 + \left(\frac{1}{4}+x_4\right)\mathbf{a}_2 + \left(\frac{1}{2}+2x_4\right)\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(8*e*)	O IV
$\mathbf{B}_{15}$	=	$\frac{3}{8}\mathbf{a}_1 + \left(\frac{1}{8}+x_5\right)\mathbf{a}_2 + \left(\frac{1}{4}+x_5\right)\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{8}c\,\hat{\mathbf{z}}$	(8*f*)	Cd
$\mathbf{B}_{16}$	=	$\frac{7}{8}\mathbf{a}_1 + \left(\frac{1}{8}-x_5\right)\mathbf{a}_2 + \left(\frac{3}{4}-x_5\right)\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{1}{8}c\,\hat{\mathbf{z}}$	(8*f*)	Cd
$\mathbf{B}_{17}$	=	$\left(\frac{7}{8}+x_5\right)\mathbf{a}_1 + \frac{1}{8}\mathbf{a}_2 + \left(\frac{1}{4}+x_5\right)\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{y}} + \frac{3}{8}c\,\hat{\mathbf{z}}$	(8*f*)	Cd
$\mathbf{B}_{18}$	=	$\left(\frac{7}{8}-x_5\right)\mathbf{a}_1 + \frac{5}{8}\mathbf{a}_2 + \left(\frac{3}{4}-x_5\right)\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{y}} + \frac{3}{8}c\,\hat{\mathbf{z}}$	(8*f*)	Cd
$\mathbf{B}_{19}$	=	$\frac{3}{8}\mathbf{a}_1 + \left(\frac{1}{8}+x_6\right)\mathbf{a}_2 + \left(\frac{1}{4}+x_6\right)\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{8}c\,\hat{\mathbf{z}}$	(8*f*)	Re
$\mathbf{B}_{20}$	=	$\frac{7}{8}\mathbf{a}_1 + \left(\frac{1}{8}-x_6\right)\mathbf{a}_2 + \left(\frac{3}{4}-x_6\right)\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{1}{8}c\,\hat{\mathbf{z}}$	(8*f*)	Re
$\mathbf{B}_{21}$	=	$\left(\frac{7}{8}+x_6\right)\mathbf{a}_1 + \frac{1}{8}\mathbf{a}_2 + \left(\frac{1}{4}+x_6\right)\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{y}} + \frac{3}{8}c\,\hat{\mathbf{z}}$	(8*f*)	Re
$\mathbf{B}_{22}$	=	$\left(\frac{7}{8}-x_6\right)\mathbf{a}_1 + \frac{5}{8}\mathbf{a}_2 + \left(\frac{3}{4}-x_6\right)\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{y}} + \frac{3}{8}c\,\hat{\mathbf{z}}$	(8*f*)	Re

References:

- S.-W. Huang, H.-T. Jeng, J.-Y. Lin, W. J. Chang, J. M. Chen, G. H. Lee, H. Berger, H. D. Yang, and K. S. Liang, *Electronic structure of pyrochlore $Cd_2Re_2O_7$*, J. Phys.: Condens. Matter **21**, 195602 (2009), doi:10.1088/0953-8984/21/19/195602.
- K. J. Kapcia, M. Reedyk, M. Hajialamdari, A. Ptok, P. Piekarz, F. S. Razavi, A. M. Oleś, and R. K. Kremer, *Discovery of a low-temperature orthorhombic phase of the $Cd_2Re_2O_7$ superconductor*, Phys. Rev. Research **2**, 033108 (2020), doi:10.1103/PhysRevResearch.2.033108.

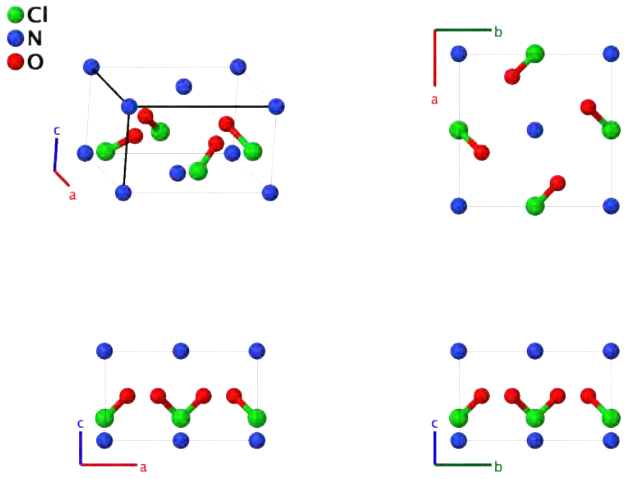
Found in:

- M. R. Norman, *The crystal structure of the inversion breaking metal* $Cd_2Re_2O_7$, Phys. Rev. B **101**, 045117 (2020), doi:10.1103/PhysRevB.101.045117.

Geometry files:
- CIF: pp. 1677
- POSCAR: pp. 1678

$F5_4$ (NH_4ClO_2) (*obsolete*) Structure: ABC2_tP8_100_b_a_c

http://aflow.org/prototype-encyclopedia/ABC2_tP8_100_b_a_c

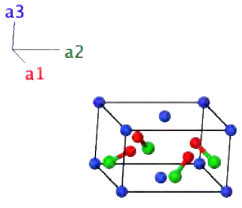


Prototype	:	$Cl(NH_4)O_2$
AFLOW prototype label	:	ABC2_tP8_100_b_a_c
***Strukturbericht* designation**	:	$F5_4$
Pearson symbol	:	tP8
Space group number	:	100
Space group symbol	:	*P*4*bm*
AFLOW prototype command	:	`aflow --proto=ABC2_tP8_100_b_a_c --params=`$a, c/a, z_1, z_2, x_3, z_3$

- (Levi, 1931) first detemined this structure, but they were actually looking at a combination of ammonium chlorite, NH_4ClO_2 and ammonium chlorate, NH_4ClO_3, which takes on the γ-KNO_3 structure. In addition, they were unable to determine the positions of the hydrogen atoms, and it appears that space group *P*4*bm* #100 is incompatible with having four hydrogen atoms in a tetrahedral arrangement about the nitrogen atom. (Smolentsev, 2005) determined the positions of the hydrogen atoms, and placed this structure in space group $P\bar{4}2_1m$ #113, making the structure of (Levi, 1931) obsolete. We present it here for historical interest.
- Since the position of the hydrogen atoms in the NH_4 ions were not determined, we only provide the nitrogen atom positions (labeled as NH_4).

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$z_1\,\mathbf{a}_3$	=	$z_1 c\,\hat{\mathbf{z}}$	($2a$)	NH_4
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	($2a$)	NH_4
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_1 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + z_2 c\,\hat{\mathbf{z}}$	($2b$)	Cl
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	($2b$)	Cl
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	($4c$)	O
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	($4c$)	O
$\mathbf{B}_7$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	($4c$)	O
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 - x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	($4c$)	O

References:

- G. R. Levi and A. Scherillo, *Ricerche cristallografiche sui sali dell'acido cloroso*, Zeitschrift für Kristallographie - Crystalline Materials **76**, 431–452 (1931), doi:10.1524/zkri.1931.76.1.431.
- A. I. Smolentsev and D. Y. Naumov, *Ammonium chlorite, NH_4ClO_2, at 150 K*, Acta Crystallogr. E **61**, i38–i40 (2005), doi:10.1107/S1600536805005088.

Found in:

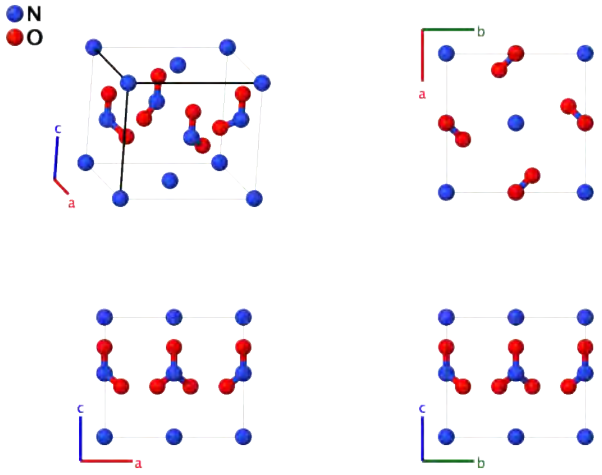
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1678
- POSCAR: pp. 1678

NH_4NO_3 II ($G0_9$) Structure: ABC3_tP10_100_b_a_bc

http://aflow.org/prototype-encyclopedia/ABC3_tP10_100_b_a_bc



Prototype	:	$N(NH_4)O_3$
AFLOW prototype label	:	ABC3_tP10_100_b_a_bc
***Strukturbericht* designation**	:	$G0_9$
Pearson symbol	:	tP10
Space group number	:	100
Space group symbol	:	*P*4*bm*
AFLOW prototype command	:	`aflow --proto=ABC3_tP10_100_b_a_bc --params=`$a, c/a, z_1, z_2, z_3, x_4, z_4$

- Ammonium Nitrate exists in a variety of forms, (Hermann, 1937) depending on the temperature:

Phase	Temperature °C	*Strukturbericht*	Page	
I	125 – 170	$G0_8$	AB_cP2_221_a_b.NH4.NO3	
II	84 – 125	$G0_9$	ABC3_tP10_100_b_a_bc	(this structure)
III	32 – 84	$G0_{10}$	ABC3_oP20_62_c_c_cd.N.NH4.O	
IV	-17 – 32	$G0_{11}$	A4B2C3_oP18_59_ef_ab_af	
V	< −17	Gwihabaite	A4B2C3_tP72_77_8d_ab2c2d_6d2	

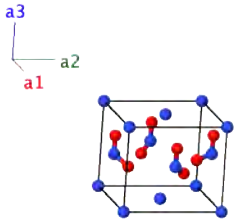
- Data for this structure was taken at 60 °C.
- The positions of the hydrogen atoms were not determined. The isolated nitrogen atoms in this structure's visualization are surrounded by four hydrogen atoms in an approximately tetrahedral arrangement. It is likely that the NH_4 ions are free to rotate (Kracek, 1937).
- Both (Shinnaka, 1956) and (Hermann, 1937) state that the available x-ray diffraction data supports a space group of either *P*4*bm* (#100) or $P\bar{4}2_1m$ (#113). The atomic positions found by Shinnaka agree with space group *P*4*bm*.
- (Shinnaka, 1956) states that the NO_3 nitrate groups are rotating, but this rotation "is almost bound in two orientations (in opposite directions)." He then gives two possible orientations for the nitrate. We present the first orientation here. The second orientation is obtained by taking $z_3 \rightarrow -z_3$ and $z_4 \rightarrow -z_4$.
- Another way of presenting this information would be to add a second nitrate group to the primitive cell, and set the occupation of all the atoms in the nitrates at 50%. This would give a structure in space group $P4/mbm$ (#127), which

might be useful as a pictorial representation but does not correctly represent the physics of the crystal, as the nitrogen and oxygen atoms in an individual nitrate ion must remain together.

- The N–O distances in this structure are about 10% smaller than the distances found in the other phases of NH_4NO_3. This suggests that the structure should be reevaluated.
- The positions of the hydrogen atoms in the NH_4 ion were not determined, so we only provide the positions of the nitrogen atoms (labeled as NH_4).

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$z_1\,\mathbf{a}_3$	=	$z_1 c\,\hat{\mathbf{z}}$	($2a$)	NH_4
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	($2a$)	NH_4
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_1 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + z_2 c\,\hat{\mathbf{z}}$	($2b$)	N
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	($2b$)	N
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_1 + z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + z_3 c\,\hat{\mathbf{z}}$	($2b$)	O I
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	($2b$)	O I
$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + x_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	($4c$)	O II
$\mathbf{B}_8$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	($4c$)	O II
$\mathbf{B}_9$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	($4c$)	O II
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 - x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	($4c$)	O II

References:

- Y. Shinnaka, *On the Metastable Transition and the Crystal Structure of Ammonium Nitrate (Tetragonal Modification)*, J. Phys. Soc. Jpn. **11**, 393–396 (1956), doi:10.1143/JPSJ.11.393.
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).
- F. C. Kracek, S. B. Hendricks, and E. Posnjak, *Group Rotation in Solid Ammonium and Calcium Nitrates*, Nature **128**, 410–411 (1931), doi:10.1038/128410b0.

Found in:

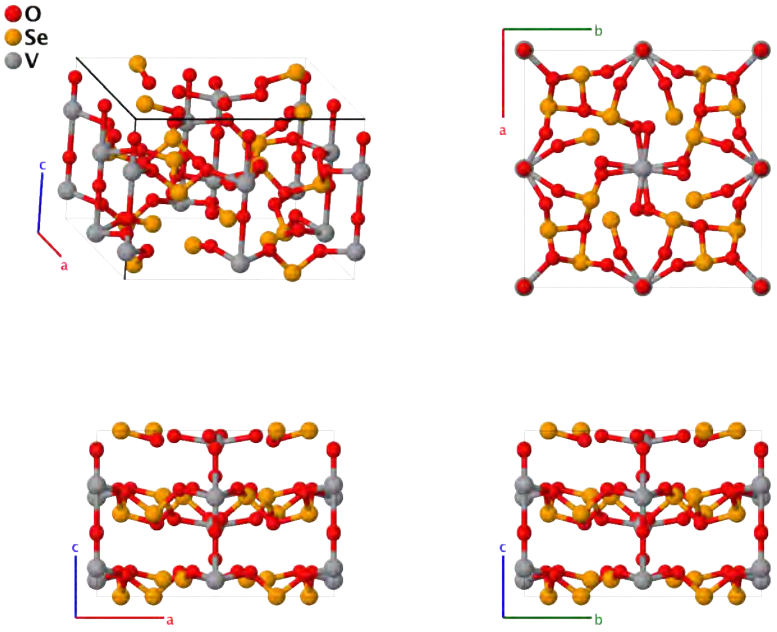
- C. S. Choi, J. E. Mapes, and E. Prince, *The structure of ammonium nitrate (IV)*, Acta Crystallogr. Sect. B Struct. Sci. **28**, 1357–1361 (1972), doi:10.1107/S0567740872004303.

Geometry files:

- CIF: pp. 1678
- POSCAR: pp. 1679

VSe_2O_6 Structure: A6B2C_tP72_103_abc5d_2d_abc

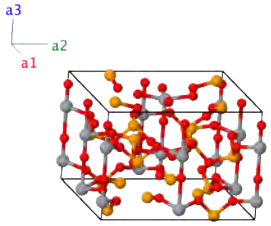
http://aflow.org/prototype-encyclopedia/A6B2C_tP72_103_abc5d_2d_abc



Prototype	:	O_6Se_2V
AFLOW prototype label	:	A6B2C_tP72_103_abc5d_2d_abc
***Strukturbericht* designation**	:	None
Pearson symbol	:	tP72
Space group number	:	103
Space group symbol	:	*P*4*cc*
AFLOW prototype command	:	`aflow --proto=A6B2C_tP72_103_abc5d_2d_abc --params=`$a, c/a, z_1, z_2, z_3, z_4, z_5, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}$

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$z_1\,\mathbf{a}_3$	=	$z_1 c\,\hat{\mathbf{z}}$	(2*a*)	O I

$\mathbf{B}_2$	=	$\left(\frac{1}{2}+z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+z_1\right)c\,\hat{\mathbf{z}}$	(2*a*)	O I
$\mathbf{B}_3$	=	$z_2\,\mathbf{a}_3$	=	$z_2 c\,\hat{\mathbf{z}}$	(2*a*)	V I
$\mathbf{B}_4$	=	$\left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+z_2\right)c\,\hat{\mathbf{z}}$	(2*a*)	V I
$\mathbf{B}_5$	=	$\frac{1}{2}\mathbf{a}_1+\frac{1}{2}\mathbf{a}_2+z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}+\frac{1}{2}a\,\hat{\mathbf{y}}+z_3 c\,\hat{\mathbf{z}}$	(2*b*)	O II
$\mathbf{B}_6$	=	$\frac{1}{2}\mathbf{a}_1+\frac{1}{2}\mathbf{a}_2+\left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}+\frac{1}{2}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_3\right)c\,\hat{\mathbf{z}}$	(2*b*)	O II
$\mathbf{B}_7$	=	$\frac{1}{2}\mathbf{a}_1+\frac{1}{2}\mathbf{a}_2+z_4\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}+\frac{1}{2}a\,\hat{\mathbf{y}}+z_4 c\,\hat{\mathbf{z}}$	(2*b*)	V II
$\mathbf{B}_8$	=	$\frac{1}{2}\mathbf{a}_1+\frac{1}{2}\mathbf{a}_2+\left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}+\frac{1}{2}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_4\right)c\,\hat{\mathbf{z}}$	(2*b*)	V II
$\mathbf{B}_9$	=	$\frac{1}{2}\mathbf{a}_2+z_5\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}}+z_5 c\,\hat{\mathbf{z}}$	(4*c*)	O III
$\mathbf{B}_{10}$	=	$\frac{1}{2}\mathbf{a}_1+z_5\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}+z_5 c\,\hat{\mathbf{z}}$	(4*c*)	O III
$\mathbf{B}_{11}$	=	$\frac{1}{2}\mathbf{a}_2+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	(4*c*)	O III
$\mathbf{B}_{12}$	=	$\frac{1}{2}\mathbf{a}_1+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	(4*c*)	O III
$\mathbf{B}_{13}$	=	$\frac{1}{2}\mathbf{a}_2+z_6\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}}+z_6 c\,\hat{\mathbf{z}}$	(4*c*)	V III
$\mathbf{B}_{14}$	=	$\frac{1}{2}\mathbf{a}_1+z_6\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}+z_6 c\,\hat{\mathbf{z}}$	(4*c*)	V III
$\mathbf{B}_{15}$	=	$\frac{1}{2}\mathbf{a}_2+\left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	(4*c*)	V III
$\mathbf{B}_{16}$	=	$\frac{1}{2}\mathbf{a}_1+\left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	(4*c*)	V III
$\mathbf{B}_{17}$	=	$x_7\,\mathbf{a}_1+y_7\,\mathbf{a}_2+z_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}}+y_7 a\,\hat{\mathbf{y}}+z_7 c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{18}$	=	$-x_7\,\mathbf{a}_1-y_7\,\mathbf{a}_2+z_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}}-y_7 a\,\hat{\mathbf{y}}+z_7 c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{19}$	=	$-y_7\,\mathbf{a}_1+x_7\,\mathbf{a}_2+z_7\,\mathbf{a}_3$	=	$-y_7 a\,\hat{\mathbf{x}}+x_7 a\,\hat{\mathbf{y}}+z_7 c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{20}$	=	$y_7\,\mathbf{a}_1-x_7\,\mathbf{a}_2+z_7\,\mathbf{a}_3$	=	$y_7 a\,\hat{\mathbf{x}}-x_7 a\,\hat{\mathbf{y}}+z_7 c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{21}$	=	$x_7\,\mathbf{a}_1-y_7\,\mathbf{a}_2+\left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}}-y_7 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{22}$	=	$-x_7\,\mathbf{a}_1+y_7\,\mathbf{a}_2+\left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}}+y_7 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{23}$	=	$-y_7\,\mathbf{a}_1-x_7\,\mathbf{a}_2+\left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	=	$-y_7 a\,\hat{\mathbf{x}}-x_7 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{24}$	=	$y_7\,\mathbf{a}_1+x_7\,\mathbf{a}_2+\left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	=	$y_7 a\,\hat{\mathbf{x}}+x_7 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	(8*d*)	O IV
$\mathbf{B}_{25}$	=	$x_8\,\mathbf{a}_1+y_8\,\mathbf{a}_2+z_8\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}}+y_8 a\,\hat{\mathbf{y}}+z_8 c\,\hat{\mathbf{z}}$	(8*d*)	O V
$\mathbf{B}_{26}$	=	$-x_8\,\mathbf{a}_1-y_8\,\mathbf{a}_2+z_8\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}}-y_8 a\,\hat{\mathbf{y}}+z_8 c\,\hat{\mathbf{z}}$	(8*d*)	O V
$\mathbf{B}_{27}$	=	$-y_8\,\mathbf{a}_1+x_8\,\mathbf{a}_2+z_8\,\mathbf{a}_3$	=	$-y_8 a\,\hat{\mathbf{x}}+x_8 a\,\hat{\mathbf{y}}+z_8 c\,\hat{\mathbf{z}}$	(8*d*)	O V
$\mathbf{B}_{28}$	=	$y_8\,\mathbf{a}_1-x_8\,\mathbf{a}_2+z_8\,\mathbf{a}_3$	=	$y_8 a\,\hat{\mathbf{x}}-x_8 a\,\hat{\mathbf{y}}+z_8 c\,\hat{\mathbf{z}}$	(8*d*)	O V
$\mathbf{B}_{29}$	=	$x_8\,\mathbf{a}_1-y_8\,\mathbf{a}_2+\left(\frac{1}{2}+z_8\right)\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}}-y_8 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_8\right)c\,\hat{\mathbf{z}}$	(8*d*)	O V
$\mathbf{B}_{30}$	=	$-x_8\,\mathbf{a}_1+y_8\,\mathbf{a}_2+\left(\frac{1}{2}+z_8\right)\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}}+y_8 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_8\right)c\,\hat{\mathbf{z}}$	(8*d*)	O V
$\mathbf{B}_{31}$	=	$-y_8\,\mathbf{a}_1-x_8\,\mathbf{a}_2+\left(\frac{1}{2}+z_8\right)\mathbf{a}_3$	=	$-y_8 a\,\hat{\mathbf{x}}-x_8 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_8\right)c\,\hat{\mathbf{z}}$	(8*d*)	O V
$\mathbf{B}_{32}$	=	$y_8\,\mathbf{a}_1+x_8\,\mathbf{a}_2+\left(\frac{1}{2}+z_8\right)\mathbf{a}_3$	=	$y_8 a\,\hat{\mathbf{x}}+x_8 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_8\right)c\,\hat{\mathbf{z}}$	(8*d*)	O V
$\mathbf{B}_{33}$	=	$x_9\,\mathbf{a}_1+y_9\,\mathbf{a}_2+z_9\,\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}}+y_9 a\,\hat{\mathbf{y}}+z_9 c\,\hat{\mathbf{z}}$	(8*d*)	O VI
$\mathbf{B}_{34}$	=	$-x_9\,\mathbf{a}_1-y_9\,\mathbf{a}_2+z_9\,\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{x}}-y_9 a\,\hat{\mathbf{y}}+z_9 c\,\hat{\mathbf{z}}$	(8*d*)	O VI
$\mathbf{B}_{35}$	=	$-y_9\,\mathbf{a}_1+x_9\,\mathbf{a}_2+z_9\,\mathbf{a}_3$	=	$-y_9 a\,\hat{\mathbf{x}}+x_9 a\,\hat{\mathbf{y}}+z_9 c\,\hat{\mathbf{z}}$	(8*d*)	O VI
$\mathbf{B}_{36}$	=	$y_9\,\mathbf{a}_1-x_9\,\mathbf{a}_2+z_9\,\mathbf{a}_3$	=	$y_9 a\,\hat{\mathbf{x}}-x_9 a\,\hat{\mathbf{y}}+z_9 c\,\hat{\mathbf{z}}$	(8*d*)	O VI
$\mathbf{B}_{37}$	=	$x_9\,\mathbf{a}_1-y_9\,\mathbf{a}_2+\left(\frac{1}{2}+z_9\right)\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}}-y_9 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_9\right)c\,\hat{\mathbf{z}}$	(8*d*)	O VI

$\mathbf{B}_{38}$	=	$-x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{x}} + y_9 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right) c\,\hat{\mathbf{z}}$	(8*d*)	O VI
$\mathbf{B}_{39}$	=	$-y_9\,\mathbf{a}_1 - x_9\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$-y_9 a\,\hat{\mathbf{x}} - x_9 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right) c\,\hat{\mathbf{z}}$	(8*d*)	O VI
$\mathbf{B}_{40}$	=	$y_9\,\mathbf{a}_1 + x_9\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$y_9 a\,\hat{\mathbf{x}} + x_9 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right) c\,\hat{\mathbf{z}}$	(8*d*)	O VI
$\mathbf{B}_{41}$	=	$x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$x_{10} a\,\hat{\mathbf{x}} + y_{10} a\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}}$	(8*d*)	O VII
$\mathbf{B}_{42}$	=	$-x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$-x_{10} a\,\hat{\mathbf{x}} - y_{10} a\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}}$	(8*d*)	O VII
$\mathbf{B}_{43}$	=	$-y_{10}\,\mathbf{a}_1 + x_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$-y_{10} a\,\hat{\mathbf{x}} + x_{10} a\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}}$	(8*d*)	O VII
$\mathbf{B}_{44}$	=	$y_{10}\,\mathbf{a}_1 - x_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$y_{10} a\,\hat{\mathbf{x}} - x_{10} a\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}}$	(8*d*)	O VII
$\mathbf{B}_{45}$	=	$x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{10}\right)\mathbf{a}_3$	=	$x_{10} a\,\hat{\mathbf{x}} - y_{10} a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{10}\right) c\,\hat{\mathbf{z}}$	(8*d*)	O VII
$\mathbf{B}_{46}$	=	$-x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{10}\right)\mathbf{a}_3$	=	$-x_{10} a\,\hat{\mathbf{x}} + y_{10} a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{10}\right) c\,\hat{\mathbf{z}}$	(8*d*)	O VII
$\mathbf{B}_{47}$	=	$-y_{10}\,\mathbf{a}_1 - x_{10}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{10}\right)\mathbf{a}_3$	=	$-y_{10} a\,\hat{\mathbf{x}} - x_{10} a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{10}\right) c\,\hat{\mathbf{z}}$	(8*d*)	O VII
$\mathbf{B}_{48}$	=	$y_{10}\,\mathbf{a}_1 + x_{10}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{10}\right)\mathbf{a}_3$	=	$y_{10} a\,\hat{\mathbf{x}} + x_{10} a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{10}\right) c\,\hat{\mathbf{z}}$	(8*d*)	O VII
$\mathbf{B}_{49}$	=	$x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$x_{11} a\,\hat{\mathbf{x}} + y_{11} a\,\hat{\mathbf{y}} + z_{11} c\,\hat{\mathbf{z}}$	(8*d*)	O VIII
$\mathbf{B}_{50}$	=	$-x_{11}\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$-x_{11} a\,\hat{\mathbf{x}} - y_{11} a\,\hat{\mathbf{y}} + z_{11} c\,\hat{\mathbf{z}}$	(8*d*)	O VIII
$\mathbf{B}_{51}$	=	$-y_{11}\,\mathbf{a}_1 + x_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$-y_{11} a\,\hat{\mathbf{x}} + x_{11} a\,\hat{\mathbf{y}} + z_{11} c\,\hat{\mathbf{z}}$	(8*d*)	O VIII
$\mathbf{B}_{52}$	=	$y_{11}\,\mathbf{a}_1 - x_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$y_{11} a\,\hat{\mathbf{x}} - x_{11} a\,\hat{\mathbf{y}} + z_{11} c\,\hat{\mathbf{z}}$	(8*d*)	O VIII
$\mathbf{B}_{53}$	=	$x_{11}\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{11}\right)\mathbf{a}_3$	=	$x_{11} a\,\hat{\mathbf{x}} - y_{11} a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{11}\right) c\,\hat{\mathbf{z}}$	(8*d*)	O VIII
$\mathbf{B}_{54}$	=	$-x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{11}\right)\mathbf{a}_3$	=	$-x_{11} a\,\hat{\mathbf{x}} + y_{11} a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{11}\right) c\,\hat{\mathbf{z}}$	(8*d*)	O VIII
$\mathbf{B}_{55}$	=	$-y_{11}\,\mathbf{a}_1 - x_{11}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{11}\right)\mathbf{a}_3$	=	$-y_{11} a\,\hat{\mathbf{x}} - x_{11} a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{11}\right) c\,\hat{\mathbf{z}}$	(8*d*)	O VIII
$\mathbf{B}_{56}$	=	$y_{11}\,\mathbf{a}_1 + x_{11}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{11}\right)\mathbf{a}_3$	=	$y_{11} a\,\hat{\mathbf{x}} + x_{11} a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{11}\right) c\,\hat{\mathbf{z}}$	(8*d*)	O VIII
$\mathbf{B}_{57}$	=	$x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3$	=	$x_{12} a\,\hat{\mathbf{x}} + y_{12} a\,\hat{\mathbf{y}} + z_{12} c\,\hat{\mathbf{z}}$	(8*d*)	Se I
$\mathbf{B}_{58}$	=	$-x_{12}\,\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3$	=	$-x_{12} a\,\hat{\mathbf{x}} - y_{12} a\,\hat{\mathbf{y}} + z_{12} c\,\hat{\mathbf{z}}$	(8*d*)	Se I
$\mathbf{B}_{59}$	=	$-y_{12}\,\mathbf{a}_1 + x_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3$	=	$-y_{12} a\,\hat{\mathbf{x}} + x_{12} a\,\hat{\mathbf{y}} + z_{12} c\,\hat{\mathbf{z}}$	(8*d*)	Se I
$\mathbf{B}_{60}$	=	$y_{12}\,\mathbf{a}_1 - x_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3$	=	$y_{12} a\,\hat{\mathbf{x}} - x_{12} a\,\hat{\mathbf{y}} + z_{12} c\,\hat{\mathbf{z}}$	(8*d*)	Se I
$\mathbf{B}_{61}$	=	$x_{12}\,\mathbf{a}_1 - y_{12}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{12}\right)\mathbf{a}_3$	=	$x_{12} a\,\hat{\mathbf{x}} - y_{12} a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{12}\right) c\,\hat{\mathbf{z}}$	(8*d*)	Se I
$\mathbf{B}_{62}$	=	$-x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{12}\right)\mathbf{a}_3$	=	$-x_{12} a\,\hat{\mathbf{x}} + y_{12} a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{12}\right) c\,\hat{\mathbf{z}}$	(8*d*)	Se I
$\mathbf{B}_{63}$	=	$-y_{12}\,\mathbf{a}_1 - x_{12}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{12}\right)\mathbf{a}_3$	=	$-y_{12} a\,\hat{\mathbf{x}} - x_{12} a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{12}\right) c\,\hat{\mathbf{z}}$	(8*d*)	Se I
$\mathbf{B}_{64}$	=	$y_{12}\,\mathbf{a}_1 + x_{12}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{12}\right)\mathbf{a}_3$	=	$y_{12} a\,\hat{\mathbf{x}} + x_{12} a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{12}\right) c\,\hat{\mathbf{z}}$	(8*d*)	Se I
$\mathbf{B}_{65}$	=	$x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3$	=	$x_{13} a\,\hat{\mathbf{x}} + y_{13} a\,\hat{\mathbf{y}} + z_{13} c\,\hat{\mathbf{z}}$	(8*d*)	Se II
$\mathbf{B}_{66}$	=	$-x_{13}\,\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3$	=	$-x_{13} a\,\hat{\mathbf{x}} - y_{13} a\,\hat{\mathbf{y}} + z_{13} c\,\hat{\mathbf{z}}$	(8*d*)	Se II
$\mathbf{B}_{67}$	=	$-y_{13}\,\mathbf{a}_1 + x_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3$	=	$-y_{13} a\,\hat{\mathbf{x}} + x_{13} a\,\hat{\mathbf{y}} + z_{13} c\,\hat{\mathbf{z}}$	(8*d*)	Se II
$\mathbf{B}_{68}$	=	$y_{13}\,\mathbf{a}_1 - x_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3$	=	$y_{13} a\,\hat{\mathbf{x}} - x_{13} a\,\hat{\mathbf{y}} + z_{13} c\,\hat{\mathbf{z}}$	(8*d*)	Se II
$\mathbf{B}_{69}$	=	$x_{13}\,\mathbf{a}_1 - y_{13}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{13}\right)\mathbf{a}_3$	=	$x_{13} a\,\hat{\mathbf{x}} - y_{13} a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{13}\right) c\,\hat{\mathbf{z}}$	(8*d*)	Se II
$\mathbf{B}_{70}$	=	$-x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{13}\right)\mathbf{a}_3$	=	$-x_{13} a\,\hat{\mathbf{x}} + y_{13} a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{13}\right) c\,\hat{\mathbf{z}}$	(8*d*)	Se II
$\mathbf{B}_{71}$	=	$-y_{13}\,\mathbf{a}_1 - x_{13}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{13}\right)\mathbf{a}_3$	=	$-y_{13} a\,\hat{\mathbf{x}} - x_{13} a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{13}\right) c\,\hat{\mathbf{z}}$	(8*d*)	Se II
$\mathbf{B}_{72}$	=	$y_{13}\,\mathbf{a}_1 + x_{13}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{13}\right)\mathbf{a}_3$	=	$y_{13} a\,\hat{\mathbf{x}} + x_{13} a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{13}\right) c\,\hat{\mathbf{z}}$	(8*d*)	Se II

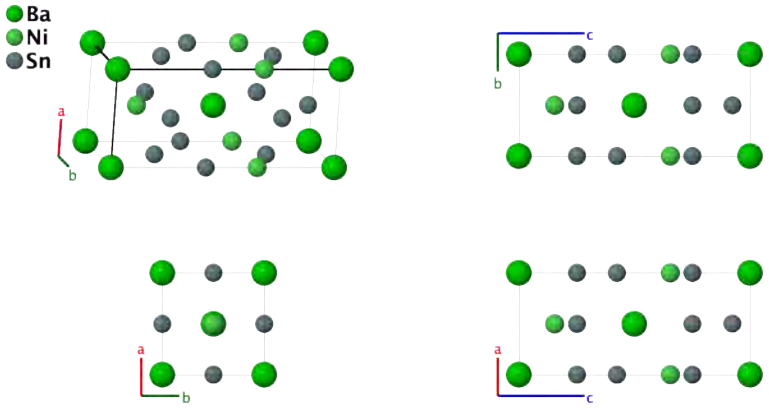
References:

- G. Meunier, M. Bertaud, and J. Galy, *Cristallochimie du sélénium(+IV). I. VSe_2O_6, une structure à trois chaînes parallèles $(VO_5)_n^{6n-}$ indépendantes pontées par des groupements $(Se_2O)^{6+}$*, Acta Crystallogr. Sect. B Struct. Sci. **30**, 2834–2839 (1974), doi:10.1107/S0567740874008260.

Geometry files:
- CIF: pp. 1679
- POSCAR: pp. 1679

$BaNiSn_3$ Structure: ABC3_tI10_107_a_a_ab

http://aflow.org/prototype-encyclopedia/ABC3_tI10_107_a_a_ab



Prototype	:	$BaNiSn_3$
AFLOW prototype label	:	ABC3_tI10_107_a_a_ab
***Strukturbericht* designation**	:	None
Pearson symbol	:	tI10
Space group number	:	107
Space group symbol	:	*I*4*mm*
AFLOW prototype command	:	`aflow --proto=ABC3_tI10_107_a_a_ab --params=`$a, c/a, z_1, z_2, z_3, z_4$

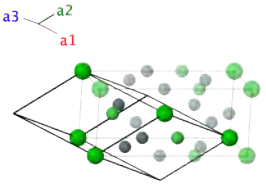
Other compounds with this structure

- $BaPdSn_3$, $CeCuAl_3$, $EuPdGe_3$, $LaOsSb_3$, $SrIrAl_3$, $SrNiSn_3$, $SrPdAl_3$, and $SrPtAl_3$

- This is a ternary form of the $D1_3$ ($BaAl_4$) structure. The atomic positions in both conventional cells are approximately the same, but the distribution of the atoms on those sites and the resulting relaxation leads to a different structure.
- Although (Dörrscheidt, 1978) give the structural information for $BaPtSn_3$ before that of $BaNiSn_3$, (Shatruk, 2019) and others list $BaNiSn_3$ as the prototype for this structure.
- Space group *I*4*mm* #107 does not specify the origin of the z-axis. (Dörrscheidt, 1978) places the barium atom at the origin.

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} - \frac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$\mathbf{B}_1$	=	$z_1\,\mathbf{a}_1 + z_1\,\mathbf{a}_2$	=	$z_1 c\,\hat{\mathbf{z}}$	($2a$)	Ba
$\mathbf{B}_2$	=	$z_2\,\mathbf{a}_1 + z_2\,\mathbf{a}_2$	=	$z_2 c\,\hat{\mathbf{z}}$	($2a$)	Ni
$\mathbf{B}_3$	=	$z_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_2$	=	$z_3 c\,\hat{\mathbf{z}}$	($2a$)	Sn I
$\mathbf{B}_4$	=	$\left(\frac{1}{2} + z_4\right)\mathbf{a}_1 + z_4\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	($4b$)	Sn II
$\mathbf{B}_5$	=	$z_4\,\mathbf{a}_1 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + z_4 c\,\hat{\mathbf{z}}$	($4b$)	Sn II

References:

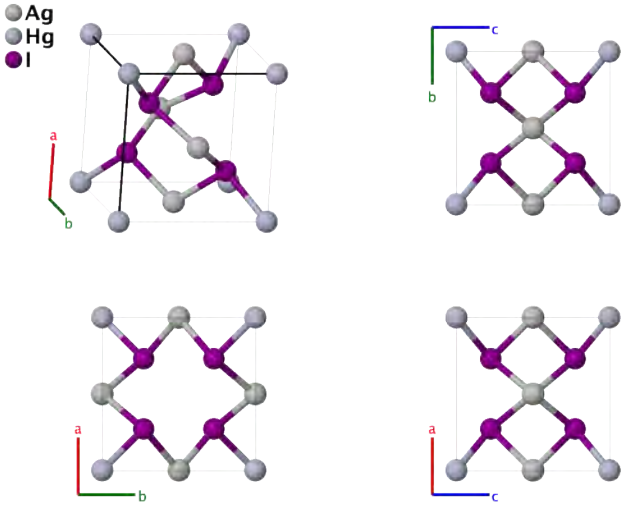
- W. Dörrscheidt and H. Schäfer, *Die Struktur des $BaPtSn_3$, $BaNiSn_3$ und $SrNiSn_3$ und ihre Verwandtschaft zum $ThCr_2Si_2$-Strukturtyp*, J. Less-Common Met. **58**, 209–216 (1978), doi:10.1016/0022-5088(78)90202-3.

Geometry files:

- CIF: pp. 1680
- POSCAR: pp. 1680

$E3_1$ (β-Ag_2HgI_4) (*obsolete*) Structure: A2BC4_tP7_111_f_a_n

http://aflow.org/prototype-encyclopedia/A2BC4_tP7_111_f_a_n

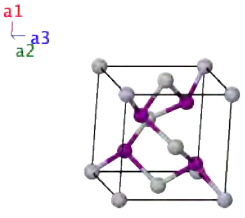


Prototype	:	Ag_2HgI_4
AFLOW prototype label	:	A2BC4_tP7_111_f_a_n
***Strukturbericht* designation**	:	$E3_1$
Pearson symbol	:	tP7
Space group number	:	111
Space group symbol	:	$P\bar{4}2m$
AFLOW prototype command	:	`aflow --proto=A2BC4_tP7_111_f_a_n --params=`$a, c/a, x_3, z_3$

- (Ketelaar, 1931) determined this pseudo-cubic structure for β-Ag_2HgI_4, and (Hermann, 1937) assigned it the *Strukturbericht* symbol $E3_1$. Later, (Browall, 1974) showed that β-Ag_2HgI_4 takes the Al_2CdS_4 structure, which some authorities give as *Strukturbericht* $E3$.

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$ =	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($1a$)	Hg

$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(2f)$	Ag
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(2f)$	Ag
$\mathbf{B}_4$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	$(4n)$	I
$\mathbf{B}_5$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	$(4n)$	I
$\mathbf{B}_6$	=	$x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	$(4n)$	I
$\mathbf{B}_7$	=	$-x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	$(4n)$	I

References:

- J. A. A. Ketelaar, *Strukturbestimmung der komplexen Quecksilberverbindungen Ag_2HgJ_4 und Cu_2HgJ_4*, Zeitschrift für Kristallographie - Crystalline Materials **80**, 190–203 (1931), doi:10.1524/zkri.1931.80.1.190.
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Found in:

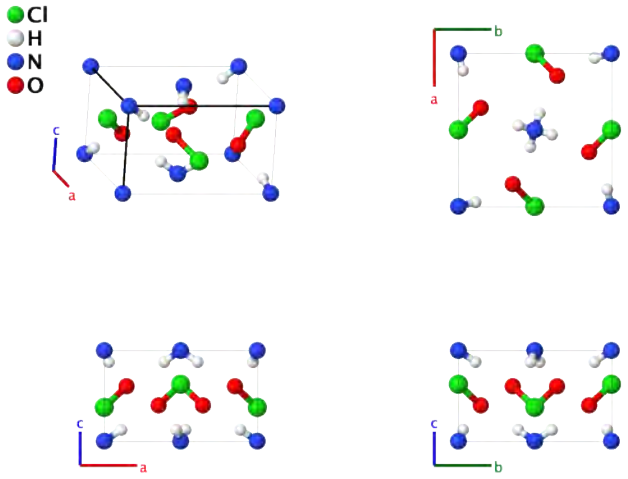
- K. W. Browall, J. S. Kasper, and H. Wiedemeier, *Single-crystal studies of β-Ag_2HgI_4*, J. Solid State Chem. **10**, 20–28 (1974), doi:10.1016/0022-4596(74)90004-8.

Geometry files:

- CIF: pp. 1680
- POSCAR: pp. 1680

Ammonium Chlorite (NH_4ClO_2) Structure: AB4CD2_tP16_113_c_f_a_e

http://aflow.org/prototype-encyclopedia/AB4CD2_tP16_113_c_f_a_e

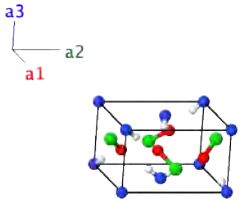


Prototype	:	ClH_4NO_2
AFLOW prototype label	:	AB4CD2_tP16_113_c_f_a_e
***Strukturbericht* designation**	:	None
Pearson symbol	:	tP16
Space group number	:	113
Space group symbol	:	$P\bar{4}2_1m$
AFLOW prototype command	:	`aflow --proto=AB4CD2_tP16_113_c_f_a_e --params=`$a, c/a, z_2, x_3, z_3, x_4, y_4, z_4$

- This is the updated version of the $F5_4$ structure, including the positions of the hydrogen atoms.
- The data for this structure was taken at 150 K.
- Not to be confused with ammonium chlorate, NH_4ClO_3, which takes on the KNO_3 structure, or NH_4ClO_4, which has the $H0_5$ structure.

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	N
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$	$(2a)$	N
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + z_2c\,\hat{\mathbf{z}}$	$(2c)$	Cl
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 - z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - z_2c\,\hat{\mathbf{z}}$	$(2c)$	Cl
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	$(4e)$	O
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	$(4e)$	O
$\mathbf{B}_7$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 - x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	$(4e)$	O
$\mathbf{B}_8$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	$(4e)$	O
$\mathbf{B}_9$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + y_4a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	$(8f)$	H
$\mathbf{B}_{10}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} - y_4a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	$(8f)$	H
$\mathbf{B}_{11}$	=	$y_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$y_4a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	$(8f)$	H
$\mathbf{B}_{12}$	=	$-y_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-y_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	$(8f)$	H
$\mathbf{B}_{13}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	$(8f)$	H
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	$(8f)$	H
$\mathbf{B}_{15}$	=	$\left(\frac{1}{2} - y_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	$(8f)$	H
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + y_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	$(8f)$	H

References:

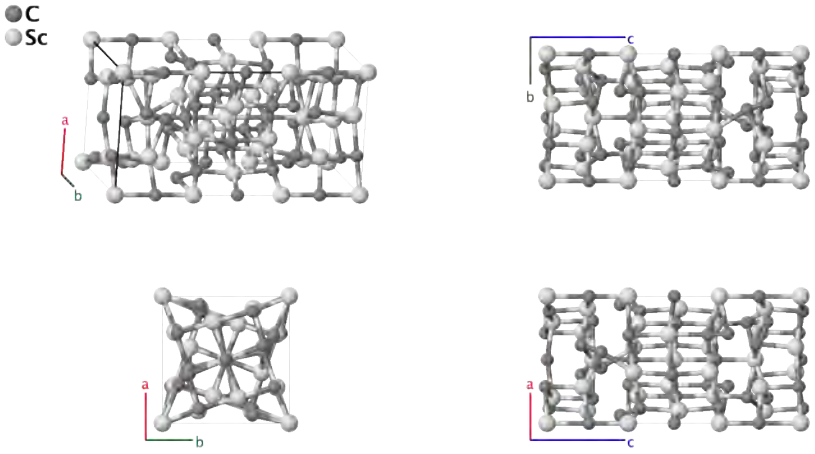
- A. I. Smolentsev and D. Y. Naumov, *Ammonium chlorite, NH_4ClO_2, at 150 K*, Acta Crystallogr. E **61**, i38–i40 (2005), doi:10.1107/S1600536805005088.

Geometry files:

- CIF: pp. 1680
- POSCAR: pp. 1681

$C_{19}Sc_{15}$ Structure: A19B15_tP68_114_bc4e_ac3e

http://aflow.org/prototype-encyclopedia/A19B15_tP68_114_bc4e_ac3e



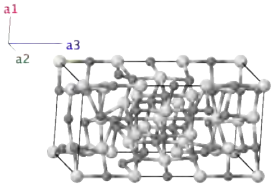
Prototype	:	$C_{19}Sc_{15}$
AFLOW prototype label	:	A19B15_tP68_114_bc4e_ac3e
***Strukturbericht* designation**	:	None
Pearson symbol	:	tP68
Space group number	:	114
Space group symbol	:	$P\bar{4}2_1c$
AFLOW prototype command	:	`aflow --proto=A19B15_tP68_114_bc4e_ac3e --params=`$a, c/a, z_3, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}$

Other compounds with this structure

- $C_{19}Er_{15}$, $C_{19}Lu_{15}$, $C_{19}Tm_{15}$, $C_{19}Y_{15}$, and $C_{19}Yb_{15}$

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(2*a*)	Sc I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(2*a*)	Sc I
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	(2*b*)	C I
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$	(2*b*)	C I
$\mathbf{B}_5$	=	$z_3\,\mathbf{a}_3$	=	$z_3 c\,\hat{\mathbf{z}}$	(4*c*)	C II

$\mathbf{B}_6$	=	$-z_3\,\mathbf{a}_3$	=	$-z_3 c\,\hat{\mathbf{z}}$	($4c$)	C II
$\mathbf{B}_7$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	($4c$)	C II
$\mathbf{B}_8$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	($4c$)	C II
$\mathbf{B}_9$	=	$z_4\,\mathbf{a}_3$	=	$z_4 c\,\hat{\mathbf{z}}$	($4c$)	Sc II
$\mathbf{B}_{10}$	=	$-z_4\,\mathbf{a}_3$	=	$-z_4 c\,\hat{\mathbf{z}}$	($4c$)	Sc II
$\mathbf{B}_{11}$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	($4c$)	Sc II
$\mathbf{B}_{12}$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	($4c$)	Sc II
$\mathbf{B}_{13}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	($8e$)	C III
$\mathbf{B}_{14}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	($8e$)	C III
$\mathbf{B}_{15}$	=	$y_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$y_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	($8e$)	C III
$\mathbf{B}_{16}$	=	$-y_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-y_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	($8e$)	C III
$\mathbf{B}_{17}$	=	$\left(\frac{1}{2} - x_5\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	($8e$)	C III
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} + x_5\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	($8e$)	C III
$\mathbf{B}_{19}$	=	$\left(\frac{1}{2} - y_5\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - x_5\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	($8e$)	C III
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} + y_5\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + x_5\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	($8e$)	C III
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + y_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	($8e$)	C IV
$\mathbf{B}_{22}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - y_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	($8e$)	C IV
$\mathbf{B}_{23}$	=	$y_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$y_6 a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	($8e$)	C IV
$\mathbf{B}_{24}$	=	$-y_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$-y_6 a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	($8e$)	C IV
$\mathbf{B}_{25}$	=	$\left(\frac{1}{2} - x_6\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\,\hat{\mathbf{z}}$	($8e$)	C IV
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} + x_6\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\,\hat{\mathbf{z}}$	($8e$)	C IV
$\mathbf{B}_{27}$	=	$\left(\frac{1}{2} - y_6\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - x_6\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\,\hat{\mathbf{z}}$	($8e$)	C IV
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} + y_6\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + x_6\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\,\hat{\mathbf{z}}$	($8e$)	C IV
$\mathbf{B}_{29}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + y_7 a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	($8e$)	C V
$\mathbf{B}_{30}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} - y_7 a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	($8e$)	C V
$\mathbf{B}_{31}$	=	$y_7\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$y_7 a\,\hat{\mathbf{x}} - x_7 a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	($8e$)	C V
$\mathbf{B}_{32}$	=	$-y_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$-y_7 a\,\hat{\mathbf{x}} + x_7 a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	($8e$)	C V
$\mathbf{B}_{33}$	=	$\left(\frac{1}{2} - x_7\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right)c\,\hat{\mathbf{z}}$	($8e$)	C V

$$\mathbf{B}_{34} = \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_7\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_7\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{C V}$$

$$\mathbf{B}_{35} = \left(\tfrac{1}{2}-y_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}-x_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-y_7\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_7\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{C V}$$

$$\mathbf{B}_{36} = \left(\tfrac{1}{2}+y_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+y_7\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_7\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{C V}$$

$$\mathbf{B}_{37} = x_8\,\mathbf{a}_1+y_8\,\mathbf{a}_2+z_8\,\mathbf{a}_3 = x_8a\,\hat{\mathbf{x}}+y_8a\,\hat{\mathbf{y}}+z_8c\,\hat{\mathbf{z}} \quad (8e) \quad \text{C VI}$$

$$\mathbf{B}_{38} = -x_8\,\mathbf{a}_1-y_8\,\mathbf{a}_2+z_8\,\mathbf{a}_3 = -x_8a\,\hat{\mathbf{x}}-y_8a\,\hat{\mathbf{y}}+z_8c\,\hat{\mathbf{z}} \quad (8e) \quad \text{C VI}$$

$$\mathbf{B}_{39} = y_8\,\mathbf{a}_1-x_8\,\mathbf{a}_2-z_8\,\mathbf{a}_3 = y_8a\,\hat{\mathbf{x}}-x_8a\,\hat{\mathbf{y}}-z_8c\,\hat{\mathbf{z}} \quad (8e) \quad \text{C VI}$$

$$\mathbf{B}_{40} = -y_8\,\mathbf{a}_1+x_8\,\mathbf{a}_2-z_8\,\mathbf{a}_3 = -y_8a\,\hat{\mathbf{x}}+x_8a\,\hat{\mathbf{y}}-z_8c\,\hat{\mathbf{z}} \quad (8e) \quad \text{C VI}$$

$$\mathbf{B}_{41} = \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_8\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_8\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{C VI}$$

$$\mathbf{B}_{42} = \left(\tfrac{1}{2}+x_8\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_8\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_8\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{C VI}$$

$$\mathbf{B}_{43} = \left(\tfrac{1}{2}-y_8\right)\mathbf{a}_1+\left(\tfrac{1}{2}-x_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-y_8\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_8\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{C VI}$$

$$\mathbf{B}_{44} = \left(\tfrac{1}{2}+y_8\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+y_8\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_8\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{C VI}$$

$$\mathbf{B}_{45} = x_9\,\mathbf{a}_1+y_9\,\mathbf{a}_2+z_9\,\mathbf{a}_3 = x_9a\,\hat{\mathbf{x}}+y_9a\,\hat{\mathbf{y}}+z_9c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Sc III}$$

$$\mathbf{B}_{46} = -x_9\,\mathbf{a}_1-y_9\,\mathbf{a}_2+z_9\,\mathbf{a}_3 = -x_9a\,\hat{\mathbf{x}}-y_9a\,\hat{\mathbf{y}}+z_9c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Sc III}$$

$$\mathbf{B}_{47} = y_9\,\mathbf{a}_1-x_9\,\mathbf{a}_2-z_9\,\mathbf{a}_3 = y_9a\,\hat{\mathbf{x}}-x_9a\,\hat{\mathbf{y}}-z_9c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Sc III}$$

$$\mathbf{B}_{48} = -y_9\,\mathbf{a}_1+x_9\,\mathbf{a}_2-z_9\,\mathbf{a}_3 = -y_9a\,\hat{\mathbf{x}}+x_9a\,\hat{\mathbf{y}}-z_9c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Sc III}$$

$$\mathbf{B}_{49} = \left(\tfrac{1}{2}-x_9\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_9\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_9\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_9\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_9\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Sc III}$$

$$\mathbf{B}_{50} = \left(\tfrac{1}{2}+x_9\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_9\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_9\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_9\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_9\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Sc III}$$

$$\mathbf{B}_{51} = \left(\tfrac{1}{2}-y_9\right)\mathbf{a}_1+\left(\tfrac{1}{2}-x_9\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-y_9\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_9\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_9\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Sc III}$$

$$\mathbf{B}_{52} = \left(\tfrac{1}{2}+y_9\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_9\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+y_9\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_9\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_9\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Sc III}$$

$$\mathbf{B}_{53} = x_{10}\,\mathbf{a}_1+y_{10}\,\mathbf{a}_2+z_{10}\,\mathbf{a}_3 = x_{10}a\,\hat{\mathbf{x}}+y_{10}a\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Sc IV}$$

$$\mathbf{B}_{54} = -x_{10}\,\mathbf{a}_1-y_{10}\,\mathbf{a}_2+z_{10}\,\mathbf{a}_3 = -x_{10}a\,\hat{\mathbf{x}}-y_{10}a\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Sc IV}$$

$$\mathbf{B}_{55} = y_{10}\,\mathbf{a}_1-x_{10}\,\mathbf{a}_2-z_{10}\,\mathbf{a}_3 = y_{10}a\,\hat{\mathbf{x}}-x_{10}a\,\hat{\mathbf{y}}-z_{10}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Sc IV}$$

$$\mathbf{B}_{56} = -y_{10}\,\mathbf{a}_1+x_{10}\,\mathbf{a}_2-z_{10}\,\mathbf{a}_3 = -y_{10}a\,\hat{\mathbf{x}}+x_{10}a\,\hat{\mathbf{y}}-z_{10}c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Sc IV}$$

$$\mathbf{B}_{57} = \left(\tfrac{1}{2}-x_{10}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{10}\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{10}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{10}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{10}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Sc IV}$$

$$\mathbf{B}_{58} = \left(\tfrac{1}{2}+x_{10}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{10}\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{10}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{10}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{10}\right)c\,\hat{\mathbf{z}} \quad (8e) \quad \text{Sc IV}$$

$\mathbf{B}_{59}$	=	$\left(\frac{1}{2}-y_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}-x_{10}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{10}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{10}\right)c\,\hat{\mathbf{z}}$	(8*e*)	Sc IV
$\mathbf{B}_{60}$	=	$\left(\frac{1}{2}+y_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}+x_{10}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{10}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{10}\right)c\,\hat{\mathbf{z}}$	(8*e*)	Sc IV
$\mathbf{B}_{61}$	=	$x_{11}\,\mathbf{a}_1+y_{11}\,\mathbf{a}_2+z_{11}\,\mathbf{a}_3$	=	$x_{11}a\,\hat{\mathbf{x}}+y_{11}a\,\hat{\mathbf{y}}+z_{11}c\,\hat{\mathbf{z}}$	(8*e*)	Sc V
$\mathbf{B}_{62}$	=	$-x_{11}\,\mathbf{a}_1-y_{11}\,\mathbf{a}_2+z_{11}\,\mathbf{a}_3$	=	$-x_{11}a\,\hat{\mathbf{x}}-y_{11}a\,\hat{\mathbf{y}}+z_{11}c\,\hat{\mathbf{z}}$	(8*e*)	Sc V
$\mathbf{B}_{63}$	=	$y_{11}\,\mathbf{a}_1-x_{11}\,\mathbf{a}_2-z_{11}\,\mathbf{a}_3$	=	$y_{11}a\,\hat{\mathbf{x}}-x_{11}a\,\hat{\mathbf{y}}-z_{11}c\,\hat{\mathbf{z}}$	(8*e*)	Sc V
$\mathbf{B}_{64}$	=	$-y_{11}\,\mathbf{a}_1+x_{11}\,\mathbf{a}_2-z_{11}\,\mathbf{a}_3$	=	$-y_{11}a\,\hat{\mathbf{x}}+x_{11}a\,\hat{\mathbf{y}}-z_{11}c\,\hat{\mathbf{z}}$	(8*e*)	Sc V
$\mathbf{B}_{65}$	=	$\left(\frac{1}{2}-x_{11}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{11}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{11}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{11}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{11}\right)c\,\hat{\mathbf{z}}$	(8*e*)	Sc V
$\mathbf{B}_{66}$	=	$\left(\frac{1}{2}+x_{11}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{11}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{11}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{11}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{11}\right)c\,\hat{\mathbf{z}}$	(8*e*)	Sc V
$\mathbf{B}_{67}$	=	$\left(\frac{1}{2}-y_{11}\right)\mathbf{a}_1+\left(\frac{1}{2}-x_{11}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{11}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_{11}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{11}\right)c\,\hat{\mathbf{z}}$	(8*e*)	Sc V
$\mathbf{B}_{68}$	=	$\left(\frac{1}{2}+y_{11}\right)\mathbf{a}_1+\left(\frac{1}{2}+x_{11}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{11}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_{11}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{11}\right)c\,\hat{\mathbf{z}}$	(8*e*)	Sc V

References:

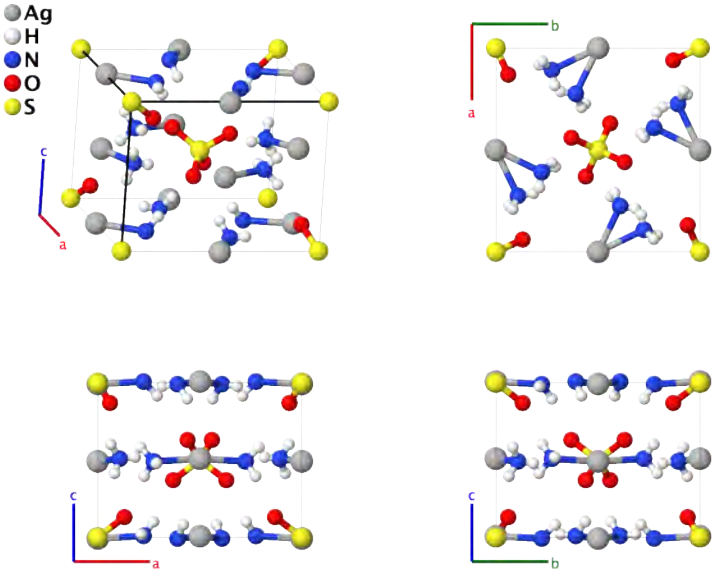
- H. Jedlicka, H. Nowotny, and F. Benesovsky, *Zum System Scandium-Kohlenstoff, 2. Mitt.: Kristallstruktur des C-reichen Carbids*, Monatshefte für Chemie - Chemical Monthly **102**, 389–403 (1971), doi:10.1007/BF00909332.

Geometry files:

- CIF: pp. 1681
- POSCAR: pp. 1681

$Ag_2SO_4{\cdot}4NH_3$ ($H4_{17}$) Structure: A2B12C4D4E_tP46_114_d_3e_e_e_a

http://aflow.org/prototype-encyclopedia/A2B12C4D4E_tP46_114_d_3e_e_e_a

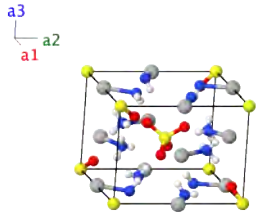


Prototype	:	$Ag_2H_{12}N_4O_4S$
AFLOW prototype label	:	A2B12C4D4E_tP46_114_d_3e_e_e_a
***Strukturbericht* designation**	:	$H4_{17}$
Pearson symbol	:	tP46
Space group number	:	114
Space group symbol	:	$P\bar{4}2_1c$
AFLOW prototype command	:	`aflow --proto=A2B12C4D4E_tP46_114_d_3e_e_e_a --params=`$a, c/a, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7$

- We use the data from (Zachwieja, 1992), which refines the data of the original $H4_{17}$ structure of (Corey, 1934) and locates the hydrogen atoms.

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(2*a*)	S
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(2*a*)	S
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(4*d*)	Ag
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 - z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - z_2 c\,\hat{\mathbf{z}}$	(4*d*)	Ag
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_1 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	(4*d*)	Ag
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	(4*d*)	Ag
$\mathbf{B}_7$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(8*e*)	H I
$\mathbf{B}_8$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(8*e*)	H I
$\mathbf{B}_9$	=	$y_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$y_3 a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(8*e*)	H I
$\mathbf{B}_{10}$	=	$-y_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-y_3 a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(8*e*)	H I
$\mathbf{B}_{11}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(8*e*)	H I
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(8*e*)	H I
$\mathbf{B}_{13}$	=	$\left(\frac{1}{2} - y_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(8*e*)	H I
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} + y_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(8*e*)	H I
$\mathbf{B}_{15}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*e*)	H II
$\mathbf{B}_{16}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(8*e*)	H II
$\mathbf{B}_{17}$	=	$y_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$y_4 a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*e*)	H II
$\mathbf{B}_{18}$	=	$-y_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-y_4 a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(8*e*)	H II
$\mathbf{B}_{19}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(8*e*)	H II
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(8*e*)	H II
$\mathbf{B}_{21}$	=	$\left(\frac{1}{2} - y_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(8*e*)	H II
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} + y_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(8*e*)	H II
$\mathbf{B}_{23}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(8*e*)	H III
$\mathbf{B}_{24}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(8*e*)	H III
$\mathbf{B}_{25}$	=	$y_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$y_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(8*e*)	H III
$\mathbf{B}_{26}$	=	$-y_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-y_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(8*e*)	H III
$\mathbf{B}_{27}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	(8*e*)	H III
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	(8*e*)	H III

$\mathbf{B}_{29}$	=	$\left(\frac{1}{2}-y_5\right)\mathbf{a}_1+\left(\frac{1}{2}-x_5\right)\mathbf{a}_2+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	(8*e*)	H III
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2}+y_5\right)\mathbf{a}_1+\left(\frac{1}{2}+x_5\right)\mathbf{a}_2+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	(8*e*)	H III
$\mathbf{B}_{31}$	=	$x_6\,\mathbf{a}_1+y_6\,\mathbf{a}_2+z_6\,\mathbf{a}_3$	=	$x_6a\,\hat{\mathbf{x}}+y_6a\,\hat{\mathbf{y}}+z_6c\,\hat{\mathbf{z}}$	(8*e*)	N
$\mathbf{B}_{32}$	=	$-x_6\,\mathbf{a}_1-y_6\,\mathbf{a}_2+z_6\,\mathbf{a}_3$	=	$-x_6a\,\hat{\mathbf{x}}-y_6a\,\hat{\mathbf{y}}+z_6c\,\hat{\mathbf{z}}$	(8*e*)	N
$\mathbf{B}_{33}$	=	$y_6\,\mathbf{a}_1-x_6\,\mathbf{a}_2-z_6\,\mathbf{a}_3$	=	$y_6a\,\hat{\mathbf{x}}-x_6a\,\hat{\mathbf{y}}-z_6c\,\hat{\mathbf{z}}$	(8*e*)	N
$\mathbf{B}_{34}$	=	$-y_6\,\mathbf{a}_1+x_6\,\mathbf{a}_2-z_6\,\mathbf{a}_3$	=	$-y_6a\,\hat{\mathbf{x}}+x_6a\,\hat{\mathbf{y}}-z_6c\,\hat{\mathbf{z}}$	(8*e*)	N
$\mathbf{B}_{35}$	=	$\left(\frac{1}{2}-x_6\right)\mathbf{a}_1+\left(\frac{1}{2}+y_6\right)\mathbf{a}_2+\left(\frac{1}{2}-z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_6\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_6\right)c\,\hat{\mathbf{z}}$	(8*e*)	N
$\mathbf{B}_{36}$	=	$\left(\frac{1}{2}+x_6\right)\mathbf{a}_1+\left(\frac{1}{2}-y_6\right)\mathbf{a}_2+\left(\frac{1}{2}-z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_6\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_6\right)c\,\hat{\mathbf{z}}$	(8*e*)	N
$\mathbf{B}_{37}$	=	$\left(\frac{1}{2}-y_6\right)\mathbf{a}_1+\left(\frac{1}{2}-x_6\right)\mathbf{a}_2+\left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	(8*e*)	N
$\mathbf{B}_{38}$	=	$\left(\frac{1}{2}+y_6\right)\mathbf{a}_1+\left(\frac{1}{2}+x_6\right)\mathbf{a}_2+\left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	(8*e*)	N
$\mathbf{B}_{39}$	=	$x_7\,\mathbf{a}_1+y_7\,\mathbf{a}_2+z_7\,\mathbf{a}_3$	=	$x_7a\,\hat{\mathbf{x}}+y_7a\,\hat{\mathbf{y}}+z_7c\,\hat{\mathbf{z}}$	(8*e*)	O
$\mathbf{B}_{40}$	=	$-x_7\,\mathbf{a}_1-y_7\,\mathbf{a}_2+z_7\,\mathbf{a}_3$	=	$-x_7a\,\hat{\mathbf{x}}-y_7a\,\hat{\mathbf{y}}+z_7c\,\hat{\mathbf{z}}$	(8*e*)	O
$\mathbf{B}_{41}$	=	$y_7\,\mathbf{a}_1-x_7\,\mathbf{a}_2-z_7\,\mathbf{a}_3$	=	$y_7a\,\hat{\mathbf{x}}-x_7a\,\hat{\mathbf{y}}-z_7c\,\hat{\mathbf{z}}$	(8*e*)	O
$\mathbf{B}_{42}$	=	$-y_7\,\mathbf{a}_1+x_7\,\mathbf{a}_2-z_7\,\mathbf{a}_3$	=	$-y_7a\,\hat{\mathbf{x}}+x_7a\,\hat{\mathbf{y}}-z_7c\,\hat{\mathbf{z}}$	(8*e*)	O
$\mathbf{B}_{43}$	=	$\left(\frac{1}{2}-x_7\right)\mathbf{a}_1+\left(\frac{1}{2}+y_7\right)\mathbf{a}_2+\left(\frac{1}{2}-z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_7\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_7\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_7\right)c\,\hat{\mathbf{z}}$	(8*e*)	O
$\mathbf{B}_{44}$	=	$\left(\frac{1}{2}+x_7\right)\mathbf{a}_1+\left(\frac{1}{2}-y_7\right)\mathbf{a}_2+\left(\frac{1}{2}-z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_7\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_7\right)c\,\hat{\mathbf{z}}$	(8*e*)	O
$\mathbf{B}_{45}$	=	$\left(\frac{1}{2}-y_7\right)\mathbf{a}_1+\left(\frac{1}{2}-x_7\right)\mathbf{a}_2+\left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_7\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_7\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	(8*e*)	O
$\mathbf{B}_{46}$	=	$\left(\frac{1}{2}+y_7\right)\mathbf{a}_1+\left(\frac{1}{2}+x_7\right)\mathbf{a}_2+\left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_7\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	(8*e*)	O

References:

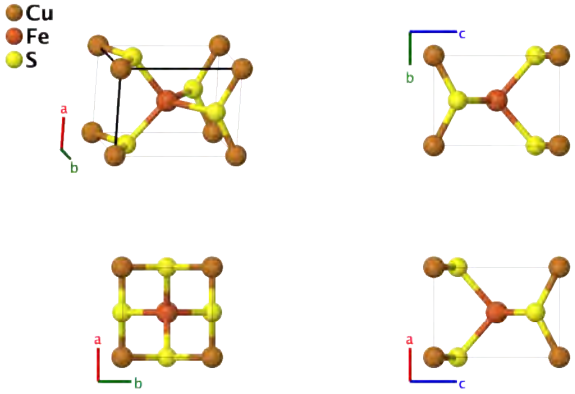
- U. Zachwieja and H. Jacobs, *Redetermination of the crystal structure of diammine silver(I)-sulfate, $[Ag(NH_3)_2]_2SO_4$*, Zeitschrift für Kristallographie - Crystalline Materials **201**, 207–212 (1992), doi:10.1524/zkri.1992.201.14.207.
- R. B. Corey and R. W. G. Wyckoff, *The Crystal Structure of Silver Sulfate Tetrammoniate*, Zeitschrift für Kristallographie - Crystalline Materials **87**, 264–274 (1934), doi:10.1524/zkri.1934.87.1.264.

Geometry files:

- CIF: pp. 1682
- POSCAR: pp. 1682

$F6_1$ (Chalcopyrite, $CuFeS_2$) (*obsolete*) Structure: ABC2_tP4_115_a_c_g

http://aflow.org/prototype-encyclopedia/ABC2_tP4_115_a_c_g

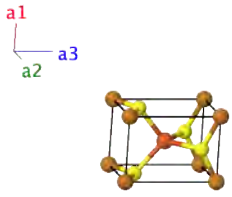


Prototype	:	$CuFeS_2$
AFLOW prototype label	:	ABC2_tP4_115_a_c_g
***Strukturbericht* designation**	:	$F6_1$
Pearson symbol	:	tP4
Space group number	:	115
Space group symbol	:	$P\bar{4}m2$
AFLOW prototype command	:	`aflow --proto=ABC2_tP4_115_a_c_g --params=`$a, c/a, z_3$

- This structure was presented as the chalcopyrite structure and given the *Strukturbericht* designation $F6_1$ in (Ewald, 1931). It was subsequently replaced with the $E1_1$ (ABC2_tI16_122_a_b_d) structure, which is now the accepted structure for chalcopyrite and similar compounds. We include the $F6_1$ structure as part of this historical record.

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(1a)$	Cu
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(1c)$	Fe
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(2g)$	S
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 - z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - z_3 c\,\hat{\mathbf{z}}$	$(2g)$	S

References:

- R. Groß and N. Groß, *Die Atomanordnung des Kupferkieses und die Struktur der Beruehrungsflaechen gesetzmaessig verwachsener Kristalle*, N. Jb. Miner. Mh., Abt. A **48**, 113–135 (1923).

Found in:

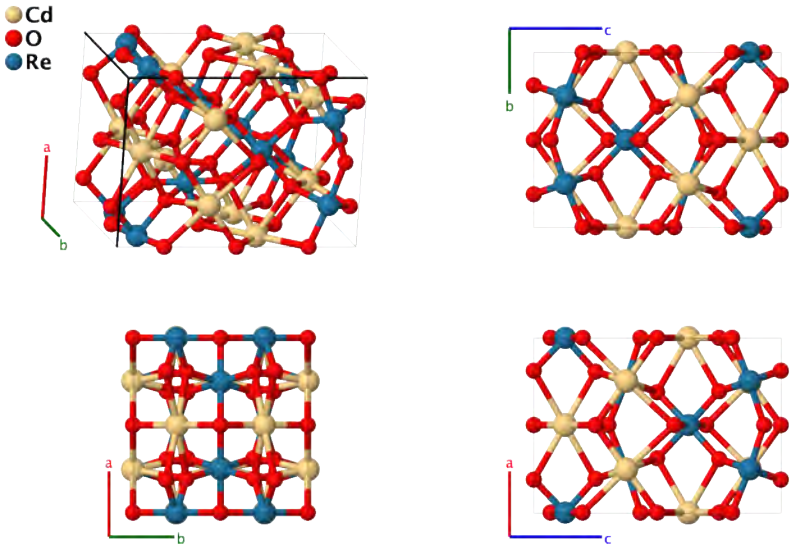
- P. P. Ewald and C. Hermann, eds., *Strukturbericht 1913-1928* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1931).

Geometry files:

- CIF: pp. 1682
- POSCAR: pp. 1683

Phase II $Cd_2Re_2O_7$ Structure: A2B7C2_tI44_119_i_bdefgh_i

http://aflow.org/prototype-encyclopedia/A2B7C2_tI44_119_i_bdefgh_i

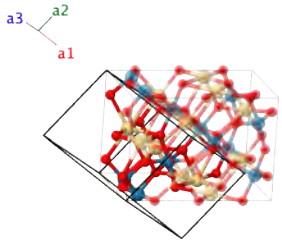


Prototype	:	$Cd_2O_7Re_2$
AFLOW prototype label	:	A2B7C2_tI44_119_i_bdefgh_i
***Strukturbericht* designation**	:	None
Pearson symbol	:	tI44
Space group number	:	119
Space group symbol	:	$I\bar{4}m2$
AFLOW prototype command	:	`aflow --proto=A2B7C2_tI44_119_i_bdefgh_i --params=`$a, c/a, z_3, z_4, x_5, x_6, x_7, z_7, x_8, z_8$

- $Cd_2Re_2O_7$ exhibits a number of phases. We will use the notation of (Kapcia, 2019) to describe them:
  - Phase I: above 200 K, the system takes on the cubic pyrochlore ($E8_1$) structure.
  - Phase II: in the range 120-200 K the system is in the tetragonal $I\bar{4}m2$ #119 structure (this structure).
  - Phase III: in the range 80-120 K the system is in the tetragonal $I4_122$ #98 structure.
  - Phase IV: (Kapcia, 2019) did a first-principles study of this system and found that below 80 K Phase III develops a soft phonon mode which transforms the system into an orthorhombic $F222$ #22 structure.
  - (Norman, 2019) points out that both Phase III and Phase IV structures have issues.
- Both Phase II and Phase III are extremely close to Phase I. If AFLOW-SYM allows a tolerance of 0.2 Å and FINDSYM allows a tolerance of 0.2 Å for both the lattice vectors and atomic positions both of the tetragonal phases become cubic.
- Phase IV is extremely close to Phase II. If AFLOW-SYM allows a tolerance of 0.002 Å and FINDSYM allows a tolerance 0.002 Å for both the lattice vectors and atomic positions the orthorhombic phase becomes tetragonal.
- Data for the Phase II structure was taken at 160 K.

Body-centered Tetragonal primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= -\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}\\ \mathbf{a}_2 &= \tfrac{1}{2}a\,\hat{\mathbf{x}} - \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}\\ \mathbf{a}_3 &= \tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} - \tfrac{1}{2}c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{2}\mathbf{a}_1 + \frac{1}{2}\mathbf{a}_2$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	(2*b*)	O I
$\mathbf{B}_2$	=	$\frac{1}{4}\mathbf{a}_1 + \frac{3}{4}\mathbf{a}_2 + \frac{1}{2}\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(2*d*)	O II
$\mathbf{B}_3$	=	$z_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_2$	=	$z_3 c\,\hat{\mathbf{z}}$	(4*e*)	O III
$\mathbf{B}_4$	=	$-z_3\,\mathbf{a}_1 - z_3\,\mathbf{a}_2$	=	$-z_3 c\,\hat{\mathbf{z}}$	(4*e*)	O III
$\mathbf{B}_5$	=	$\left(\frac{1}{2}+z_4\right)\mathbf{a}_1 + z_4\,\mathbf{a}_2 + \frac{1}{2}\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(4*f*)	O IV
$\mathbf{B}_6$	=	$-z_4\,\mathbf{a}_1 + \left(\frac{1}{2}-z_4\right)\mathbf{a}_2 + \frac{1}{2}\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - z_4 c\,\hat{\mathbf{z}}$	(4*f*)	O IV
$\mathbf{B}_7$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + 2x_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}}$	(8*g*)	O V
$\mathbf{B}_8$	=	$-x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - 2x_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}}$	(8*g*)	O V
$\mathbf{B}_9$	=	$-x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2$	=	$x_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}}$	(8*g*)	O V
$\mathbf{B}_{10}$	=	$x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2$	=	$-x_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}}$	(8*g*)	O V
$\mathbf{B}_{11}$	=	$\left(\frac{3}{4}+x_6\right)\mathbf{a}_1 + \left(\frac{1}{4}+x_6\right)\mathbf{a}_2 + \left(\frac{1}{2}+2x_6\right)\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(8*h*)	O VI
$\mathbf{B}_{12}$	=	$\left(\frac{3}{4}-x_6\right)\mathbf{a}_1 + \left(\frac{1}{4}-x_6\right)\mathbf{a}_2 + \left(\frac{1}{2}-2x_6\right)\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(8*h*)	O VI
$\mathbf{B}_{13}$	=	$\left(\frac{3}{4}-x_6\right)\mathbf{a}_1 + \left(\frac{1}{4}+x_6\right)\mathbf{a}_2 + \frac{1}{2}\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(8*h*)	O VI
$\mathbf{B}_{14}$	=	$\left(\frac{3}{4}+x_6\right)\mathbf{a}_1 + \left(\frac{1}{4}-x_6\right)\mathbf{a}_2 + \frac{1}{2}\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(8*h*)	O VI
$\mathbf{B}_{15}$	=	$z_7\,\mathbf{a}_1 + (x_7+z_7)\,\mathbf{a}_2 + x_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + z_7 c\,\hat{\mathbf{z}}$	(8*i*)	Cd
$\mathbf{B}_{16}$	=	$z_7\,\mathbf{a}_1 + (-x_7+z_7)\,\mathbf{a}_2 - x_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} + z_7 c\,\hat{\mathbf{z}}$	(8*i*)	Cd
$\mathbf{B}_{17}$	=	$(-x_7-z_7)\,\mathbf{a}_1 - z_7\,\mathbf{a}_2 - x_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	(8*i*)	Cd
$\mathbf{B}_{18}$	=	$(x_7-z_7)\,\mathbf{a}_1 - z_7\,\mathbf{a}_2 + x_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	(8*i*)	Cd
$\mathbf{B}_{19}$	=	$z_8\,\mathbf{a}_1 + (x_8+z_8)\,\mathbf{a}_2 + x_8\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + z_8 c\,\hat{\mathbf{z}}$	(8*i*)	Re
$\mathbf{B}_{20}$	=	$z_8\,\mathbf{a}_1 + (-x_8+z_8)\,\mathbf{a}_2 - x_8\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} + z_8 c\,\hat{\mathbf{z}}$	(8*i*)	Re
$\mathbf{B}_{21}$	=	$(-x_8-z_8)\,\mathbf{a}_1 - z_8\,\mathbf{a}_2 - x_8\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(8*i*)	Re
$\mathbf{B}_{22}$	=	$(x_8-z_8)\,\mathbf{a}_1 - z_8\,\mathbf{a}_2 + x_8\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(8*i*)	Re

References:

- S.-W. Huang, H.-T. Jeng, J.-Y. Lin, W. J. Chang, J. M. Chen, G. H. Lee, H. Berger, H. D. Yang, and K. S. Liang, *Electronic structure of pyrochlore $Cd_2Re_2O_7$*, J. Phys.: Condens. Matter **21**, 195602 (2009), doi:10.1088/0953-8984/21/19/195602.
- K. J. Kapcia, M. Reedyk, M. Hajialamdari, A. Ptok, P. Piekarz, F. S. Razavi, A. M. Oleś, and R. K. Kremer, *Discovery of a low-temperature orthorhombic phase of the $Cd_2Re_2O_7$ superconductor*, Phys. Rev. Research **2**, 033108 (2020), doi:10.1103/PhysRevResearch.2.033108.

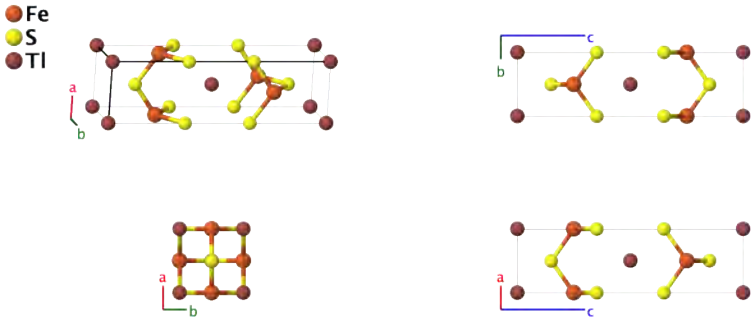
Found in:

- M. R. Norman, *The crystal structure of the inversion breaking metal* $Cd_2Re_2O_7$, Phys. Rev. B **101**, 045117 (2020), doi:10.1103/PhysRevB.101.045117.

Geometry files:
- CIF: pp. 1683
- POSCAR: pp. 1683

Tetragonal $TlFeS_2$ Structure: AB2C_tI8_119_c_e_a

http://aflow.org/prototype-encyclopedia/AB2C_tI8_119_c_e_a



Prototype	:	FeS_2Tl
AFLOW prototype label	:	AB2C_tI8_119_c_e_a
***Strukturbericht* designation**	:	None
Pearson symbol	:	tI8
Space group number	:	119
Space group symbol	:	$I\bar{4}m2$
AFLOW prototype command	:	`aflow --proto=AB2C_tI8_119_c_e_a --params=`$a, c/a, z_3$

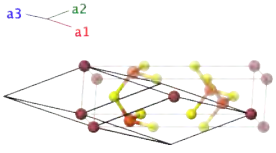
Other compounds with this structure

- $CsFeSe_2$, $KFeSe_2$, $RbFeSe_2$, and $TlFeSe_2$

- This is the high-temperature form of $TlFeS_2$/$TlFeSe_2$, formed above 300 °C. Below that temperature, it transforms into the monoclinic $TlFeSe_2$ structure.

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} - \frac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	Tl
$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(2c)$	Fe
$\mathbf{B}_3$	=	$z_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_2$	=	$z_3 c\,\hat{\mathbf{z}}$	$(4e)$	S
$\mathbf{B}_4$	=	$-z_3\,\mathbf{a}_1 - z_3\,\mathbf{a}_2$	=	$-z_3 c\,\hat{\mathbf{z}}$	$(4e)$	S

References:

- A. Kutoglu, *Synthese und Kristallstrukturen von $TlFeS_2$ und $TlFeSe_2$*, Naturwissenschaften **61**, 125–126 (1974), doi:10.1007/BF00606283.

Found in:

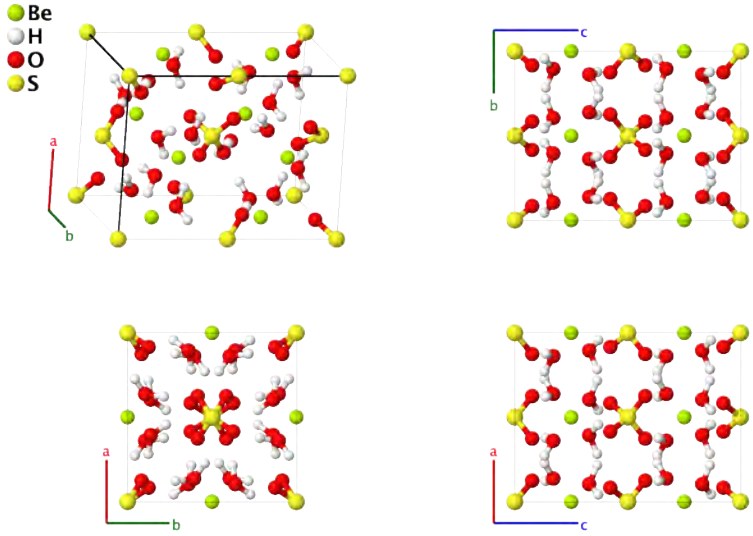
- E. B. Asgerov, A. I. Madadzada, A. I. Beskrovnyy, D. I. Ismayilov, R. N. Mehdieva, S. H. Jabarov, E. M. Kerimova, and D. Neov, *Neutron-Diffraction Study in $TlFeS_2$ and $TlFeSe_2$ at Low Temperatures*, J. Surf. Invest.: X-Ray, Synchrotron Neutron Tech. **8**, 1193–1197 (2014), doi:10.1134/S1027451014060238.

Geometry files:

- CIF: pp. 1683
- POSCAR: pp. 1684

$BeSO_4{\cdot}4H_2O$ ($H4_3$) Structure: AB8C8D_tI72_120_c_2i_2i_b

http://aflow.org/prototype-encyclopedia/AB8C8D_tI72_120_c_2i_2i_b

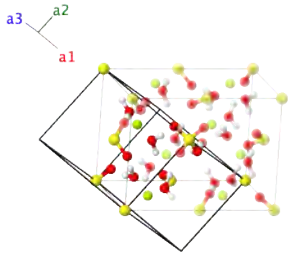


Prototype	:	BeH_8O_8S
AFLOW prototype label	:	AB8C8D_tI72_120_c_2i_2i_b
***Strukturbericht* designation**	:	$H4_3$
Pearson symbol	:	tI72
Space group number	:	120
Space group symbol	:	$I\bar{4}c2$
AFLOW prototype command	:	`aflow --proto=AB8C8D_tI72_120_c_2i_2i_b --params=`$a, c/a, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6$

- The original determination of the $H4_3$ structure did not locate the positions of the hydrogen atoms. Since (Sikka, 1969) showed that the placement of the hydrogen atoms did not substantially affect the positions of the other atoms in the primitive cell, nor change the space group, we retain the original *Strukturbericht* designation for the improved structure.

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}a\,\hat{\mathbf{x}} - \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} - \tfrac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($4b$)	S

$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	(4*b*)	S
$\mathbf{B}_3$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*c*)	Be
$\mathbf{B}_4$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*c*)	Be
$\mathbf{B}_5$	=	$(y_3 + z_3)\,\mathbf{a}_1 + (x_3 + z_3)\,\mathbf{a}_2 + (x_3 + y_3)\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + y_3a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	(16*i*)	H I
$\mathbf{B}_6$	=	$(-y_3 + z_3)\,\mathbf{a}_1 + (-x_3 + z_3)\,\mathbf{a}_2 + (-x_3 - y_3)\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - y_3a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	(16*i*)	H I
$\mathbf{B}_7$	=	$(-x_3 - z_3)\,\mathbf{a}_1 + (y_3 - z_3)\,\mathbf{a}_2 + (-x_3 + y_3)\,\mathbf{a}_3$	=	$y_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	(16*i*)	H I
$\mathbf{B}_8$	=	$(x_3 - z_3)\,\mathbf{a}_1 + (-y_3 - z_3)\,\mathbf{a}_2 + (x_3 - y_3)\,\mathbf{a}_3$	=	$-y_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	(16*i*)	H I
$\mathbf{B}_9$	=	$\left(\frac{1}{2} - y_3 + z_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_3 + z_3\right)\mathbf{a}_2 + (x_3 - y_3)\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} - y_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(16*i*)	H I
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} + y_3 + z_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_3 + z_3\right)\mathbf{a}_2 + (-x_3 + y_3)\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + y_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(16*i*)	H I
$\mathbf{B}_{11}$	=	$\left(\frac{1}{2} + x_3 - z_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3 - z_3\right)\mathbf{a}_2 + (x_3 + y_3)\,\mathbf{a}_3$	=	$y_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(16*i*)	H I
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} - x_3 - z_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3 - z_3\right)\mathbf{a}_2 + (-x_3 - y_3)\,\mathbf{a}_3$	=	$-y_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(16*i*)	H I
$\mathbf{B}_{13}$	=	$(y_4 + z_4)\,\mathbf{a}_1 + (x_4 + z_4)\,\mathbf{a}_2 + (x_4 + y_4)\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + y_4a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	(16*i*)	H II
$\mathbf{B}_{14}$	=	$(-y_4 + z_4)\,\mathbf{a}_1 + (-x_4 + z_4)\,\mathbf{a}_2 + (-x_4 - y_4)\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} - y_4a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	(16*i*)	H II
$\mathbf{B}_{15}$	=	$(-x_4 - z_4)\,\mathbf{a}_1 + (y_4 - z_4)\,\mathbf{a}_2 + (-x_4 + y_4)\,\mathbf{a}_3$	=	$y_4a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	(16*i*)	H II
$\mathbf{B}_{16}$	=	$(x_4 - z_4)\,\mathbf{a}_1 + (-y_4 - z_4)\,\mathbf{a}_2 + (x_4 - y_4)\,\mathbf{a}_3$	=	$-y_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	(16*i*)	H II
$\mathbf{B}_{17}$	=	$\left(\frac{1}{2} - y_4 + z_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_4 + z_4\right)\mathbf{a}_2 + (x_4 - y_4)\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} - y_4a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(16*i*)	H II
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} + y_4 + z_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_4 + z_4\right)\mathbf{a}_2 + (-x_4 + y_4)\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + y_4a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(16*i*)	H II
$\mathbf{B}_{19}$	=	$\left(\frac{1}{2} + x_4 - z_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4 - z_4\right)\mathbf{a}_2 + (x_4 + y_4)\,\mathbf{a}_3$	=	$y_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(16*i*)	H II
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} - x_4 - z_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4 - z_4\right)\mathbf{a}_2 + (-x_4 - y_4)\,\mathbf{a}_3$	=	$-y_4a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(16*i*)	H II
$\mathbf{B}_{21}$	=	$(y_5 + z_5)\,\mathbf{a}_1 + (x_5 + z_5)\,\mathbf{a}_2 + (x_5 + y_5)\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}} + y_5a\,\hat{\mathbf{y}} + z_5c\,\hat{\mathbf{z}}$	(16*i*)	O I
$\mathbf{B}_{22}$	=	$(-y_5 + z_5)\,\mathbf{a}_1 + (-x_5 + z_5)\,\mathbf{a}_2 + (-x_5 - y_5)\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}} - y_5a\,\hat{\mathbf{y}} + z_5c\,\hat{\mathbf{z}}$	(16*i*)	O I
$\mathbf{B}_{23}$	=	$(-x_5 - z_5)\,\mathbf{a}_1 + (y_5 - z_5)\,\mathbf{a}_2 + (-x_5 + y_5)\,\mathbf{a}_3$	=	$y_5a\,\hat{\mathbf{x}} - x_5a\,\hat{\mathbf{y}} - z_5c\,\hat{\mathbf{z}}$	(16*i*)	O I
$\mathbf{B}_{24}$	=	$(x_5 - z_5)\,\mathbf{a}_1 + (-y_5 - z_5)\,\mathbf{a}_2 + (x_5 - y_5)\,\mathbf{a}_3$	=	$-y_5a\,\hat{\mathbf{x}} + x_5a\,\hat{\mathbf{y}} - z_5c\,\hat{\mathbf{z}}$	(16*i*)	O I
$\mathbf{B}_{25}$	=	$\left(\frac{1}{2} - y_5 + z_5\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_5 + z_5\right)\mathbf{a}_2 + (x_5 - y_5)\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}} - y_5a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	(16*i*)	O I
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} + y_5 + z_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_5 + z_5\right)\mathbf{a}_2 + (-x_5 + y_5)\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}} + y_5a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	(16*i*)	O I
$\mathbf{B}_{27}$	=	$\left(\frac{1}{2} + x_5 - z_5\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_5 - z_5\right)\mathbf{a}_2 + (x_5 + y_5)\,\mathbf{a}_3$	=	$y_5a\,\hat{\mathbf{x}} + x_5a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	(16*i*)	O I
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} - x_5 - z_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_5 - z_5\right)\mathbf{a}_2 + (-x_5 - y_5)\,\mathbf{a}_3$	=	$-y_5a\,\hat{\mathbf{x}} - x_5a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	(16*i*)	O I
$\mathbf{B}_{29}$	=	$(y_6 + z_6)\,\mathbf{a}_1 + (x_6 + z_6)\,\mathbf{a}_2 + (x_6 + y_6)\,\mathbf{a}_3$	=	$x_6a\,\hat{\mathbf{x}} + y_6a\,\hat{\mathbf{y}} + z_6c\,\hat{\mathbf{z}}$	(16*i*)	O II

$\mathbf{B}_{30}$	=	$(-y_6 + z_6)\ \mathbf{a}_1 + (-x_6 + z_6)\ \mathbf{a}_2 + (-x_6 - y_6)\ \mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - y_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(16*i*)	O II
$\mathbf{B}_{31}$	=	$(-x_6 - z_6)\ \mathbf{a}_1 + (y_6 - z_6)\ \mathbf{a}_2 + (-x_6 + y_6)\ \mathbf{a}_3$	=	$y_6 a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(16*i*)	O II
$\mathbf{B}_{32}$	=	$(x_6 - z_6)\ \mathbf{a}_1 + (-y_6 - z_6)\ \mathbf{a}_2 + (x_6 - y_6)\ \mathbf{a}_3$	=	$-y_6 a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(16*i*)	O II
$\mathbf{B}_{33}$	=	$\left(\frac{1}{2} - y_6 + z_6\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_6 + z_6\right)\mathbf{a}_2 + (x_6 - y_6)\ \mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} - y_6 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	(16*i*)	O II
$\mathbf{B}_{34}$	=	$\left(\frac{1}{2} + y_6 + z_6\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_6 + z_6\right)\mathbf{a}_2 + (-x_6 + y_6)\ \mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + y_6 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	(16*i*)	O II
$\mathbf{B}_{35}$	=	$\left(\frac{1}{2} + x_6 - z_6\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_6 - z_6\right)\mathbf{a}_2 + (x_6 + y_6)\ \mathbf{a}_3$	=	$y_6 a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right) c\,\hat{\mathbf{z}}$	(16*i*)	O II
$\mathbf{B}_{36}$	=	$\left(\frac{1}{2} - x_6 - z_6\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_6 - z_6\right)\mathbf{a}_2 + (-x_6 - y_6)\ \mathbf{a}_3$	=	$-y_6 a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right) c\,\hat{\mathbf{z}}$	(16*i*)	O II

References:

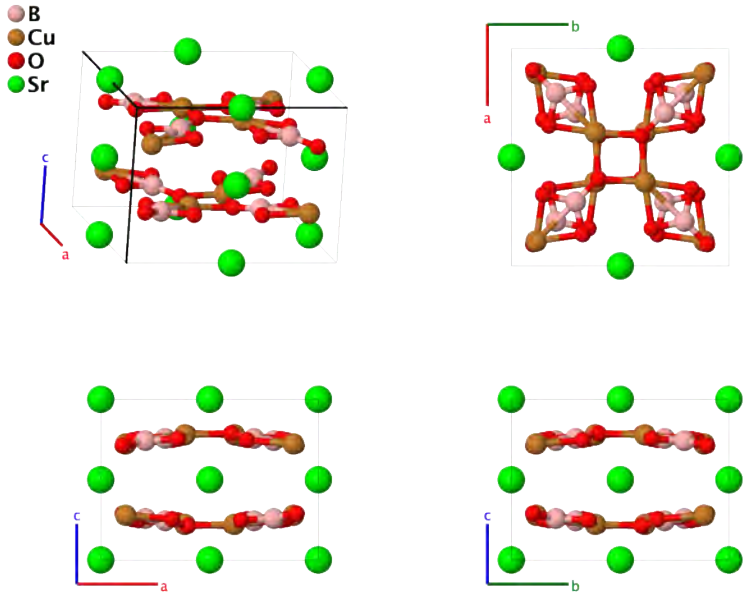
- S. K. Sikka and R. Chidambaram, *A neutron diffraction determination of the structure of beryllium sulphate tetrahydrate, $BeSO_4{\cdot}4H_2O$*, Acta Crystallogr. Sect. B Struct. Sci. **25**, 310–315 (1969), doi:10.1107/S0567740869002160.

Geometry files:

- CIF: pp. 1684
- POSCAR: pp. 1684

$SrCu_2(BO_3)_2$ Structure: A2B2C6D_tI44_121_i_i_ij_c

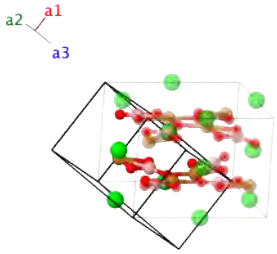
http://aflow.org/prototype-encyclopedia/A2B2C6D_tI44_121_i_i_ij_c



Prototype	:	$B_2Cu_2O_6Sr$
AFLOW prototype label	:	A2B2C6D_tI44_121_i_i_ij_c
***Strukturbericht* designation**	:	None
Pearson symbol	:	tI44
Space group number	:	121
Space group symbol	:	$I\bar{4}2m$
AFLOW prototype command	:	`aflow --proto=A2B2C6D_tI44_121_i_i_ij_c --params=`$a, c/a, x_2, z_2, x_3, z_3, x_4, z_4, x_5, y_5, z_5$

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}a\,\hat{\mathbf{x}} - \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} - \tfrac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\tfrac{1}{2}\,\mathbf{a}_1 + \tfrac{1}{2}\,\mathbf{a}_3$	=	$\tfrac{1}{2}a\,\hat{\mathbf{y}}$	(4*c*)	Sr
$\mathbf{B}_2$	=	$\tfrac{1}{2}\,\mathbf{a}_2 + \tfrac{1}{2}\,\mathbf{a}_3$	=	$\tfrac{1}{2}a\,\hat{\mathbf{x}}$	(4*c*)	Sr
$\mathbf{B}_3$	=	$(x_2 + z_2)\,\mathbf{a}_1 + (x_2 + z_2)\,\mathbf{a}_2 + 2x_2\,\mathbf{a}_3$	=	$x_2a\,\hat{\mathbf{x}} + x_2a\,\hat{\mathbf{y}} + z_2c\,\hat{\mathbf{z}}$	(8*i*)	B

$\mathbf{B}_4$	=	$(-x_2 + z_2)\,\mathbf{a}_1 + (-x_2 + z_2)\,\mathbf{a}_2 - 2x_2\,\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}} - x_2a\,\hat{\mathbf{y}} + z_2c\,\hat{\mathbf{z}}$	(8*i*)	B
$\mathbf{B}_5$	=	$(-x_2 - z_2)\,\mathbf{a}_1 + (x_2 - z_2)\,\mathbf{a}_2$	=	$x_2a\,\hat{\mathbf{x}} - x_2a\,\hat{\mathbf{y}} - z_2c\,\hat{\mathbf{z}}$	(8*i*)	B
$\mathbf{B}_6$	=	$(x_2 - z_2)\,\mathbf{a}_1 + (-x_2 - z_2)\,\mathbf{a}_2$	=	$-x_2a\,\hat{\mathbf{x}} + x_2a\,\hat{\mathbf{y}} - z_2c\,\hat{\mathbf{z}}$	(8*i*)	B
$\mathbf{B}_7$	=	$(x_3 + z_3)\,\mathbf{a}_1 + (x_3 + z_3)\,\mathbf{a}_2 + 2x_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	(8*i*)	Cu
$\mathbf{B}_8$	=	$(-x_3 + z_3)\,\mathbf{a}_1 + (-x_3 + z_3)\,\mathbf{a}_2 - 2x_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	(8*i*)	Cu
$\mathbf{B}_9$	=	$(-x_3 - z_3)\,\mathbf{a}_1 + (x_3 - z_3)\,\mathbf{a}_2$	=	$x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	(8*i*)	Cu
$\mathbf{B}_{10}$	=	$(x_3 - z_3)\,\mathbf{a}_1 + (-x_3 - z_3)\,\mathbf{a}_2$	=	$-x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	(8*i*)	Cu
$\mathbf{B}_{11}$	=	$(x_4 + z_4)\,\mathbf{a}_1 + (x_4 + z_4)\,\mathbf{a}_2 + 2x_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	(8*i*)	O I
$\mathbf{B}_{12}$	=	$(-x_4 + z_4)\,\mathbf{a}_1 + (-x_4 + z_4)\,\mathbf{a}_2 - 2x_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	(8*i*)	O I
$\mathbf{B}_{13}$	=	$(-x_4 - z_4)\,\mathbf{a}_1 + (x_4 - z_4)\,\mathbf{a}_2$	=	$x_4a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	(8*i*)	O I
$\mathbf{B}_{14}$	=	$(x_4 - z_4)\,\mathbf{a}_1 + (-x_4 - z_4)\,\mathbf{a}_2$	=	$-x_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	(8*i*)	O I
$\mathbf{B}_{15}$	=	$(y_5 + z_5)\,\mathbf{a}_1 + (x_5 + z_5)\,\mathbf{a}_2 + (x_5 + y_5)\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}} + y_5a\,\hat{\mathbf{y}} + z_5c\,\hat{\mathbf{z}}$	(16*j*)	O II
$\mathbf{B}_{16}$	=	$(-y_5 + z_5)\,\mathbf{a}_1 + (-x_5 + z_5)\,\mathbf{a}_2 + (-x_5 - y_5)\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}} - y_5a\,\hat{\mathbf{y}} + z_5c\,\hat{\mathbf{z}}$	(16*j*)	O II
$\mathbf{B}_{17}$	=	$(-x_5 - z_5)\,\mathbf{a}_1 + (y_5 - z_5)\,\mathbf{a}_2 + (-x_5 + y_5)\,\mathbf{a}_3$	=	$y_5a\,\hat{\mathbf{x}} - x_5a\,\hat{\mathbf{y}} - z_5c\,\hat{\mathbf{z}}$	(16*j*)	O II
$\mathbf{B}_{18}$	=	$(x_5 - z_5)\,\mathbf{a}_1 + (-y_5 - z_5)\,\mathbf{a}_2 + (x_5 - y_5)\,\mathbf{a}_3$	=	$-y_5a\,\hat{\mathbf{x}} + x_5a\,\hat{\mathbf{y}} - z_5c\,\hat{\mathbf{z}}$	(16*j*)	O II
$\mathbf{B}_{19}$	=	$(y_5 - z_5)\,\mathbf{a}_1 + (-x_5 - z_5)\,\mathbf{a}_2 + (-x_5 + y_5)\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}} + y_5a\,\hat{\mathbf{y}} - z_5c\,\hat{\mathbf{z}}$	(16*j*)	O II
$\mathbf{B}_{20}$	=	$(-y_5 - z_5)\,\mathbf{a}_1 + (x_5 - z_5)\,\mathbf{a}_2 + (x_5 - y_5)\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}} - y_5a\,\hat{\mathbf{y}} - z_5c\,\hat{\mathbf{z}}$	(16*j*)	O II
$\mathbf{B}_{21}$	=	$(-x_5 + z_5)\,\mathbf{a}_1 + (-y_5 + z_5)\,\mathbf{a}_2 + (-x_5 - y_5)\,\mathbf{a}_3$	=	$-y_5a\,\hat{\mathbf{x}} - x_5a\,\hat{\mathbf{y}} + z_5c\,\hat{\mathbf{z}}$	(16*j*)	O II
$\mathbf{B}_{22}$	=	$(x_5 + z_5)\,\mathbf{a}_1 + (y_5 + z_5)\,\mathbf{a}_2 + (x_5 + y_5)\,\mathbf{a}_3$	=	$y_5a\,\hat{\mathbf{x}} + x_5a\,\hat{\mathbf{y}} + z_5c\,\hat{\mathbf{z}}$	(16*j*)	O II

References:

- R. W. Smith and D. A. Keszler, *Synthesis, structure, and properties of the orthoborate $SrCu_2(BO_3)_2$*, J. Solid State Chem. **93**, 430–435 (1991), doi:10.1016/0022-4596(91)90316-A.

Found in:

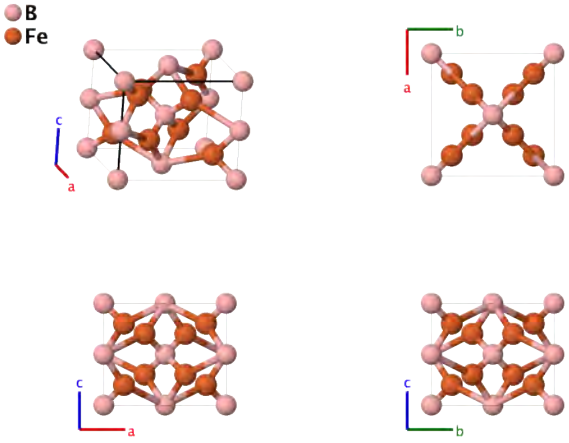
- H. Kageyama, K. Yoshimura, R. Stern, N. V. Mushnikov, K. Onizuka, M. Kato, K. Kosuge, C. P. Slichter, T. Goto, and Y. Ueda, *Exact Dimer Ground State and Quantized Magnetization Plateaus in the Two-Dimensional Spin System $SrCu_2(BO_3)_2$*, Phys. Rev. Lett. **82**, 3168–3171 (1999), doi:10.1103/PhysRevLett.82.3168.

Geometry files:

- CIF: pp. 1685
- POSCAR: pp. 1685

*C*17 (Fe_2B) (*obsolete*) Structure: AB2_tI12_121_ab_i

http://aflow.org/prototype-encyclopedia/AB2_tI12_121_ab_i

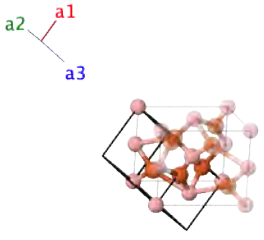


Prototype	:	BFe_2
AFLOW prototype label	:	AB2_tI12_121_ab_i
***Strukturbericht* designation**	:	*C*17
Pearson symbol	:	tI12
Space group number	:	121
Space group symbol	:	$I\bar{4}2m$
AFLOW prototype command	:	`aflow --proto=AB2_tI12_121_ab_i --params=`$a, c/a, x_3, z_3$

- (Wever, 1930) placed this structure in space group #121, and this was used to designate the structure as *C*17 in (Ewald, 1931). However, even then it was recognized that changing z_3 from the value of 0.2 assigned by Wever to 0.25 would add an inversion symmetry to the crystal, change the space group to "$I4/mcm$ #140", and make the structure look very much like Al_2Cu, *Strukturbericht* *C*16. This has become the accepted structure, as in (Kapfenberger, 2006), and so *C*17 is an obsolete designation. We retain it here for historical interest.

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}a\,\hat{\mathbf{x}} - \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} - \tfrac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(2*a*)	B I

$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	($2b$)	B II
$\mathbf{B}_3$	=	$(x_3 + z_3)\,\mathbf{a}_1 + (x_3 + z_3)\,\mathbf{a}_2 + 2x_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	($8i$)	Fe
$\mathbf{B}_4$	=	$(-x_3 + z_3)\,\mathbf{a}_1 + (-x_3 + z_3)\,\mathbf{a}_2 - 2x_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	($8i$)	Fe
$\mathbf{B}_5$	=	$(-x_3 - z_3)\,\mathbf{a}_1 + (x_3 - z_3)\,\mathbf{a}_2$	=	$x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	($8i$)	Fe
$\mathbf{B}_6$	=	$(x_3 - z_3)\,\mathbf{a}_1 + (-x_3 - z_3)\,\mathbf{a}_2$	=	$-x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	($8i$)	Fe

References:

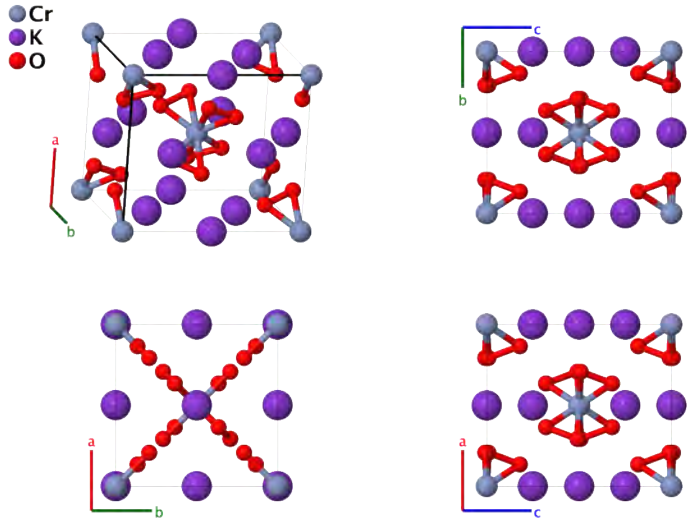
- F. Wever and A. Müller, *Über das Zweistoffsystem Eisen-Bor und über die Struktur des Eisenborides Fe_4B_2*, Z. Anorg. Allg. Chem. **192**, 317–336 (1930), doi:10.1002/zaac.19301920125.
- P. P. Ewald and C. Hermann, eds., *Strukturbericht 1913-1928* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1931).
- C. Kapfenberger, B. Albert, R. Pöttgen, and H. Huppertz, *Structure refinements of iron borides Fe_2B and FeB*, Zeitschrift für Kristallographie - Crystalline Materials **221**, 477–481 (2006), doi:10.1524/zkri.2006.221.5-7.477.

Geometry files:

- CIF: pp. 1685
- POSCAR: pp. 1685

K_3CrO_8 Structure: AB3C8_tI24_121_a_bd_2i

http://aflow.org/prototype-encyclopedia/AB3C8_tI24_121_a_bd_2i



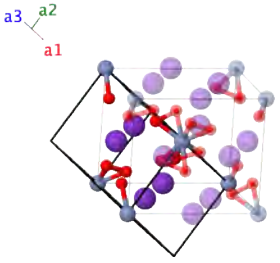
Prototype	:	CrK_3O_8
AFLOW prototype label	:	AB3C8_tI24_121_a_bd_2i
***Strukturbericht* designation**	:	None
Pearson symbol	:	tI24
Space group number	:	121
Space group symbol	:	$I\bar{4}2m$
AFLOW prototype command	:	`aflow --proto=AB3C8_tI24_121_a_bd_2i --params=`$a, c/a, x_4, z_4, x_5, z_5$

Other compounds with this structure

- Zr_3GeO_8

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}a\,\hat{\mathbf{x}} - \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} - \tfrac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(2*a*)	Cr
$\mathbf{B}_2$	=	$\tfrac{1}{2}\,\mathbf{a}_1 + \tfrac{1}{2}\,\mathbf{a}_2$	=	$\tfrac{1}{2}c\,\hat{\mathbf{z}}$	(2*b*)	K I
$\mathbf{B}_3$	=	$\tfrac{3}{4}\,\mathbf{a}_1 + \tfrac{1}{4}\,\mathbf{a}_2 + \tfrac{1}{2}\,\mathbf{a}_3$	=	$\tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{4}c\,\hat{\mathbf{z}}$	(4*d*)	K II

$\mathbf{B}_4$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*d*)	K II
$\mathbf{B}_5$	=	$(x_4 + z_4)\,\mathbf{a}_1 + (x_4 + z_4)\,\mathbf{a}_2 + 2x_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	(8*i*)	O I
$\mathbf{B}_6$	=	$(-x_4 + z_4)\,\mathbf{a}_1 + (-x_4 + z_4)\,\mathbf{a}_2 - 2x_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	(8*i*)	O I
$\mathbf{B}_7$	=	$(-x_4 - z_4)\,\mathbf{a}_1 + (x_4 - z_4)\,\mathbf{a}_2$	=	$x_4a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	(8*i*)	O I
$\mathbf{B}_8$	=	$(x_4 - z_4)\,\mathbf{a}_1 + (-x_4 - z_4)\,\mathbf{a}_2$	=	$-x_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	(8*i*)	O I
$\mathbf{B}_9$	=	$(x_5 + z_5)\,\mathbf{a}_1 + (x_5 + z_5)\,\mathbf{a}_2 + 2x_5\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}} + x_5a\,\hat{\mathbf{y}} + z_5c\,\hat{\mathbf{z}}$	(8*i*)	O II
$\mathbf{B}_{10}$	=	$(-x_5 + z_5)\,\mathbf{a}_1 + (-x_5 + z_5)\,\mathbf{a}_2 - 2x_5\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}} - x_5a\,\hat{\mathbf{y}} + z_5c\,\hat{\mathbf{z}}$	(8*i*)	O II
$\mathbf{B}_{11}$	=	$(-x_5 - z_5)\,\mathbf{a}_1 + (x_5 - z_5)\,\mathbf{a}_2$	=	$x_5a\,\hat{\mathbf{x}} - x_5a\,\hat{\mathbf{y}} - z_5c\,\hat{\mathbf{z}}$	(8*i*)	O II
$\mathbf{B}_{12}$	=	$(x_5 - z_5)\,\mathbf{a}_1 + (-x_5 - z_5)\,\mathbf{a}_2$	=	$-x_5a\,\hat{\mathbf{x}} + x_5a\,\hat{\mathbf{y}} - z_5c\,\hat{\mathbf{z}}$	(8*i*)	O II

References:

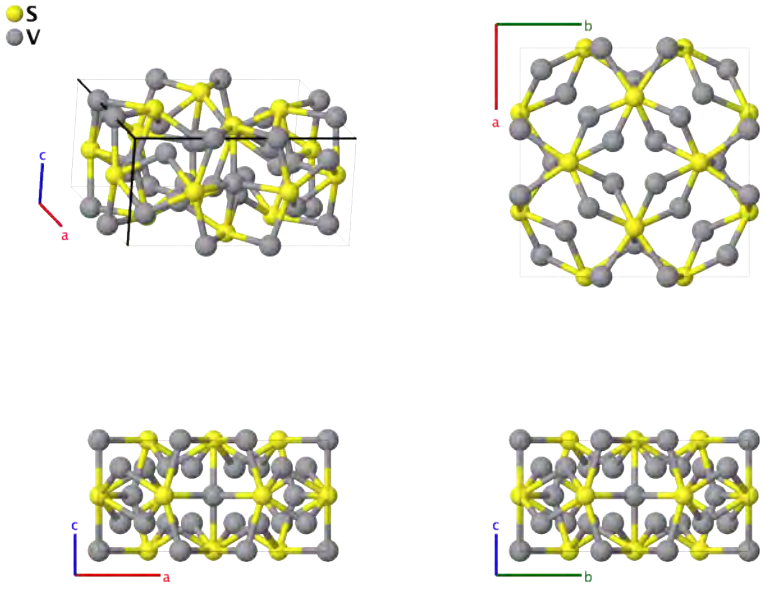
- R. Stomberg, *Least-Squares Refinement of the Crystal Structure of Potassium Peroxochromate*, Acta Chem. Scand. **17**, 1563–1566 (1963), doi:10.3891/acta.chem.scand.17-1563.

Geometry files:

- CIF: pp. 1686
- POSCAR: pp. 1686

α-V_3S Structure: AB3_tI32_121_g_f2i

http://aflow.org/prototype-encyclopedia/AB3_tI32_121_g_f2i



Prototype	:	SV_3
AFLOW prototype label	:	AB3_tI32_121_g_f2i
***Strukturbericht* designation**	:	None
Pearson symbol	:	tI32
Space group number	:	121
Space group symbol	:	$I\bar{4}2m$
AFLOW prototype command	:	`aflow --proto=AB3_tI32_121_g_f2i --params=`$a, c/a, x_1, x_2, x_3, z_3, x_4, z_4$

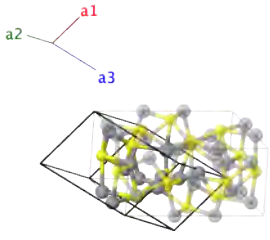
Other compounds with this structure

- Zr_3Ir

- α-V_3S is stable above 950 °C, but metastable at 25 °C, where this data was taken. Below 825 °C, the system transforms to the β-V_3S structure.

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} - \frac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_2 + x_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}}$	$(8f)$	V I
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_2 - x_1\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}}$	$(8f)$	V I
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 - x_1\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{y}}$	$(8f)$	V I
$\mathbf{B}_4$	=	$x_1\,\mathbf{a}_1 + x_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{y}}$	$(8f)$	V I
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_1 + \left(\frac{1}{2} + x_2\right)\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(8g)$	S
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_1 + \left(\frac{1}{2} - x_2\right)\,\mathbf{a}_2 - x_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(8g)$	S
$\mathbf{B}_7$	=	$\left(\frac{1}{2} - x_2\right)\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 - x_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(8g)$	S
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(8g)$	S
$\mathbf{B}_9$	=	$(x_3 + z_3)\,\mathbf{a}_1 + (x_3 + z_3)\,\mathbf{a}_2 + 2x_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(8i)$	V II
$\mathbf{B}_{10}$	=	$(-x_3 + z_3)\,\mathbf{a}_1 + (-x_3 + z_3)\,\mathbf{a}_2 - 2x_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(8i)$	V II
$\mathbf{B}_{11}$	=	$(-x_3 - z_3)\,\mathbf{a}_1 + (x_3 - z_3)\,\mathbf{a}_2$	=	$x_3 a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(8i)$	V II
$\mathbf{B}_{12}$	=	$(x_3 - z_3)\,\mathbf{a}_1 + (-x_3 - z_3)\,\mathbf{a}_2$	=	$-x_3 a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(8i)$	V II
$\mathbf{B}_{13}$	=	$(x_4 + z_4)\,\mathbf{a}_1 + (x_4 + z_4)\,\mathbf{a}_2 + 2x_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(8i)$	V III
$\mathbf{B}_{14}$	=	$(-x_4 + z_4)\,\mathbf{a}_1 + (-x_4 + z_4)\,\mathbf{a}_2 - 2x_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(8i)$	V III
$\mathbf{B}_{15}$	=	$(-x_4 - z_4)\,\mathbf{a}_1 + (x_4 - z_4)\,\mathbf{a}_2$	=	$x_4 a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	$(8i)$	V III
$\mathbf{B}_{16}$	=	$(x_4 - z_4)\,\mathbf{a}_1 + (-x_4 - z_4)\,\mathbf{a}_2$	=	$-x_4 a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	$(8i)$	V III

References:

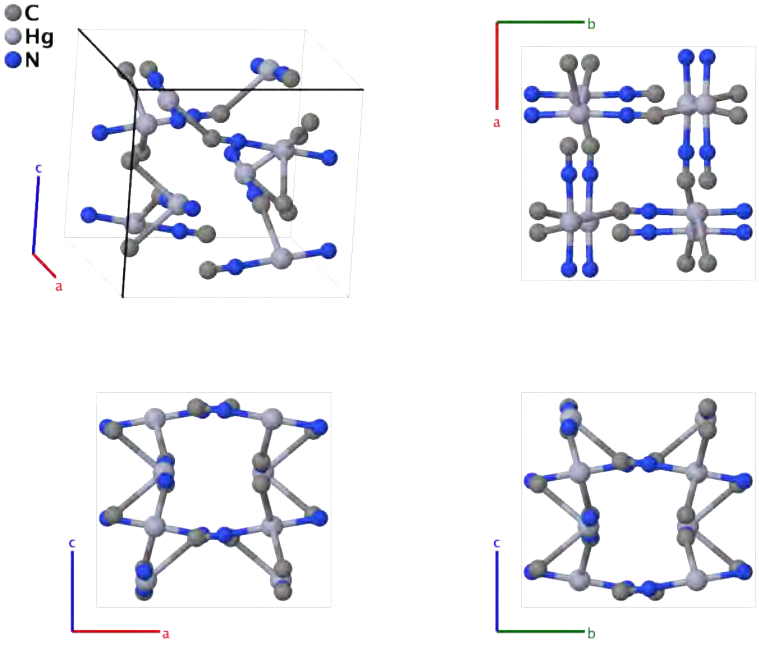
- B. Pedersen and F. Grønvold, *The Crystal Structures of α-V_3S and β-V_3S*, Acta Cryst. **12**, 1022–1027 (1959), doi:10.1107/S0365110X59002869.

Geometry files:

- CIF: pp. 1686
- POSCAR: pp. 1686

Mercury Cyanide [$Hg(CN)_2$, $F1_1$] Structure: A2BC2_tI40_122_e_d_e

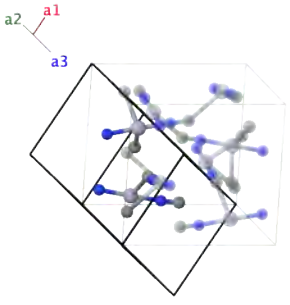
http://aflow.org/prototype-encyclopedia/A2BC2_tI40_122_e_d_e



Prototype	:	C_2HgN_2
AFLOW prototype label	:	A2BC2_tI40_122_e_d_e
***Strukturbericht* designation**	:	$F1_1$
Pearson symbol	:	tI40
Space group number	:	122
Space group symbol	:	$I\bar{4}2d$
AFLOW prototype command	:	`aflow --proto=A2BC2_tI40_122_e_d_e --params=`$a, c/a, x_1, x_2, y_2, z_2, x_3, y_3, z_3$

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}a\,\hat{\mathbf{x}} - \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} - \tfrac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$ =	$\frac{3}{8}\mathbf{a}_1 + \left(\frac{1}{8} + x_1\right)\mathbf{a}_2 + \left(\frac{1}{4} + x_1\right)\mathbf{a}_3$ =	$x_1 a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{8}c\,\hat{\mathbf{z}}$	($8d$)	Hg

$\mathbf{B}_2$	=	$\frac{7}{8}\,\mathbf{a}_1 + \left(\frac{1}{8} - x_1\right)\,\mathbf{a}_2 + \left(\frac{3}{4} - x_1\right)\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \frac{3}{4} a\,\hat{\mathbf{y}} + \frac{1}{8} c\,\hat{\mathbf{z}}$	(8*d*)	Hg
$\mathbf{B}_3$	=	$\left(\frac{7}{8} - x_1\right)\,\mathbf{a}_1 + \frac{1}{8}\,\mathbf{a}_2 + \left(\frac{1}{4} - x_1\right)\,\mathbf{a}_3$	=	$-\frac{1}{4} a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_1\right) a\,\hat{\mathbf{y}} + \frac{3}{8} c\,\hat{\mathbf{z}}$	(8*d*)	Hg
$\mathbf{B}_4$	=	$\left(\frac{7}{8} + x_1\right)\,\mathbf{a}_1 + \frac{5}{8}\,\mathbf{a}_2 + \left(\frac{3}{4} + x_1\right)\,\mathbf{a}_3$	=	$\frac{1}{4} a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_1\right) a\,\hat{\mathbf{y}} + \frac{3}{8} c\,\hat{\mathbf{z}}$	(8*d*)	Hg
$\mathbf{B}_5$	=	$(y_2 + z_2)\,\mathbf{a}_1 + (x_2 + z_2)\,\mathbf{a}_2 + (x_2 + y_2)\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + y_2 a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(16*e*)	C
$\mathbf{B}_6$	=	$(-y_2 + z_2)\,\mathbf{a}_1 + (-x_2 + z_2)\,\mathbf{a}_2 + (-x_2 - y_2)\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} - y_2 a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(16*e*)	C
$\mathbf{B}_7$	=	$(-x_2 - z_2)\,\mathbf{a}_1 + (y_2 - z_2)\,\mathbf{a}_2 + (-x_2 + y_2)\,\mathbf{a}_3$	=	$y_2 a\,\hat{\mathbf{x}} - x_2 a\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(16*e*)	C
$\mathbf{B}_8$	=	$(x_2 - z_2)\,\mathbf{a}_1 + (-y_2 - z_2)\,\mathbf{a}_2 + (x_2 - y_2)\,\mathbf{a}_3$	=	$-y_2 a\,\hat{\mathbf{x}} + x_2 a\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(16*e*)	C
$\mathbf{B}_9$	=	$\left(\frac{3}{4} + y_2 - z_2\right)\,\mathbf{a}_1 + \left(\frac{1}{4} - x_2 - z_2\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - x_2 + y_2\right)\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_2\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_2\right) c\,\hat{\mathbf{z}}$	(16*e*)	C
$\mathbf{B}_{10}$	=	$\left(\frac{3}{4} - y_2 - z_2\right)\,\mathbf{a}_1 + \left(\frac{1}{4} + x_2 - z_2\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + x_2 - y_2\right)\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_2\right) c\,\hat{\mathbf{z}}$	(16*e*)	C
$\mathbf{B}_{11}$	=	$\left(\frac{3}{4} - x_2 + z_2\right)\,\mathbf{a}_1 + \left(\frac{1}{4} - y_2 + z_2\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - x_2 - y_2\right)\,\mathbf{a}_3$	=	$-y_2 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_2\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_2\right) c\,\hat{\mathbf{z}}$	(16*e*)	C
$\mathbf{B}_{12}$	=	$\left(\frac{3}{4} + x_2 + z_2\right)\,\mathbf{a}_1 + \left(\frac{1}{4} + y_2 + z_2\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + x_2 + y_2\right)\,\mathbf{a}_3$	=	$y_2 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_2\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_2\right) c\,\hat{\mathbf{z}}$	(16*e*)	C
$\mathbf{B}_{13}$	=	$(y_3 + z_3)\,\mathbf{a}_1 + (x_3 + z_3)\,\mathbf{a}_2 + (x_3 + y_3)\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(16*e*)	N
$\mathbf{B}_{14}$	=	$(-y_3 + z_3)\,\mathbf{a}_1 + (-x_3 + z_3)\,\mathbf{a}_2 + (-x_3 - y_3)\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(16*e*)	N
$\mathbf{B}_{15}$	=	$(-x_3 - z_3)\,\mathbf{a}_1 + (y_3 - z_3)\,\mathbf{a}_2 + (-x_3 + y_3)\,\mathbf{a}_3$	=	$y_3 a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(16*e*)	N
$\mathbf{B}_{16}$	=	$(x_3 - z_3)\,\mathbf{a}_1 + (-y_3 - z_3)\,\mathbf{a}_2 + (x_3 - y_3)\,\mathbf{a}_3$	=	$-y_3 a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(16*e*)	N
$\mathbf{B}_{17}$	=	$\left(\frac{3}{4} + y_3 - z_3\right)\,\mathbf{a}_1 + \left(\frac{1}{4} - x_3 - z_3\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - x_3 + y_3\right)\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_3\right) c\,\hat{\mathbf{z}}$	(16*e*)	N
$\mathbf{B}_{18}$	=	$\left(\frac{3}{4} - y_3 - z_3\right)\,\mathbf{a}_1 + \left(\frac{1}{4} + x_3 - z_3\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + x_3 - y_3\right)\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_3\right) c\,\hat{\mathbf{z}}$	(16*e*)	N
$\mathbf{B}_{19}$	=	$\left(\frac{3}{4} - x_3 + z_3\right)\,\mathbf{a}_1 + \left(\frac{1}{4} - y_3 + z_3\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - x_3 - y_3\right)\,\mathbf{a}_3$	=	$-y_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_3\right) c\,\hat{\mathbf{z}}$	(16*e*)	N
$\mathbf{B}_{20}$	=	$\left(\frac{3}{4} + x_3 + z_3\right)\,\mathbf{a}_1 + \left(\frac{1}{4} + y_3 + z_3\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + x_3 + y_3\right)\,\mathbf{a}_3$	=	$y_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_3\right) c\,\hat{\mathbf{z}}$	(16*e*)	N

References:

- O. Reckeweg and A. Simon, *X-Ray and Raman Investigations on Cyanides of Mono- and Divalent Metals and Synthesis, Crystal Structure and Raman Spectrum of $Tl_5(CO_3)_2(CN)$*, Z. Naturforsch. B **57**, 895–900 (2002), doi:10.1515/znb-2002-0809.

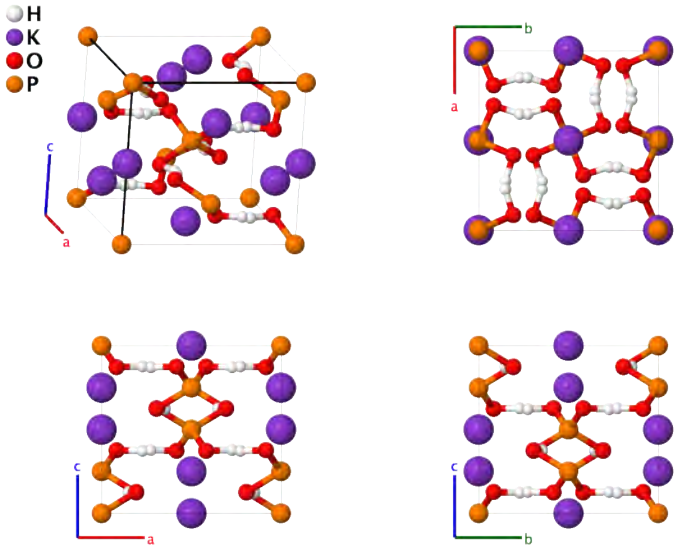
Found in:

- P. Villars (Chief Editor), *$Hg(CN)_2$ ($Hg[CN]_2$) Crystal Structure*, http://materials.springer.com/isp/crystallographic/docs/sd_1415375 (2016). PAULING FILE in: Inorganic Solid Phases, SpringerMaterials (online database).

Geometry files:
- CIF: pp. 1687
- POSCAR: pp. 1687

KH_2PO_4 ($H2_2$) Structure: A4BC4D_tI40_122_e_b_e_a

http://aflow.org/prototype-encyclopedia/A4BC4D_tI40_122_e_b_e_a



Prototype	:	H_2KO_4P
AFLOW prototype label	:	A4BC4D_tI40_122_e_b_e_a
***Strukturbericht* designation**	:	$H2_2$
Pearson symbol	:	tI40
Space group number	:	122
Space group symbol	:	$I\bar{4}2d$
AFLOW prototype command	:	`aflow --proto=A4BC4D_tI40_122_e_b_e_a --params=`$a, c/a, x_3, y_3, z_3, x_4, y_4, z_4$

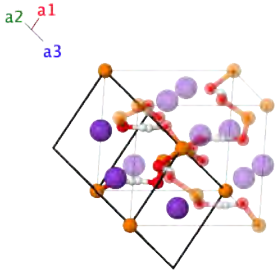
Other compounds with this structure

- RbH_2PO_4, KH_2AsO_4, RbH_2AsO_4, CsH_2AsO_4, $NH_4H_2PO_4$, and $NH_4H_2AsO_4$

- The hydrogen ($16e$) sites are half-occupied. Given the closeness of pairs of hydrogen positions, presumably only one site in each pair is ever occupied.
- This partial occupancy gives a structure that differs slightly from the $H2_2$ structure described by (Ewald, 1928). There, the hydrogen atoms are at their averaged positions, Wyckoff position ($8d$).

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} - \frac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(4*a*)	P
$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*a*)	P
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	(4*b*)	K
$\mathbf{B}_4$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*b*)	K
$\mathbf{B}_5$	=	$(y_3 + z_3)\,\mathbf{a}_1 + (x_3 + z_3)\,\mathbf{a}_2 + (x_3 + y_3)\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + y_3a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	(16*e*)	H
$\mathbf{B}_6$	=	$(-y_3 + z_3)\,\mathbf{a}_1 + (-x_3 + z_3)\,\mathbf{a}_2 + (-x_3 - y_3)\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - y_3a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	(16*e*)	H
$\mathbf{B}_7$	=	$(-x_3 - z_3)\,\mathbf{a}_1 + (y_3 - z_3)\,\mathbf{a}_2 + (-x_3 + y_3)\,\mathbf{a}_3$	=	$y_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	(16*e*)	H
$\mathbf{B}_8$	=	$(x_3 - z_3)\,\mathbf{a}_1 + (-y_3 - z_3)\,\mathbf{a}_2 + (x_3 - y_3)\,\mathbf{a}_3$	=	$-y_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	(16*e*)	H
$\mathbf{B}_9$	=	$\left(\frac{3}{4} + y_3 - z_3\right)\mathbf{a}_1 + \left(\frac{1}{4} - x_3 - z_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_3 + y_3\right)\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_3\right)c\,\hat{\mathbf{z}}$	(16*e*)	H
$\mathbf{B}_{10}$	=	$\left(\frac{3}{4} - y_3 - z_3\right)\mathbf{a}_1 + \left(\frac{1}{4} + x_3 - z_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_3 - y_3\right)\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_3\right)c\,\hat{\mathbf{z}}$	(16*e*)	H
$\mathbf{B}_{11}$	=	$\left(\frac{3}{4} - x_3 + z_3\right)\mathbf{a}_1 + \left(\frac{1}{4} - y_3 + z_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_3 - y_3\right)\mathbf{a}_3$	=	$-y_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_3\right)c\,\hat{\mathbf{z}}$	(16*e*)	H
$\mathbf{B}_{12}$	=	$\left(\frac{3}{4} + x_3 + z_3\right)\mathbf{a}_1 + \left(\frac{1}{4} + y_3 + z_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_3 + y_3\right)\mathbf{a}_3$	=	$y_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_3\right)c\,\hat{\mathbf{z}}$	(16*e*)	H
$\mathbf{B}_{13}$	=	$(y_4 + z_4)\,\mathbf{a}_1 + (x_4 + z_4)\,\mathbf{a}_2 + (x_4 + y_4)\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + y_4a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	(16*e*)	O
$\mathbf{B}_{14}$	=	$(-y_4 + z_4)\,\mathbf{a}_1 + (-x_4 + z_4)\,\mathbf{a}_2 + (-x_4 - y_4)\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} - y_4a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	(16*e*)	O
$\mathbf{B}_{15}$	=	$(-x_4 - z_4)\,\mathbf{a}_1 + (y_4 - z_4)\,\mathbf{a}_2 + (-x_4 + y_4)\,\mathbf{a}_3$	=	$y_4a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	(16*e*)	O
$\mathbf{B}_{16}$	=	$(x_4 - z_4)\,\mathbf{a}_1 + (-y_4 - z_4)\,\mathbf{a}_2 + (x_4 - y_4)\,\mathbf{a}_3$	=	$-y_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	(16*e*)	O
$\mathbf{B}_{17}$	=	$\left(\frac{3}{4} + y_4 - z_4\right)\mathbf{a}_1 + \left(\frac{1}{4} - x_4 - z_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_4 + y_4\right)\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_4\right)c\,\hat{\mathbf{z}}$	(16*e*)	O
$\mathbf{B}_{18}$	=	$\left(\frac{3}{4} - y_4 - z_4\right)\mathbf{a}_1 + \left(\frac{1}{4} + x_4 - z_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_4 - y_4\right)\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_4\right)c\,\hat{\mathbf{z}}$	(16*e*)	O
$\mathbf{B}_{19}$	=	$\left(\frac{3}{4} - x_4 + z_4\right)\mathbf{a}_1 + \left(\frac{1}{4} - y_4 + z_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_4 - y_4\right)\mathbf{a}_3$	=	$-y_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_4\right)c\,\hat{\mathbf{z}}$	(16*e*)	O
$\mathbf{B}_{20}$	=	$\left(\frac{3}{4} + x_4 + z_4\right)\mathbf{a}_1 + \left(\frac{1}{4} + y_4 + z_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_4 + y_4\right)\mathbf{a}_3$	=	$y_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_4\right)c\,\hat{\mathbf{z}}$	(16*e*)	O

References:

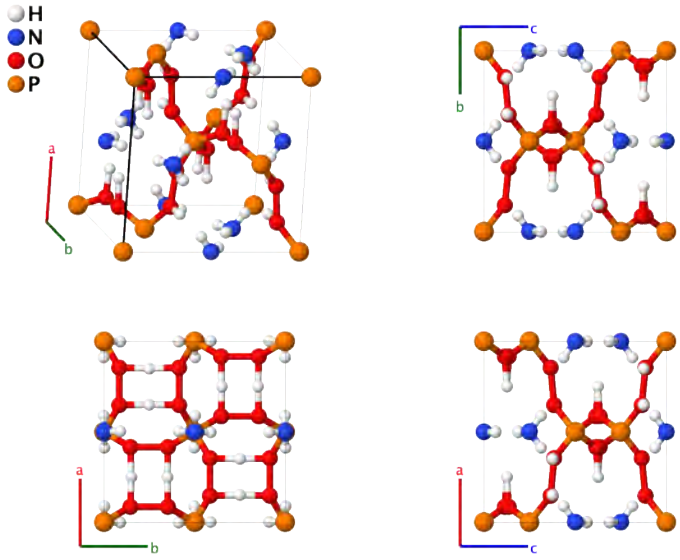
- R. J. Nelmes, G. M. Meyer, and J. E. Tibballs, *The crystal structure of tetragonal KH_2PO_4 and KD_2PO_4 as a function of temperature*, J. Phys. C: Solid State Phys. **15**, 59–75 (1982), doi:10.1088/0022-3719/15/1/005. Corrigendum: R. J. Nelmes and G. M. Meyer and J. E. Tibballs, J. Phys. C 15, 3040 (1982).
- R. J. Nelmes, G. M. Meyer, and J. E. Tibballs, *The crystal structure of tetragonal KH_2PO_4 and KD_2PO_4 as a function of temperature*, J. Phys. C: Solid State Phys. **15**, 3040 (1982), doi:10.1088/0022-3719/15/13/531. Corrigendum to R. J.

Nelmes and G. M. Meyer and J. E. Tibballs, J. Phys. C 15, 59 (1982).
- P. P. Ewald and C. Hermann, eds., *Strukturbericht 1913-1928* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1931).

Geometry files:
- CIF: pp. 1687
- POSCAR: pp. 1688

$NH_4H_2PO_4$ Structure: A8BC4D_tI56_122_2e_b_e_a

http://aflow.org/prototype-encyclopedia/A8BC4D_tI56_122_2e_b_e_a



Prototype	:	H_6NO_4P
AFLOW prototype label	:	A8BC4D_tI56_122_2e_b_e_a
***Strukturbericht* designation**	:	None
Pearson symbol	:	tI56
Space group number	:	122
Space group symbol	:	$I\bar{4}2d$
AFLOW prototype command	:	`aflow --proto=A8BC4D_tI56_122_2e_b_e_a --params=`$a, c/a, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5$

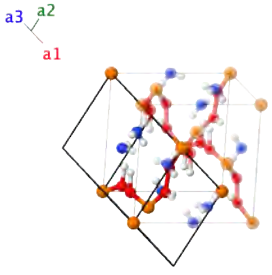
Other compounds with this structure

- $NH_4H_2AsO_4$

- $NH_4H_2PO_4$ and $NH_4H_2AsO_4$ are usually considered to be isomorphous with the $H2_2$ KH_2PO_4 structure, but (Khan, 1973) and (Fukami, 1987) were able to locate the hydrogen atoms in the NH_4 ion, so we include this as a new structure.
- As in KH_2PO_4, the H-I site, which is associated with the PO_4 ion, is 50% occupied.
- Below 148 K the H-I atoms become locked in place, and $NH_4H_2PO_4$ distorts in to a orthorhombic ferroelectric phase.

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} - \frac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(4*a*)	P
$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*a*)	P
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	(4*b*)	N
$\mathbf{B}_4$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*b*)	N
$\mathbf{B}_5$	=	$(y_3 + z_3)\,\mathbf{a}_1 + (x_3 + z_3)\,\mathbf{a}_2 + (x_3 + y_3)\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + y_3a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	(16*e*)	H I
$\mathbf{B}_6$	=	$(-y_3 + z_3)\,\mathbf{a}_1 + (-x_3 + z_3)\,\mathbf{a}_2 + (-x_3 - y_3)\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - y_3a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	(16*e*)	H I
$\mathbf{B}_7$	=	$(-x_3 - z_3)\,\mathbf{a}_1 + (y_3 - z_3)\,\mathbf{a}_2 + (-x_3 + y_3)\,\mathbf{a}_3$	=	$y_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	(16*e*)	H I
$\mathbf{B}_8$	=	$(x_3 - z_3)\,\mathbf{a}_1 + (-y_3 - z_3)\,\mathbf{a}_2 + (x_3 - y_3)\,\mathbf{a}_3$	=	$-y_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	(16*e*)	H I
$\mathbf{B}_9$	=	$\left(\frac{3}{4} + y_3 - z_3\right)\mathbf{a}_1 + \left(\frac{1}{4} - x_3 - z_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_3 + y_3\right)\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_3\right)c\,\hat{\mathbf{z}}$	(16*e*)	H I
$\mathbf{B}_{10}$	=	$\left(\frac{3}{4} - y_3 - z_3\right)\mathbf{a}_1 + \left(\frac{1}{4} + x_3 - z_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_3 - y_3\right)\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_3\right)c\,\hat{\mathbf{z}}$	(16*e*)	H I
$\mathbf{B}_{11}$	=	$\left(\frac{3}{4} - x_3 + z_3\right)\mathbf{a}_1 + \left(\frac{1}{4} - y_3 + z_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_3 - y_3\right)\mathbf{a}_3$	=	$-y_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_3\right)c\,\hat{\mathbf{z}}$	(16*e*)	H I
$\mathbf{B}_{12}$	=	$\left(\frac{3}{4} + x_3 + z_3\right)\mathbf{a}_1 + \left(\frac{1}{4} + y_3 + z_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_3 + y_3\right)\mathbf{a}_3$	=	$y_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_3\right)c\,\hat{\mathbf{z}}$	(16*e*)	H I
$\mathbf{B}_{13}$	=	$(y_4 + z_4)\,\mathbf{a}_1 + (x_4 + z_4)\,\mathbf{a}_2 + (x_4 + y_4)\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + y_4a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	(16*e*)	H II
$\mathbf{B}_{14}$	=	$(-y_4 + z_4)\,\mathbf{a}_1 + (-x_4 + z_4)\,\mathbf{a}_2 + (-x_4 - y_4)\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} - y_4a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	(16*e*)	H II
$\mathbf{B}_{15}$	=	$(-x_4 - z_4)\,\mathbf{a}_1 + (y_4 - z_4)\,\mathbf{a}_2 + (-x_4 + y_4)\,\mathbf{a}_3$	=	$y_4a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	(16*e*)	H II
$\mathbf{B}_{16}$	=	$(x_4 - z_4)\,\mathbf{a}_1 + (-y_4 - z_4)\,\mathbf{a}_2 + (x_4 - y_4)\,\mathbf{a}_3$	=	$-y_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	(16*e*)	H II
$\mathbf{B}_{17}$	=	$\left(\frac{3}{4} + y_4 - z_4\right)\mathbf{a}_1 + \left(\frac{1}{4} - x_4 - z_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_4 + y_4\right)\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_4\right)c\,\hat{\mathbf{z}}$	(16*e*)	H II
$\mathbf{B}_{18}$	=	$\left(\frac{3}{4} - y_4 - z_4\right)\mathbf{a}_1 + \left(\frac{1}{4} + x_4 - z_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_4 - y_4\right)\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_4\right)c\,\hat{\mathbf{z}}$	(16*e*)	H II
$\mathbf{B}_{19}$	=	$\left(\frac{3}{4} - x_4 + z_4\right)\mathbf{a}_1 + \left(\frac{1}{4} - y_4 + z_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_4 - y_4\right)\mathbf{a}_3$	=	$-y_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_4\right)c\,\hat{\mathbf{z}}$	(16*e*)	H II
$\mathbf{B}_{20}$	=	$\left(\frac{3}{4} + x_4 + z_4\right)\mathbf{a}_1 + \left(\frac{1}{4} + y_4 + z_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_4 + y_4\right)\mathbf{a}_3$	=	$y_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_4\right)c\,\hat{\mathbf{z}}$	(16*e*)	H II
$\mathbf{B}_{21}$	=	$(y_5 + z_5)\,\mathbf{a}_1 + (x_5 + z_5)\,\mathbf{a}_2 + (x_5 + y_5)\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}} + y_5a\,\hat{\mathbf{y}} + z_5c\,\hat{\mathbf{z}}$	(16*e*)	O
$\mathbf{B}_{22}$	=	$(-y_5 + z_5)\,\mathbf{a}_1 + (-x_5 + z_5)\,\mathbf{a}_2 + (-x_5 - y_5)\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}} - y_5a\,\hat{\mathbf{y}} + z_5c\,\hat{\mathbf{z}}$	(16*e*)	O

$\mathbf{B}_{23}$	=	$(-x_5 - z_5)\ \mathbf{a}_1 + (y_5 - z_5)\ \mathbf{a}_2 + (-x_5 + y_5)\ \mathbf{a}_3$	=	$y_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(16*e*)	O
$\mathbf{B}_{24}$	=	$(x_5 - z_5)\ \mathbf{a}_1 + (-y_5 - z_5)\ \mathbf{a}_2 + (x_5 - y_5)\ \mathbf{a}_3$	=	$-y_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(16*e*)	O
$\mathbf{B}_{25}$	=	$\left(\frac{3}{4} + y_5 - z_5\right)\mathbf{a}_1 + \left(\frac{1}{4} - x_5 - z_5\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_5 + y_5\right)\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_5\right) c\,\hat{\mathbf{z}}$	(16*e*)	O
$\mathbf{B}_{26}$	=	$\left(\frac{3}{4} - y_5 - z_5\right)\mathbf{a}_1 + \left(\frac{1}{4} + x_5 - z_5\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_5 - y_5\right)\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_5\right) c\,\hat{\mathbf{z}}$	(16*e*)	O
$\mathbf{B}_{27}$	=	$\left(\frac{3}{4} - x_5 + z_5\right)\mathbf{a}_1 + \left(\frac{1}{4} - y_5 + z_5\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_5 - y_5\right)\mathbf{a}_3$	=	$-y_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_5\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_5\right) c\,\hat{\mathbf{z}}$	(16*e*)	O
$\mathbf{B}_{28}$	=	$\left(\frac{3}{4} + x_5 + z_5\right)\mathbf{a}_1 + \left(\frac{1}{4} + y_5 + z_5\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_5 + y_5\right)\mathbf{a}_3$	=	$y_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_5\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_5\right) c\,\hat{\mathbf{z}}$	(16*e*)	O

References:

- A. A. Khan and W. H. Baur, *Refinement of the crystal structures of ammonium dihydrogen phosphate and ammonium dihydrogen arsenate*, Acta Crystallogr. Sect. B Struct. Sci. **29**, 2721–2726 (1973), doi:10.1107/S0567740873007442.

Found in:

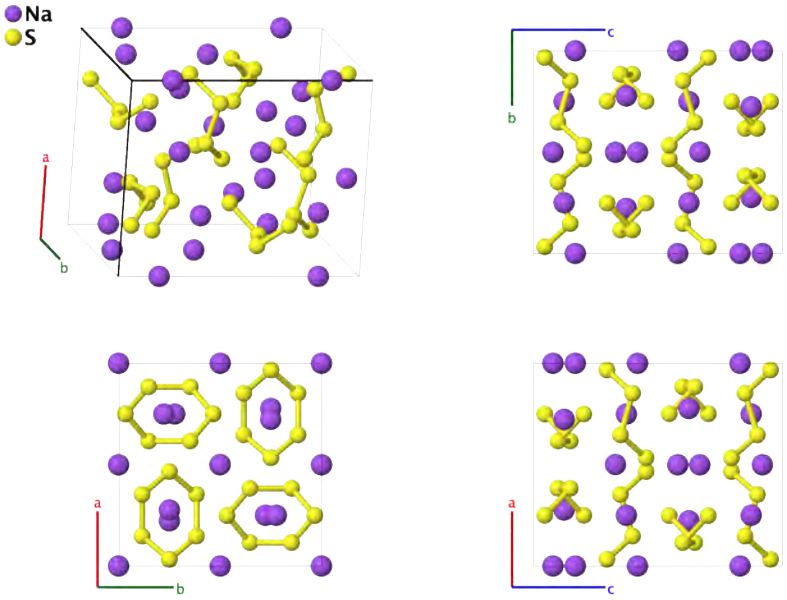
- T. Fukami, S. Akahoshi, K. Hukuda, and T. Yagi, *Refinement of the Crystal Structure of $NH_4H_2PO_4$ above and below Antiferroelectric Phase Transition Temperature*, J. Phys. Soc. Jpn. **56**, 2223–2224 (1987), doi:10.1143/JPSJ.56.2223.

Geometry files:

- CIF: pp. 1688
- POSCAR: pp. 1688

NaS_2 Structure: AB2_tI48_122_cd_2e

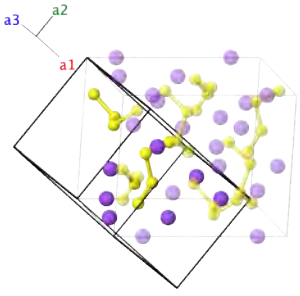
http://aflow.org/prototype-encyclopedia/AB2_tI48_122_cd_2e



Prototype	:	NaS_2
AFLOW prototype label	:	AB2_tI48_122_cd_2e
***Strukturbericht* designation**	:	None
Pearson symbol	:	tI48
Space group number	:	122
Space group symbol	:	$I\bar{4}2d$
AFLOW prototype command	:	`aflow --proto=AB2_tI48_122_cd_2e --params=`$a, c/a, z_1, x_2, x_3, y_3, z_3, x_4, y_4, z_4$

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}a\,\hat{\mathbf{x}} - \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} - \tfrac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$z_1\,\mathbf{a}_1 + z_1\,\mathbf{a}_2$	=	$z_1 c\,\hat{\mathbf{z}}$	($8c$)	Na I
$\mathbf{B}_2$	=	$-z_1\,\mathbf{a}_1 - z_1\,\mathbf{a}_2$	=	$-z_1 c\,\hat{\mathbf{z}}$	($8c$)	Na I
$\mathbf{B}_3$	=	$\left(\frac{3}{4} - z_1\right)\mathbf{a}_1 + \left(\frac{1}{4} - z_1\right)\mathbf{a}_2 + \frac{1}{2}\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_1\right)c\,\hat{\mathbf{z}}$	($8c$)	Na I
$\mathbf{B}_4$	=	$\left(\frac{3}{4} + z_1\right)\mathbf{a}_1 + \left(\frac{1}{4} + z_1\right)\mathbf{a}_2 + \frac{1}{2}\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_1\right)c\,\hat{\mathbf{z}}$	($8c$)	Na I

$$
\begin{array}{lcccccl}
\mathbf{B}_5 & = & \frac{3}{8}\,\mathbf{a}_1 + \left(\frac{1}{8} + x_2\right)\mathbf{a}_2 + \left(\frac{1}{4} + x_2\right)\mathbf{a}_3 & = & x_2 a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{8}c\,\hat{\mathbf{z}} & (8d) & \text{Na II} \\
\mathbf{B}_6 & = & \frac{7}{8}\,\mathbf{a}_1 + \left(\frac{1}{8} - x_2\right)\mathbf{a}_2 + \left(\frac{3}{4} - x_2\right)\mathbf{a}_3 & = & -x_2 a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{1}{8}c\,\hat{\mathbf{z}} & (8d) & \text{Na II} \\
\mathbf{B}_7 & = & \left(\frac{7}{8} - x_2\right)\mathbf{a}_1 + \frac{1}{8}\,\mathbf{a}_2 + \left(\frac{1}{4} - x_2\right)\mathbf{a}_3 & = & -\frac{1}{4}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{y}} + \frac{3}{8}c\,\hat{\mathbf{z}} & (8d) & \text{Na II} \\
\mathbf{B}_8 & = & \left(\frac{7}{8} + x_2\right)\mathbf{a}_1 + \frac{5}{8}\,\mathbf{a}_2 + \left(\frac{3}{4} + x_2\right)\mathbf{a}_3 & = & \frac{1}{4}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{y}} + \frac{3}{8}c\,\hat{\mathbf{z}} & (8d) & \text{Na II} \\
\mathbf{B}_9 & = & (y_3 + z_3)\,\mathbf{a}_1 + (x_3 + z_3)\,\mathbf{a}_2 + (x_3 + y_3)\,\mathbf{a}_3 & = & x_3 a\,\hat{\mathbf{x}} + y_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}} & (16e) & \text{S I} \\
\mathbf{B}_{10} & = & (-y_3 + z_3)\,\mathbf{a}_1 + (-x_3 + z_3)\,\mathbf{a}_2 + (-x_3 - y_3)\,\mathbf{a}_3 & = & -x_3 a\,\hat{\mathbf{x}} - y_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}} & (16e) & \text{S I} \\
\mathbf{B}_{11} & = & (-x_3 - z_3)\,\mathbf{a}_1 + (y_3 - z_3)\,\mathbf{a}_2 + (-x_3 + y_3)\,\mathbf{a}_3 & = & y_3 a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}} & (16e) & \text{S I} \\
\mathbf{B}_{12} & = & (x_3 - z_3)\,\mathbf{a}_1 + (-y_3 - z_3)\,\mathbf{a}_2 + (x_3 - y_3)\,\mathbf{a}_3 & = & -y_3 a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}} & (16e) & \text{S I} \\
\mathbf{B}_{13} & = & \left(\frac{3}{4} + y_3 - z_3\right)\mathbf{a}_1 + \left(\frac{1}{4} - x_3 - z_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_3 + y_3\right)\mathbf{a}_3 & = & -x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_3\right)c\,\hat{\mathbf{z}} & (16e) & \text{S I} \\
\mathbf{B}_{14} & = & \left(\frac{3}{4} - y_3 - z_3\right)\mathbf{a}_1 + \left(\frac{1}{4} + x_3 - z_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_3 - y_3\right)\mathbf{a}_3 & = & x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_3\right)c\,\hat{\mathbf{z}} & (16e) & \text{S I} \\
\mathbf{B}_{15} & = & \left(\frac{3}{4} - x_3 + z_3\right)\mathbf{a}_1 + \left(\frac{1}{4} - y_3 + z_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_3 - y_3\right)\mathbf{a}_3 & = & -y_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_3\right)c\,\hat{\mathbf{z}} & (16e) & \text{S I} \\
\mathbf{B}_{16} & = & \left(\frac{3}{4} + x_3 + z_3\right)\mathbf{a}_1 + \left(\frac{1}{4} + y_3 + z_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_3 + y_3\right)\mathbf{a}_3 & = & y_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_3\right)c\,\hat{\mathbf{z}} & (16e) & \text{S I} \\
\mathbf{B}_{17} & = & (y_4 + z_4)\,\mathbf{a}_1 + (x_4 + z_4)\,\mathbf{a}_2 + (x_4 + y_4)\,\mathbf{a}_3 & = & x_4 a\,\hat{\mathbf{x}} + y_4 a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}} & (16e) & \text{S II} \\
\mathbf{B}_{18} & = & (-y_4 + z_4)\,\mathbf{a}_1 + (-x_4 + z_4)\,\mathbf{a}_2 + (-x_4 - y_4)\,\mathbf{a}_3 & = & -x_4 a\,\hat{\mathbf{x}} - y_4 a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}} & (16e) & \text{S II} \\
\mathbf{B}_{19} & = & (-x_4 - z_4)\,\mathbf{a}_1 + (y_4 - z_4)\,\mathbf{a}_2 + (-x_4 + y_4)\,\mathbf{a}_3 & = & y_4 a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}} & (16e) & \text{S II} \\
\mathbf{B}_{20} & = & (x_4 - z_4)\,\mathbf{a}_1 + (-y_4 - z_4)\,\mathbf{a}_2 + (x_4 - y_4)\,\mathbf{a}_3 & = & -y_4 a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}} & (16e) & \text{S II} \\
\mathbf{B}_{21} & = & \left(\frac{3}{4} + y_4 - z_4\right)\mathbf{a}_1 + \left(\frac{1}{4} - x_4 - z_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_4 + y_4\right)\mathbf{a}_3 & = & -x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_4\right)c\,\hat{\mathbf{z}} & (16e) & \text{S II} \\
\mathbf{B}_{22} & = & \left(\frac{3}{4} - y_4 - z_4\right)\mathbf{a}_1 + \left(\frac{1}{4} + x_4 - z_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_4 - y_4\right)\mathbf{a}_3 & = & x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_4\right)c\,\hat{\mathbf{z}} & (16e) & \text{S II} \\
\mathbf{B}_{23} & = & \left(\frac{3}{4} - x_4 + z_4\right)\mathbf{a}_1 + \left(\frac{1}{4} - y_4 + z_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_4 - y_4\right)\mathbf{a}_3 & = & -y_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_4\right)c\,\hat{\mathbf{z}} & (16e) & \text{S II} \\
\mathbf{B}_{24} & = & \left(\frac{3}{4} + x_4 + z_4\right)\mathbf{a}_1 + \left(\frac{1}{4} + y_4 + z_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_4 + y_4\right)\mathbf{a}_3 & = & y_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_4\right)c\,\hat{\mathbf{z}} & (16e) & \text{S II}
\end{array}
$$

References:

- R. Tegman, *The Crystal Structure of Sodium Tetrasulphide, Na_2S_4*, Acta Crystallogr. Sect. B Struct. Sci. **29**, 1463–1469 (1973), doi:10.1107/S0567740873004735.

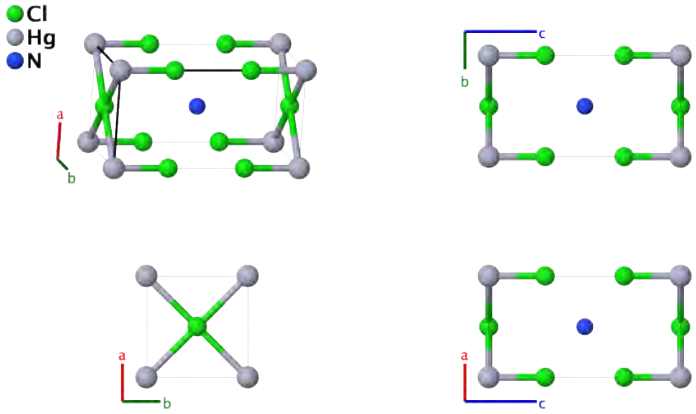
Found in:

- P. Villars and L. Calvert, *Pearson's Handbook of Crystallographic Data for Intermetallic Phases* (ASM International, Materials Park, OH, 1991), 2nd edn.

Geometry files:
- CIF: pp. 1688
- POSCAR: pp. 1689

NH_4HgCl_3 ($E2_5$) Structure: A3BC_tP5_123_cg_a_d

http://aflow.org/prototype-encyclopedia/A3BC_tP5_123_cg_a_d

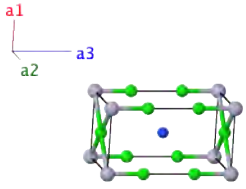


Prototype	:	$Cl_3Hg(NH_4)$
AFLOW prototype label	:	A3BC_tP5_123_cg_a_d
***Strukturbericht* designation**	:	$E2_5$
Pearson symbol	:	tP5
Space group number	:	123
Space group symbol	:	$P4/mmm$
AFLOW prototype command	:	`aflow --proto=A3BC_tP5_123_cg_a_d --params=`$a, c/a, z_4$

- The positions of the hydrogen atoms in the NH_4 ion are not given, so we only provide the nitrogen positions (labeled as NH_4). It is likely that the hydrogen atoms are freely rotating around the nitrogen, as any reasonable fixed positions would destroy both the inversion symmetry and the four-fold rotation axis exhibited by space group $P4/mmm$.
- (Harmsen, 1939) and (Herrmann, 1943) give multiple possible space groups for this structure. We have chosen the highest symmetry representation, space group $P4/mmm$ #123.

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($1a$)	Hg
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$	($1c$)	Cl I
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($1d$)	NH_4
$\mathbf{B}_4$	=	$z_4\,\mathbf{a}_3$	=	$z_4 c\,\hat{\mathbf{z}}$	($2g$)	Cl II

$\mathbf{B}_5$	=	$-z_4\,\mathbf{a}_3$	=	$-z_4 c\,\hat{\mathbf{z}}$	$(2g)$	Cl II

References:

- E. J. Harmsen, *The Crystal Structure of NH_4HgCl_3*, Zeitschrift für Kristallographie - Crystalline Materials **100**, 208–211 (1939), doi:10.1524/zkri.1939.100.1.208.

Found in:

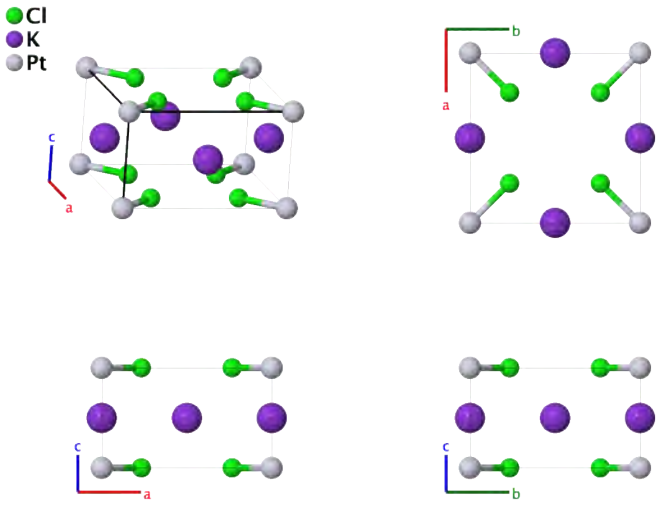
- K. Herrmann, ed., *Strukturbericht Band VII 1939* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1943).

Geometry files:

- CIF: pp. 1689
- POSCAR: pp. 1689

K_2PtCl_4 ($H1_5$) Structure: A4B2C_tP7_123_j_e_a

http://aflow.org/prototype-encyclopedia/A4B2C_tP7_123_j_e_a



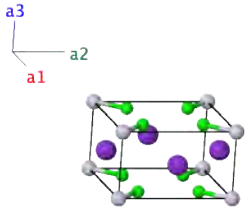
Prototype	:	Cl_4K_2Pt
AFLOW prototype label	:	A4B2C_tP7_123_j_e_a
***Strukturbericht* designation**	:	$H1_5$
Pearson symbol	:	tP7
Space group number	:	123
Space group symbol	:	$P4/mmm$
AFLOW prototype command	:	`aflow --proto=A4B2C_tP7_123_j_e_a --params=`$a, c/a, x_3$

Other compounds with this structure

- K_2PdCl_4

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(1a)$	Pt
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(2e)$	K

$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(2e)$	K
$\mathbf{B}_4$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2$	=	$x_3 a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}}$	$(4j)$	Cl
$\mathbf{B}_5$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2$	=	$-x_3 a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}}$	$(4j)$	Cl
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2$	=	$-x_3 a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}}$	$(4j)$	Cl
$\mathbf{B}_7$	=	$x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2$	=	$x_3 a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}}$	$(4j)$	Cl

References:

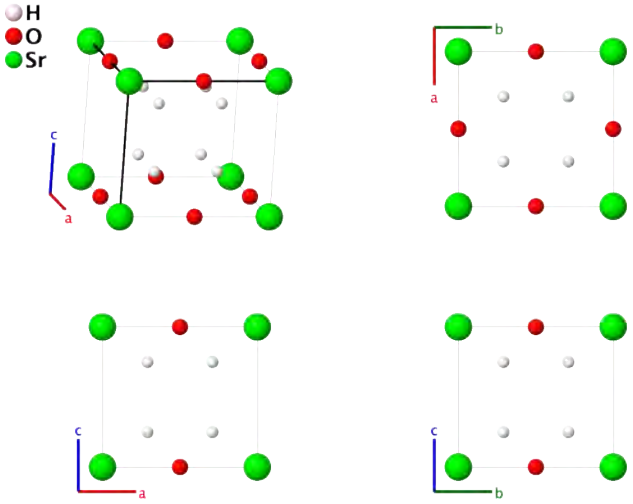
- R. H. B. Mais, P. G. Owston, and A. Wood, *The crystal structure of K_2PtCl_4 and K_2PdCl_4 with estimates of the factors affecting accuracy*, Acta Crystallogr. Sect. B Struct. Sci. **28**, 393–399 (1972), doi:10.1107/S0567740872002468.

Geometry files:

- CIF: pp. 1689
- POSCAR: pp. 1690

$E6_1$ ($Sr(OH)_2(H_2O)_8$) (*obsolete*) Structure: A8B2C_tP11_123_r_f_a

http://aflow.org/prototype-encyclopedia/A8B2C_tP11_123_r_f_a

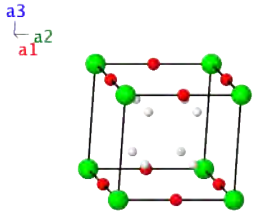


Prototype	:	$Sr(OH)_2(H_2O)_8$
AFLOW prototype label	:	A8B2C_tP11_123_r_f_a
***Strukturbericht* designation**	:	$E6_1$
Pearson symbol	:	tP11
Space group number	:	123
Space group symbol	:	$P4/mmm$
AFLOW prototype command	:	`aflow --proto=A8B2C_tP11_123_r_f_a --params=`$a, c/a, x_3, z_3$

- This structure was designated $E6_1$ by (Hermann, 1937). It has been superseded by the structure determined by (Ricci, 2005). We present it for historical interest.

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(1a)$	Sr
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{y}}$	$(2f)$	OH

$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	$(2f)$	OH
$\mathbf{B}_4$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	$(8r)$	H_2O
$\mathbf{B}_5$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	$(8r)$	H_2O
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	$(8r)$	H_2O
$\mathbf{B}_7$	=	$x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	$(8r)$	H_2O
$\mathbf{B}_8$	=	$-x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	$(8r)$	H_2O
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	$(8r)$	H_2O
$\mathbf{B}_{10}$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	$(8r)$	H_2O
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	$(8r)$	H_2O

References:

- G. Natta, *Constitution of hydroxides and of hydrates. III. Octahydrated strontium hydroxide*, Gazz. Chim. Ital. **58**, 870–872 (1928).

Found in:

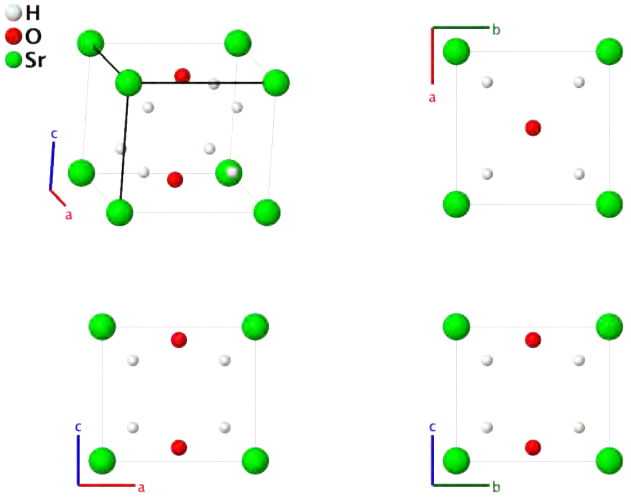
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).
- J. S. Ricci, R. C. Stevens, R. K. McMullan, and W. T. Klooster, *Structure of strontium hydroxide octahydrate, $Sr(OH)_2{\cdot}8H_2O$, at 20, 100 and 200 K from neutron diffraction*, Acta Crystallogr. Sect. B Struct. Sci. **61**, 381–386 (2005), doi:10.1107/S0108768105013480.

Geometry files:

- CIF: pp. 1690
- POSCAR: pp. 1690

$E6_2$ [$SrO_2(H_2O)_8$] (*possibly obsolete*) Structure: A8B2C_tP11_123_r_h_a

http://aflow.org/prototype-encyclopedia/A8B2C_tP11_123_r_h_a



Prototype	:	$(H_2O)_8SrO_2$
AFLOW prototype label	:	A8B2C_tP11_123_r_h_a
***Strukturbericht* designation**	:	$E6_2$
Pearson symbol	:	tP11
Space group number	:	123
Space group symbol	:	$P4/mmm$
AFLOW prototype command	:	`aflow --proto=A8B2C_tP11_123_r_h_a --params=`$a, c/a, z_2, x_3, z_3$

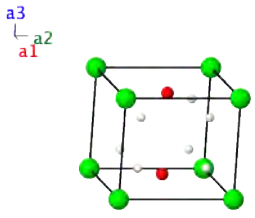
Other compounds with this structure

- $CaO_2(H_2O)_8$ and $BrO_2(H_2O)_8$

- This structure was designated $E6_2$ by (Hermann, 1937). (Shineman, 1951) suggest that this should be replaced by their $CaO_2(H_2O)_8$ structure.
- The positions of the hydrogen atoms in the water molecule were not determined, so we only provide the positions of the oyxgen atoms (labeled as H_2O).

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($1a$)	Sr
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	($2h$)	O
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	($2h$)	O
$\mathbf{B}_4$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	($8r$)	H_2O
$\mathbf{B}_5$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	($8r$)	H_2O
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	($8r$)	H_2O
$\mathbf{B}_7$	=	$x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	($8r$)	H_2O
$\mathbf{B}_8$	=	$-x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	($8r$)	H_2O
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	($8r$)	H_2O
$\mathbf{B}_{10}$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	($8r$)	H_2O
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	($8r$)	H_2O

References:

- G. Natta, *Struttura degli idrossidi ed idrati - Nota IV. Sul perossido di stronzio ottoidrato*, Gazz. Chim. Ital. **62**, 444 (1932).

Found in:

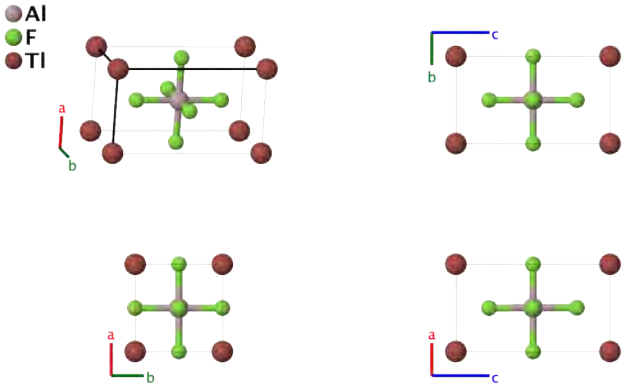
- R. S. Shineman and A. J. King, *The space group of calcium peroxide octahydrate*, Acta Cryst. **4**, 67–68 (1951), doi:10.1107/S0365110X5100012X.

Geometry files:

- CIF: pp. 1690
- POSCAR: pp. 1691

$TlAlF_4$ ($H0_8$) Structure: AB4C_tP6_123_d_eh_a

http://aflow.org/prototype-encyclopedia/AB4C_tP6_123_d_eh_a



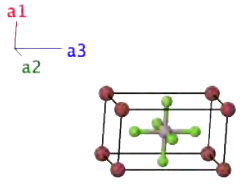
Prototype	:	AlF_4Tl
AFLOW prototype label	:	AB4C_tP6_123_d_eh_a
***Strukturbericht* designation**	:	$H0_8$
Pearson symbol	:	tP6
Space group number	:	123
Space group symbol	:	$P4/mmm$
AFLOW prototype command	:	`aflow --proto=AB4C_tP6_123_d_eh_a --params=`$a, c/a, z_4$

Other compounds with this structure

- β-$RbFeF_4$, $CsFeF_4$, $KAlF_4$, and $RbAlF_4$

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($1a$)	Tl
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($1d$)	Al
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($2e$)	F I
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($2e$)	F I
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	($2h$)	F II
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	($2h$)	F II

References:

- C. Brosset, *Herstellung und Kristallbau der Verbindungen $TlAlF_4$ und Tl_2AlF_5*, Z. Anorg. Allg. Chem. **235**, 139–147 (1937), doi:10.1002/zaac.19372350119.

Found in:

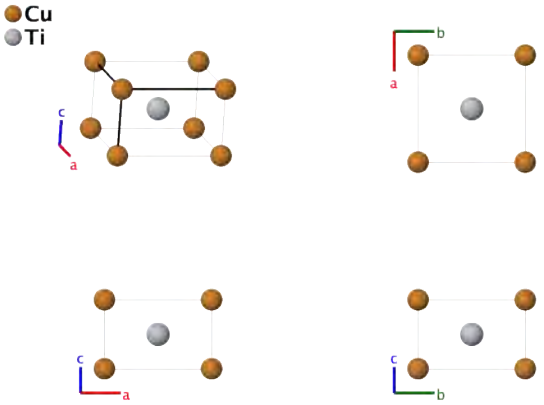
- A. Pabst, *A Structural Classification of Fluoaluminates*, Am. Mineral. **35**, 149–165 (1950).

Geometry files:

- CIF: pp. 1691
- POSCAR: pp. 1691

δ-CuTi ($L2_a$) Structure: AB_tP2_123_a_d

http://aflow.org/prototype-encyclopedia/AB_tP2_123_a_d.CuTi



Prototype	:	CuTi
AFLOW prototype label	:	AB_tP2_123_a_d
***Strukturbericht* designation**	:	$L2_a$
Pearson symbol	:	tP2
Space group number	:	123
Space group symbol	:	$P4/mmm$
AFLOW prototype command	:	`aflow --proto=AB_tP2_123_a_d --params=`$a, c/a$

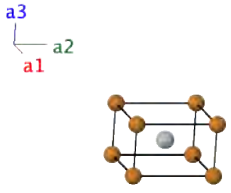
Other compounds with this structure

- HgMn

- This structure has the same AFLOW designation, AB_tP2_123_a_d, as CuAu ($L1_0$). The only difference in the structures is the c/a ratio. $L1_0$ has $c/a \approx \sqrt{2}$, characteristitic of face-centered cubic ordering, while $L2_a$ has $c/a \approx 1$, a body-centered cubic-like system.

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(1a)$	Cu
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(1d)$	Ti

References:

- N. Karlsson, *An X-Ray Study of the Phases in the Copper-Titanium System*, J. Inst. Met. **79**, 391–405 (1951).

Found in:

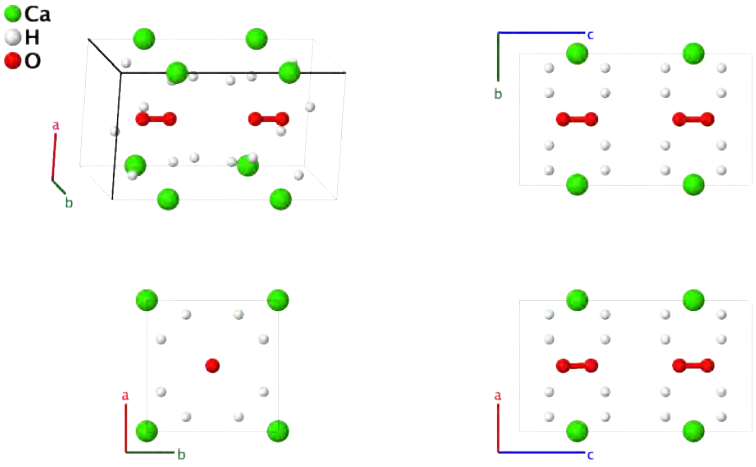
- W. B. Pearson, *A Handbook of Lattice Spacings and Structures of Metals and Alloys*, *International Series of Monographs on Metal Physics and Physical Metallurgy*, vol. 4 (Pergamon Press, Oxford, London, Edinburgh, New York, Paris, Frankfort, 1958), 1964 reprint with corrections edn. N. R. C. No. 4303.

Geometry files:

- CIF: pp. 1691
- POSCAR: pp. 1692

$CaO_2(H_2O)_8$ Structure: AB8C2_tP22_124_a_n_h

http://aflow.org/prototype-encyclopedia/AB8C2_tP22_124_a_n_h



Prototype	:	$Ca(H_2O)_8O_2$
AFLOW prototype label	:	AB8C2_tP22_124_a_n_h
***Strukturbericht* designation**	:	None
Pearson symbol	:	tP22
Space group number	:	124
Space group symbol	:	$P4/mcc$
AFLOW prototype command	:	`aflow --proto=AB8C2_tP22_124_a_n_h --params=`$a, c/a, z_2, x_3, y_3, z_3$

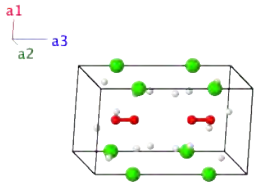
Other compounds with this structure

- $BrO_2(H_2O)_8$ and $SaO_2(H_2O)_8$

- This structure was proposed by (Shineman, 1951) to replace the $E6_2$ $SrO_2(H_2O)_8$ structure. The unit cell is doubled in the (001) direction, and the oxygens are now molecular.
- The positions of the hydrogen atom in the water molecule were not determined, so we only provide the oxygen position (labeled as H_2O).

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\,\hat{\mathbf{z}}$	$(2a)$	Ca

$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\,\hat{\mathbf{z}}$	$(2a)$	Ca
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(4h)$	O
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	$(4h)$	O
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	$(4h)$	O
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	$(4h)$	O
$\mathbf{B}_7$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(16n)$	H_2O
$\mathbf{B}_8$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(16n)$	H_2O
$\mathbf{B}_9$	=	$-y_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$-y_3 a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(16n)$	H_2O
$\mathbf{B}_{10}$	=	$y_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$y_3 a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(16n)$	H_2O
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + y_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	$(16n)$	H_2O
$\mathbf{B}_{12}$	=	$x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} - y_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	$(16n)$	H_2O
$\mathbf{B}_{13}$	=	$y_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$y_3 a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	$(16n)$	H_2O
$\mathbf{B}_{14}$	=	$-y_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$-y_3 a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	$(16n)$	H_2O
$\mathbf{B}_{15}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(16n)$	H_2O
$\mathbf{B}_{16}$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(16n)$	H_2O
$\mathbf{B}_{17}$	=	$y_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$y_3 a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(16n)$	H_2O
$\mathbf{B}_{18}$	=	$-y_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-y_3 a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(16n)$	H_2O
$\mathbf{B}_{19}$	=	$x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} - y_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	$(16n)$	H_2O
$\mathbf{B}_{20}$	=	$-x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + y_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	$(16n)$	H_2O
$\mathbf{B}_{21}$	=	$-y_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$-y_3 a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	$(16n)$	H_2O
$\mathbf{B}_{22}$	=	$y_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$y_3 a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	$(16n)$	H_2O

References:

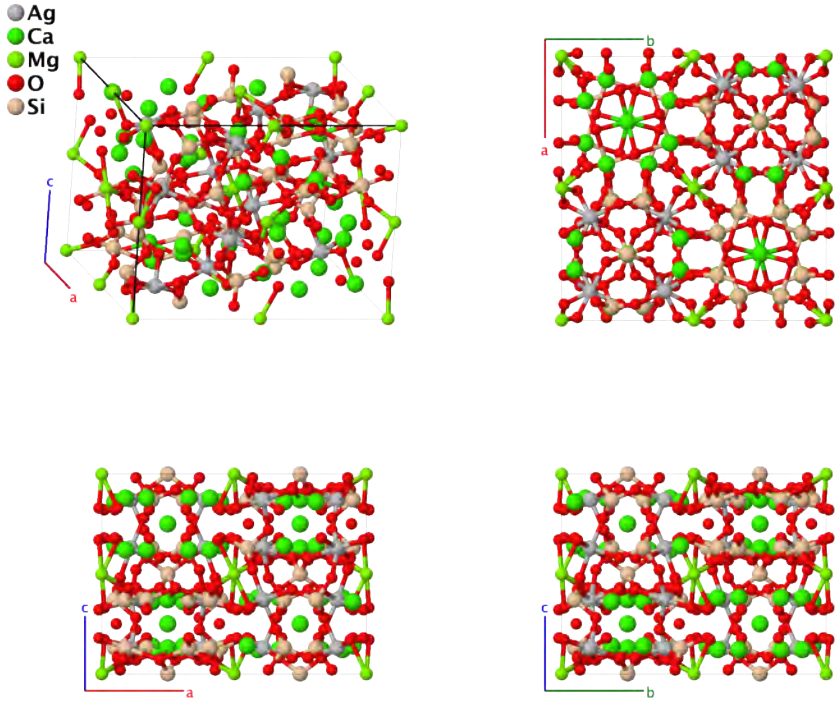
- R. S. Shineman and A. J. King, *The space group of calcium peroxide octahydrate*, Acta Cryst. **4**, 67–68 (1951), doi:10.1107/S0365110X5100012X.

Geometry files:

- CIF: pp. 1692
- POSCAR: pp. 1692

Vesuvianite ($Ca_{10}Al_4(Mg,Fe)_2Si_9O_{34}(OH)_4$, $S2_3$) Structure: A4B10C2D34E4F9_tP252_126_k_ce2k_f_h8k_k_d2k

http://aflow.org/prototype-encyclopedia/A4B10C2D34E4F9_tP252_126_k_ce2k_f_h8k_k_d2k

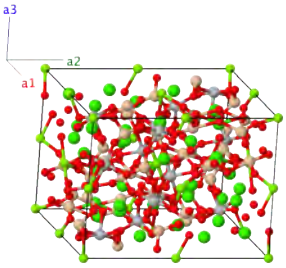


Prototype	:	$Al_4Ca_{10}(Mg,Fe)_2O_{34}(OH)_4Si_9$
AFLOW prototype label	:	A4B10C2D34E4F9_tP252_126_k_ce2k_f_h8k_k_d2k
***Strukturbericht* designation**	:	$S2_3$
Pearson symbol	:	tP252
Space group number	:	126
Space group symbol	:	$P4/nnc$
AFLOW prototype command	:	`aflow --proto=A4B10C2D34E4F9_tP252_126_k_ce2k_f_h8k_k_d2k --params=`$a, c/a, z_3, x_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}, x_{16}, y_{16}, z_{16}, x_{17}, y_{17}, z_{17}, x_{18}, y_{18}, z_{18}, x_{19}, y_{19}, z_{19}$

- Vesuvianite, also known as idocrase, comes in a variety of compositions and structures, see *e.g.*, (Allen, 1992) and (Rucklidge, 1975) and references therein. For our example we use the original structure of (Warren, 1931), where the magnesium ($8f$) site is filled by a random (Mg,Fe) alloy. The positions of the hydrogen atoms in the OH ions were not determined, so we only give the positions of the oyxgen atoms (labeled as OH).

Simple Tetragonal primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= a\,\hat{\mathbf{x}}\\ \mathbf{a}_2 &= a\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(4c)$	Ca I
$\mathbf{B}_2$	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(4c)$	Ca I
$\mathbf{B}_3$	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(4c)$	Ca I
$\mathbf{B}_4$	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(4c)$	Ca I
$\mathbf{B}_5$	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2$	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}}$	$(4d)$	Si I
$\mathbf{B}_6$	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2$	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}}$	$(4d)$	Si I
$\mathbf{B}_7$	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(4d)$	Si I
$\mathbf{B}_8$	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(4d)$	Si I
$\mathbf{B}_9$	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(4e)$	Ca II
$\mathbf{B}_{10}$	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	$(4e)$	Ca II
$\mathbf{B}_{11}$	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(4e)$	Ca II
$\mathbf{B}_{12}$	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	$(4e)$	Ca II
$\mathbf{B}_{13}$	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(8f)$	Mg
$\mathbf{B}_{14}$	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$	$(8f)$	Mg
$\mathbf{B}_{15}$	$\frac{1}{2}\,\mathbf{a}_1$	$\frac{1}{2}a\,\hat{\mathbf{x}}$	$(8f)$	Mg
$\mathbf{B}_{16}$	$\frac{1}{2}\,\mathbf{a}_2$	$\frac{1}{2}a\,\hat{\mathbf{y}}$	$(8f)$	Mg
$\mathbf{B}_{17}$	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(8f)$	Mg
$\mathbf{B}_{18}$	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(8f)$	Mg
$\mathbf{B}_{19}$	$\frac{1}{2}\,\mathbf{a}_3$	$\frac{1}{2}c\,\hat{\mathbf{z}}$	$(8f)$	Mg
$\mathbf{B}_{20}$	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(8f)$	Mg
$\mathbf{B}_{21}$	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	$x_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(8h)$	O I
$\mathbf{B}_{22}$	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_5\right)\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(8h)$	O I
$\mathbf{B}_{23}$	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(8h)$	O I
$\mathbf{B}_{24}$	$x_5\,\mathbf{a}_1 + \left(\frac{1}{2} - x_5\right)\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	$x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(8h)$	O I
$\mathbf{B}_{25}$	$-x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	$-x_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(8h)$	O I
$\mathbf{B}_{26}$	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_5\right)\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(8h)$	O I
$\mathbf{B}_{27}$	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(8h)$	O I

$\mathbf{B}_{28}$	$= -x_5\,\mathbf{a}_1 + \left(\frac{1}{2} + x_5\right)\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	$= -x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_5\right) a\,\hat{\mathbf{y}} + \frac{3}{4} c\,\hat{\mathbf{z}}$	$(8h)$	O I
$\mathbf{B}_{29}$	$= x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	$= x_6 a\,\hat{\mathbf{x}} + y_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(16k)$	Ag
$\mathbf{B}_{30}$	$= \left(\frac{1}{2} - x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + z_6\,\mathbf{a}_3$	$= \left(\frac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right) a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(16k)$	Ag
$\mathbf{B}_{31}$	$= \left(\frac{1}{2} - y_6\right)\mathbf{a}_1 + x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	$= \left(\frac{1}{2} - y_6\right) a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(16k)$	Ag
$\mathbf{B}_{32}$	$= y_6\,\mathbf{a}_1 + \left(\frac{1}{2} - x_6\right)\mathbf{a}_2 + z_6\,\mathbf{a}_3$	$= y_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_6\right) a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(16k)$	Ag
$\mathbf{B}_{33}$	$= \left(\frac{1}{2} - x_6\right)\mathbf{a}_1 + y_6\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	$= \left(\frac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} + y_6 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right) c\,\hat{\mathbf{z}}$	$(16k)$	Ag
$\mathbf{B}_{34}$	$= x_6\,\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	$= x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right) c\,\hat{\mathbf{z}}$	$(16k)$	Ag
$\mathbf{B}_{35}$	$= y_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	$= y_6 a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right) c\,\hat{\mathbf{z}}$	$(16k)$	Ag
$\mathbf{B}_{36}$	$= \left(\frac{1}{2} - y_6\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_6\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	$= \left(\frac{1}{2} - y_6\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_6\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right) c\,\hat{\mathbf{z}}$	$(16k)$	Ag
$\mathbf{B}_{37}$	$= -x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	$= -x_6 a\,\hat{\mathbf{x}} - y_6 a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	$(16k)$	Ag
$\mathbf{B}_{38}$	$= \left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3$	$= \left(\frac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right) a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	$(16k)$	Ag
$\mathbf{B}_{39}$	$= \left(\frac{1}{2} + y_6\right)\mathbf{a}_1 - x_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	$= \left(\frac{1}{2} + y_6\right) a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	$(16k)$	Ag
$\mathbf{B}_{40}$	$= -y_6\,\mathbf{a}_1 + \left(\frac{1}{2} + x_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3$	$= -y_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_6\right) a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	$(16k)$	Ag
$\mathbf{B}_{41}$	$= \left(\frac{1}{2} + x_6\right)\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	$= \left(\frac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} - y_6 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	$(16k)$	Ag
$\mathbf{B}_{42}$	$= -x_6\,\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	$= -x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	$(16k)$	Ag
$\mathbf{B}_{43}$	$= -y_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	$= -y_6 a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	$(16k)$	Ag
$\mathbf{B}_{44}$	$= \left(\frac{1}{2} + y_6\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_6\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	$= \left(\frac{1}{2} + y_6\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_6\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right) c\,\hat{\mathbf{z}}$	$(16k)$	Ag
$\mathbf{B}_{45}$	$= x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	$= x_7 a\,\hat{\mathbf{x}} + y_7 a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	$(16k)$	Ca III
$\mathbf{B}_{46}$	$= \left(\frac{1}{2} - x_7\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_7\right)\mathbf{a}_2 + z_7\,\mathbf{a}_3$	$= \left(\frac{1}{2} - x_7\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_7\right) a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	$(16k)$	Ca III
$\mathbf{B}_{47}$	$= \left(\frac{1}{2} - y_7\right)\mathbf{a}_1 + x_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	$= \left(\frac{1}{2} - y_7\right) a\,\hat{\mathbf{x}} + x_7 a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	$(16k)$	Ca III
$\mathbf{B}_{48}$	$= y_7\,\mathbf{a}_1 + \left(\frac{1}{2} - x_7\right)\mathbf{a}_2 + z_7\,\mathbf{a}_3$	$= y_7 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_7\right) a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	$(16k)$	Ca III
$\mathbf{B}_{49}$	$= \left(\frac{1}{2} - x_7\right)\mathbf{a}_1 + y_7\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\mathbf{a}_3$	$= \left(\frac{1}{2} - x_7\right) a\,\hat{\mathbf{x}} + y_7 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right) c\,\hat{\mathbf{z}}$	$(16k)$	Ca III
$\mathbf{B}_{50}$	$= x_7\,\mathbf{a}_1 + \left(\frac{1}{2} - y_7\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\mathbf{a}_3$	$= x_7 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_7\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right) c\,\hat{\mathbf{z}}$	$(16k)$	Ca III
$\mathbf{B}_{51}$	$= y_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\mathbf{a}_3$	$= y_7 a\,\hat{\mathbf{x}} + x_7 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right) c\,\hat{\mathbf{z}}$	$(16k)$	Ca III
$\mathbf{B}_{52}$	$= \left(\frac{1}{2} - y_7\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_7\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\mathbf{a}_3$	$= \left(\frac{1}{2} - y_7\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_7\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right) c\,\hat{\mathbf{z}}$	$(16k)$	Ca III
$\mathbf{B}_{53}$	$= -x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	$= -x_7 a\,\hat{\mathbf{x}} - y_7 a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	$(16k)$	Ca III
$\mathbf{B}_{54}$	$= \left(\frac{1}{2} + x_7\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\mathbf{a}_2 - z_7\,\mathbf{a}_3$	$= \left(\frac{1}{2} + x_7\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right) a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	$(16k)$	Ca III
$\mathbf{B}_{55}$	$= \left(\frac{1}{2} + y_7\right)\mathbf{a}_1 - x_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	$= \left(\frac{1}{2} + y_7\right) a\,\hat{\mathbf{x}} - x_7 a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	$(16k)$	Ca III
$\mathbf{B}_{56}$	$= -y_7\,\mathbf{a}_1 + \left(\frac{1}{2} + x_7\right)\mathbf{a}_2 - z_7\,\mathbf{a}_3$	$= -y_7 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_7\right) a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	$(16k)$	Ca III
$\mathbf{B}_{57}$	$= \left(\frac{1}{2} + x_7\right)\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	$= \left(\frac{1}{2} + x_7\right) a\,\hat{\mathbf{x}} - y_7 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right) c\,\hat{\mathbf{z}}$	$(16k)$	Ca III
$\mathbf{B}_{58}$	$= -x_7\,\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	$= -x_7 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right) c\,\hat{\mathbf{z}}$	$(16k)$	Ca III
$\mathbf{B}_{59}$	$= -y_7\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	$= -y_7 a\,\hat{\mathbf{x}} - x_7 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right) c\,\hat{\mathbf{z}}$	$(16k)$	Ca III
$\mathbf{B}_{60}$	$= \left(\frac{1}{2} + y_7\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_7\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	$= \left(\frac{1}{2} + y_7\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_7\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right) c\,\hat{\mathbf{z}}$	$(16k)$	Ca III

$\mathbf{B}_{61}$	=	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + y_8 a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(16k)	Ca IV
$\mathbf{B}_{62}$	=	$\left(\frac{1}{2} - x_8\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_8\right)\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_8\right)a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(16k)	Ca IV
$\mathbf{B}_{63}$	=	$\left(\frac{1}{2} - y_8\right)\mathbf{a}_1 + x_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_8\right)a\,\hat{\mathbf{x}} + x_8 a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(16k)	Ca IV
$\mathbf{B}_{64}$	=	$y_8\,\mathbf{a}_1 + \left(\frac{1}{2} - x_8\right)\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$y_8 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_8\right)a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(16k)	Ca IV
$\mathbf{B}_{65}$	=	$\left(\frac{1}{2} - x_8\right)\mathbf{a}_1 + y_8\,\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_8\right)a\,\hat{\mathbf{x}} + y_8 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right)c\,\hat{\mathbf{z}}$	(16k)	Ca IV
$\mathbf{B}_{66}$	=	$x_8\,\mathbf{a}_1 + \left(\frac{1}{2} - y_8\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_8\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right)c\,\hat{\mathbf{z}}$	(16k)	Ca IV
$\mathbf{B}_{67}$	=	$y_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\mathbf{a}_3$	=	$y_8 a\,\hat{\mathbf{x}} + x_8 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right)c\,\hat{\mathbf{z}}$	(16k)	Ca IV
$\mathbf{B}_{68}$	=	$\left(\frac{1}{2} - y_8\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_8\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_8\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right)c\,\hat{\mathbf{z}}$	(16k)	Ca IV
$\mathbf{B}_{69}$	=	$-x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} - y_8 a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(16k)	Ca IV
$\mathbf{B}_{70}$	=	$\left(\frac{1}{2} + x_8\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_8\right)\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_8\right)a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(16k)	Ca IV
$\mathbf{B}_{71}$	=	$\left(\frac{1}{2} + y_8\right)\mathbf{a}_1 - x_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_8\right)a\,\hat{\mathbf{x}} - x_8 a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(16k)	Ca IV
$\mathbf{B}_{72}$	=	$-y_8\,\mathbf{a}_1 + \left(\frac{1}{2} + x_8\right)\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$-y_8 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_8\right)a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(16k)	Ca IV
$\mathbf{B}_{73}$	=	$\left(\frac{1}{2} + x_8\right)\mathbf{a}_1 - y_8\,\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_8\right)a\,\hat{\mathbf{x}} - y_8 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right)c\,\hat{\mathbf{z}}$	(16k)	Ca IV
$\mathbf{B}_{74}$	=	$-x_8\,\mathbf{a}_1 + \left(\frac{1}{2} + y_8\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_8\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right)c\,\hat{\mathbf{z}}$	(16k)	Ca IV
$\mathbf{B}_{75}$	=	$-y_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$-y_8 a\,\hat{\mathbf{x}} - x_8 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right)c\,\hat{\mathbf{z}}$	(16k)	Ca IV
$\mathbf{B}_{76}$	=	$\left(\frac{1}{2} + y_8\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_8\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_8\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right)c\,\hat{\mathbf{z}}$	(16k)	Ca IV
$\mathbf{B}_{77}$	=	$x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}} + y_9 a\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(16k)	O II
$\mathbf{B}_{78}$	=	$\left(\frac{1}{2} - x_9\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_9\right)\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_9\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_9\right)a\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(16k)	O II
$\mathbf{B}_{79}$	=	$\left(\frac{1}{2} - y_9\right)\mathbf{a}_1 + x_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_9\right)a\,\hat{\mathbf{x}} + x_9 a\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(16k)	O II
$\mathbf{B}_{80}$	=	$y_9\,\mathbf{a}_1 + \left(\frac{1}{2} - x_9\right)\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$y_9 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_9\right)a\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(16k)	O II
$\mathbf{B}_{81}$	=	$\left(\frac{1}{2} - x_9\right)\mathbf{a}_1 + y_9\,\mathbf{a}_2 + \left(\frac{1}{2} - z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_9\right)a\,\hat{\mathbf{x}} + y_9 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_9\right)c\,\hat{\mathbf{z}}$	(16k)	O II
$\mathbf{B}_{82}$	=	$x_9\,\mathbf{a}_1 + \left(\frac{1}{2} - y_9\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_9\right)\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_9\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_9\right)c\,\hat{\mathbf{z}}$	(16k)	O II
$\mathbf{B}_{83}$	=	$y_9\,\mathbf{a}_1 + x_9\,\mathbf{a}_2 + \left(\frac{1}{2} - z_9\right)\mathbf{a}_3$	=	$y_9 a\,\hat{\mathbf{x}} + x_9 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_9\right)c\,\hat{\mathbf{z}}$	(16k)	O II
$\mathbf{B}_{84}$	=	$\left(\frac{1}{2} - y_9\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_9\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_9\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_9\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_9\right)c\,\hat{\mathbf{z}}$	(16k)	O II
$\mathbf{B}_{85}$	=	$-x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{x}} - y_9 a\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}}$	(16k)	O II
$\mathbf{B}_{86}$	=	$\left(\frac{1}{2} + x_9\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_9\right)\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_9\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_9\right)a\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}}$	(16k)	O II
$\mathbf{B}_{87}$	=	$\left(\frac{1}{2} + y_9\right)\mathbf{a}_1 - x_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_9\right)a\,\hat{\mathbf{x}} - x_9 a\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}}$	(16k)	O II
$\mathbf{B}_{88}$	=	$-y_9\,\mathbf{a}_1 + \left(\frac{1}{2} + x_9\right)\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$-y_9 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_9\right)a\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}}$	(16k)	O II
$\mathbf{B}_{89}$	=	$\left(\frac{1}{2} + x_9\right)\mathbf{a}_1 - y_9\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_9\right)a\,\hat{\mathbf{x}} - y_9 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right)c\,\hat{\mathbf{z}}$	(16k)	O II
$\mathbf{B}_{90}$	=	$-x_9\,\mathbf{a}_1 + \left(\frac{1}{2} + y_9\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_9\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right)c\,\hat{\mathbf{z}}$	(16k)	O II
$\mathbf{B}_{91}$	=	$-y_9\,\mathbf{a}_1 - x_9\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$-y_9 a\,\hat{\mathbf{x}} - x_9 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right)c\,\hat{\mathbf{z}}$	(16k)	O II
$\mathbf{B}_{92}$	=	$\left(\frac{1}{2} + y_9\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_9\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_9\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_9\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right)c\,\hat{\mathbf{z}}$	(16k)	O II
$\mathbf{B}_{93}$	=	$x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$x_{10} a\,\hat{\mathbf{x}} + y_{10} a\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}}$	(16k)	O III

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\begin{array}{lcccll}
\mathbf{B}_{94} & = & \left(\frac{1}{2}-x_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{10}\right)\mathbf{a}_2+z_{10}\,\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{10}\right)a\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}} & (16k)\quad \text{O III}\\
\mathbf{B}_{95} & = & \left(\frac{1}{2}-y_{10}\right)\mathbf{a}_1+x_{10}\,\mathbf{a}_2+z_{10}\,\mathbf{a}_3 & = & \left(\frac{1}{2}-y_{10}\right)a\,\hat{\mathbf{x}}+x_{10}a\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}} & (16k)\quad \text{O III}\\
\mathbf{B}_{96} & = & y_{10}\,\mathbf{a}_1+\left(\frac{1}{2}-x_{10}\right)\mathbf{a}_2+z_{10}\,\mathbf{a}_3 & = & y_{10}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}} & (16k)\quad \text{O III}\\
\mathbf{B}_{97} & = & \left(\frac{1}{2}-x_{10}\right)\mathbf{a}_1+y_{10}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{10}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{x}}+y_{10}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{10}\right)c\,\hat{\mathbf{z}} & (16k)\quad \text{O III}\\
\mathbf{B}_{98} & = & x_{10}\,\mathbf{a}_1+\left(\frac{1}{2}-y_{10}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{10}\right)\mathbf{a}_3 & = & x_{10}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{10}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{10}\right)c\,\hat{\mathbf{z}} & (16k)\quad \text{O III}\\
\mathbf{B}_{99} & = & y_{10}\,\mathbf{a}_1+x_{10}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{10}\right)\mathbf{a}_3 & = & y_{10}a\,\hat{\mathbf{x}}+x_{10}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{10}\right)c\,\hat{\mathbf{z}} & (16k)\quad \text{O III}\\
\mathbf{B}_{100} & = & \left(\frac{1}{2}-y_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}-x_{10}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{10}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-y_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{10}\right)c\,\hat{\mathbf{z}} & (16k)\quad \text{O III}\\
\mathbf{B}_{101} & = & -x_{10}\,\mathbf{a}_1-y_{10}\,\mathbf{a}_2-z_{10}\,\mathbf{a}_3 & = & -x_{10}a\,\hat{\mathbf{x}}-y_{10}a\,\hat{\mathbf{y}}-z_{10}c\,\hat{\mathbf{z}} & (16k)\quad \text{O III}\\
\mathbf{B}_{102} & = & \left(\frac{1}{2}+x_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{10}\right)\mathbf{a}_2-z_{10}\,\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{10}\right)a\,\hat{\mathbf{y}}-z_{10}c\,\hat{\mathbf{z}} & (16k)\quad \text{O III}\\
\mathbf{B}_{103} & = & \left(\frac{1}{2}+y_{10}\right)\mathbf{a}_1-x_{10}\,\mathbf{a}_2-z_{10}\,\mathbf{a}_3 & = & \left(\frac{1}{2}+y_{10}\right)a\,\hat{\mathbf{x}}-x_{10}a\,\hat{\mathbf{y}}-z_{10}c\,\hat{\mathbf{z}} & (16k)\quad \text{O III}\\
\mathbf{B}_{104} & = & -y_{10}\,\mathbf{a}_1+\left(\frac{1}{2}+x_{10}\right)\mathbf{a}_2-z_{10}\,\mathbf{a}_3 & = & -y_{10}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{y}}-z_{10}c\,\hat{\mathbf{z}} & (16k)\quad \text{O III}\\
\mathbf{B}_{105} & = & \left(\frac{1}{2}+x_{10}\right)\mathbf{a}_1-y_{10}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{10}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{x}}-y_{10}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{10}\right)c\,\hat{\mathbf{z}} & (16k)\quad \text{O III}\\
\mathbf{B}_{106} & = & -x_{10}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{10}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{10}\right)\mathbf{a}_3 & = & -x_{10}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{10}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{10}\right)c\,\hat{\mathbf{z}} & (16k)\quad \text{O III}\\
\mathbf{B}_{107} & = & -y_{10}\,\mathbf{a}_1-x_{10}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{10}\right)\mathbf{a}_3 & = & -y_{10}a\,\hat{\mathbf{x}}-x_{10}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{10}\right)c\,\hat{\mathbf{z}} & (16k)\quad \text{O III}\\
\mathbf{B}_{108} & = & \left(\frac{1}{2}+y_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}+x_{10}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{10}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+y_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{10}\right)c\,\hat{\mathbf{z}} & (16k)\quad \text{O III}\\
\mathbf{B}_{109} & = & x_{11}\,\mathbf{a}_1+y_{11}\,\mathbf{a}_2+z_{11}\,\mathbf{a}_3 & = & x_{11}a\,\hat{\mathbf{x}}+y_{11}a\,\hat{\mathbf{y}}+z_{11}c\,\hat{\mathbf{z}} & (16k)\quad \text{O IV}\\
\mathbf{B}_{110} & = & \left(\frac{1}{2}-x_{11}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{11}\right)\mathbf{a}_2+z_{11}\,\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{11}\right)a\,\hat{\mathbf{y}}+z_{11}c\,\hat{\mathbf{z}} & (16k)\quad \text{O IV}\\
\mathbf{B}_{111} & = & \left(\frac{1}{2}-y_{11}\right)\mathbf{a}_1+x_{11}\,\mathbf{a}_2+z_{11}\,\mathbf{a}_3 & = & \left(\frac{1}{2}-y_{11}\right)a\,\hat{\mathbf{x}}+x_{11}a\,\hat{\mathbf{y}}+z_{11}c\,\hat{\mathbf{z}} & (16k)\quad \text{O IV}\\
\mathbf{B}_{112} & = & y_{11}\,\mathbf{a}_1+\left(\frac{1}{2}-x_{11}\right)\mathbf{a}_2+z_{11}\,\mathbf{a}_3 & = & y_{11}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_{11}\right)a\,\hat{\mathbf{y}}+z_{11}c\,\hat{\mathbf{z}} & (16k)\quad \text{O IV}\\
\mathbf{B}_{113} & = & \left(\frac{1}{2}-x_{11}\right)\mathbf{a}_1+y_{11}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{11}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{11}\right)a\,\hat{\mathbf{x}}+y_{11}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{11}\right)c\,\hat{\mathbf{z}} & (16k)\quad \text{O IV}\\
\mathbf{B}_{114} & = & x_{11}\,\mathbf{a}_1+\left(\frac{1}{2}-y_{11}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{11}\right)\mathbf{a}_3 & = & x_{11}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{11}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{11}\right)c\,\hat{\mathbf{z}} & (16k)\quad \text{O IV}\\
\mathbf{B}_{115} & = & y_{11}\,\mathbf{a}_1+x_{11}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{11}\right)\mathbf{a}_3 & = & y_{11}a\,\hat{\mathbf{x}}+x_{11}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{11}\right)c\,\hat{\mathbf{z}} & (16k)\quad \text{O IV}\\
\mathbf{B}_{116} & = & \left(\frac{1}{2}-y_{11}\right)\mathbf{a}_1+\left(\frac{1}{2}-x_{11}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{11}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-y_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_{11}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{11}\right)c\,\hat{\mathbf{z}} & (16k)\quad \text{O IV}\\
\mathbf{B}_{117} & = & -x_{11}\,\mathbf{a}_1-y_{11}\,\mathbf{a}_2-z_{11}\,\mathbf{a}_3 & = & -x_{11}a\,\hat{\mathbf{x}}-y_{11}a\,\hat{\mathbf{y}}-z_{11}c\,\hat{\mathbf{z}} & (16k)\quad \text{O IV}\\
\mathbf{B}_{118} & = & \left(\frac{1}{2}+x_{11}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{11}\right)\mathbf{a}_2-z_{11}\,\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{11}\right)a\,\hat{\mathbf{y}}-z_{11}c\,\hat{\mathbf{z}} & (16k)\quad \text{O IV}\\
\mathbf{B}_{119} & = & \left(\frac{1}{2}+y_{11}\right)\mathbf{a}_1-x_{11}\,\mathbf{a}_2-z_{11}\,\mathbf{a}_3 & = & \left(\frac{1}{2}+y_{11}\right)a\,\hat{\mathbf{x}}-x_{11}a\,\hat{\mathbf{y}}-z_{11}c\,\hat{\mathbf{z}} & (16k)\quad \text{O IV}\\
\mathbf{B}_{120} & = & -y_{11}\,\mathbf{a}_1+\left(\frac{1}{2}+x_{11}\right)\mathbf{a}_2-z_{11}\,\mathbf{a}_3 & = & -y_{11}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_{11}\right)a\,\hat{\mathbf{y}}-z_{11}c\,\hat{\mathbf{z}} & (16k)\quad \text{O IV}\\
\mathbf{B}_{121} & = & \left(\frac{1}{2}+x_{11}\right)\mathbf{a}_1-y_{11}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{11}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{11}\right)a\,\hat{\mathbf{x}}-y_{11}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{11}\right)c\,\hat{\mathbf{z}} & (16k)\quad \text{O IV}\\
\mathbf{B}_{122} & = & -x_{11}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{11}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{11}\right)\mathbf{a}_3 & = & -x_{11}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{11}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{11}\right)c\,\hat{\mathbf{z}} & (16k)\quad \text{O IV}\\
\mathbf{B}_{123} & = & -y_{11}\,\mathbf{a}_1-x_{11}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{11}\right)\mathbf{a}_3 & = & -y_{11}a\,\hat{\mathbf{x}}-x_{11}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{11}\right)c\,\hat{\mathbf{z}} & (16k)\quad \text{O IV}\\
\mathbf{B}_{124} & = & \left(\frac{1}{2}+y_{11}\right)\mathbf{a}_1+\left(\frac{1}{2}+x_{11}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{11}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+y_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_{11}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{11}\right)c\,\hat{\mathbf{z}} & (16k)\quad \text{O IV}\\
\mathbf{B}_{125} & = & x_{12}\,\mathbf{a}_1+y_{12}\,\mathbf{a}_2+z_{12}\,\mathbf{a}_3 & = & x_{12}a\,\hat{\mathbf{x}}+y_{12}a\,\hat{\mathbf{y}}+z_{12}c\,\hat{\mathbf{z}} & (16k)\quad \text{O V}
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\begin{array}{lcccll}
\mathbf{B}_{126} & = & \left(\frac{1}{2}-x_{12}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{12}\right)\mathbf{a}_2+z_{12}\,\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{12}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{12}\right)a\,\hat{\mathbf{y}}+z_{12}c\,\hat{\mathbf{z}} & (16k)\quad \text{O V}\\
\mathbf{B}_{127} & = & \left(\frac{1}{2}-y_{12}\right)\mathbf{a}_1+x_{12}\,\mathbf{a}_2+z_{12}\,\mathbf{a}_3 & = & \left(\frac{1}{2}-y_{12}\right)a\,\hat{\mathbf{x}}+x_{12}a\,\hat{\mathbf{y}}+z_{12}c\,\hat{\mathbf{z}} & (16k)\quad \text{O V}\\
\mathbf{B}_{128} & = & y_{12}\,\mathbf{a}_1+\left(\frac{1}{2}-x_{12}\right)\mathbf{a}_2+z_{12}\,\mathbf{a}_3 & = & y_{12}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_{12}\right)a\,\hat{\mathbf{y}}+z_{12}c\,\hat{\mathbf{z}} & (16k)\quad \text{O V}\\
\mathbf{B}_{129} & = & \left(\frac{1}{2}-x_{12}\right)\mathbf{a}_1+y_{12}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{12}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{12}\right)a\,\hat{\mathbf{x}}+y_{12}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{12}\right)c\,\hat{\mathbf{z}} & (16k)\quad \text{O V}\\
\mathbf{B}_{130} & = & x_{12}\,\mathbf{a}_1+\left(\frac{1}{2}-y_{12}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{12}\right)\mathbf{a}_3 & = & x_{12}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{12}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{12}\right)c\,\hat{\mathbf{z}} & (16k)\quad \text{O V}\\
\mathbf{B}_{131} & = & y_{12}\,\mathbf{a}_1+x_{12}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{12}\right)\mathbf{a}_3 & = & y_{12}a\,\hat{\mathbf{x}}+x_{12}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{12}\right)c\,\hat{\mathbf{z}} & (16k)\quad \text{O V}\\
\mathbf{B}_{132} & = & \left(\frac{1}{2}-y_{12}\right)\mathbf{a}_1+\left(\frac{1}{2}-x_{12}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{12}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-y_{12}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_{12}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{12}\right)c\,\hat{\mathbf{z}} & (16k)\quad \text{O V}\\
\mathbf{B}_{133} & = & -x_{12}\,\mathbf{a}_1-y_{12}\,\mathbf{a}_2-z_{12}\,\mathbf{a}_3 & = & -x_{12}a\,\hat{\mathbf{x}}-y_{12}a\,\hat{\mathbf{y}}-z_{12}c\,\hat{\mathbf{z}} & (16k)\quad \text{O V}\\
\mathbf{B}_{134} & = & \left(\frac{1}{2}+x_{12}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{12}\right)\mathbf{a}_2-z_{12}\,\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{12}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{12}\right)a\,\hat{\mathbf{y}}-z_{12}c\,\hat{\mathbf{z}} & (16k)\quad \text{O V}\\
\mathbf{B}_{135} & = & \left(\frac{1}{2}+y_{12}\right)\mathbf{a}_1-x_{12}\,\mathbf{a}_2-z_{12}\,\mathbf{a}_3 & = & \left(\frac{1}{2}+y_{12}\right)a\,\hat{\mathbf{x}}-x_{12}a\,\hat{\mathbf{y}}-z_{12}c\,\hat{\mathbf{z}} & (16k)\quad \text{O V}\\
\mathbf{B}_{136} & = & -y_{12}\,\mathbf{a}_1+\left(\frac{1}{2}+x_{12}\right)\mathbf{a}_2-z_{12}\,\mathbf{a}_3 & = & -y_{12}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_{12}\right)a\,\hat{\mathbf{y}}-z_{12}c\,\hat{\mathbf{z}} & (16k)\quad \text{O V}\\
\mathbf{B}_{137} & = & \left(\frac{1}{2}+x_{12}\right)\mathbf{a}_1-y_{12}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{12}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{12}\right)a\,\hat{\mathbf{x}}-y_{12}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{12}\right)c\,\hat{\mathbf{z}} & (16k)\quad \text{O V}\\
\mathbf{B}_{138} & = & -x_{12}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{12}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{12}\right)\mathbf{a}_3 & = & -x_{12}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{12}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{12}\right)c\,\hat{\mathbf{z}} & (16k)\quad \text{O V}\\
\mathbf{B}_{139} & = & -y_{12}\,\mathbf{a}_1-x_{12}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{12}\right)\mathbf{a}_3 & = & -y_{12}a\,\hat{\mathbf{x}}-x_{12}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{12}\right)c\,\hat{\mathbf{z}} & (16k)\quad \text{O V}\\
\mathbf{B}_{140} & = & \left(\frac{1}{2}+y_{12}\right)\mathbf{a}_1+\left(\frac{1}{2}+x_{12}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{12}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+y_{12}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_{12}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{12}\right)c\,\hat{\mathbf{z}} & (16k)\quad \text{O V}\\
\mathbf{B}_{141} & = & x_{13}\,\mathbf{a}_1+y_{13}\,\mathbf{a}_2+z_{13}\,\mathbf{a}_3 & = & x_{13}a\,\hat{\mathbf{x}}+y_{13}a\,\hat{\mathbf{y}}+z_{13}c\,\hat{\mathbf{z}} & (16k)\quad \text{O VI}\\
\mathbf{B}_{142} & = & \left(\frac{1}{2}-x_{13}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{13}\right)\mathbf{a}_2+z_{13}\,\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{13}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{13}\right)a\,\hat{\mathbf{y}}+z_{13}c\,\hat{\mathbf{z}} & (16k)\quad \text{O VI}\\
\mathbf{B}_{143} & = & \left(\frac{1}{2}-y_{13}\right)\mathbf{a}_1+x_{13}\,\mathbf{a}_2+z_{13}\,\mathbf{a}_3 & = & \left(\frac{1}{2}-y_{13}\right)a\,\hat{\mathbf{x}}+x_{13}a\,\hat{\mathbf{y}}+z_{13}c\,\hat{\mathbf{z}} & (16k)\quad \text{O VI}\\
\mathbf{B}_{144} & = & y_{13}\,\mathbf{a}_1+\left(\frac{1}{2}-x_{13}\right)\mathbf{a}_2+z_{13}\,\mathbf{a}_3 & = & y_{13}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_{13}\right)a\,\hat{\mathbf{y}}+z_{13}c\,\hat{\mathbf{z}} & (16k)\quad \text{O VI}\\
\mathbf{B}_{145} & = & \left(\frac{1}{2}-x_{13}\right)\mathbf{a}_1+y_{13}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{13}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{13}\right)a\,\hat{\mathbf{x}}+y_{13}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{13}\right)c\,\hat{\mathbf{z}} & (16k)\quad \text{O VI}\\
\mathbf{B}_{146} & = & x_{13}\,\mathbf{a}_1+\left(\frac{1}{2}-y_{13}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{13}\right)\mathbf{a}_3 & = & x_{13}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{13}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{13}\right)c\,\hat{\mathbf{z}} & (16k)\quad \text{O VI}\\
\mathbf{B}_{147} & = & y_{13}\,\mathbf{a}_1+x_{13}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{13}\right)\mathbf{a}_3 & = & y_{13}a\,\hat{\mathbf{x}}+x_{13}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{13}\right)c\,\hat{\mathbf{z}} & (16k)\quad \text{O VI}\\
\mathbf{B}_{148} & = & \left(\frac{1}{2}-y_{13}\right)\mathbf{a}_1+\left(\frac{1}{2}-x_{13}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{13}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-y_{13}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_{13}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{13}\right)c\,\hat{\mathbf{z}} & (16k)\quad \text{O VI}\\
\mathbf{B}_{149} & = & -x_{13}\,\mathbf{a}_1-y_{13}\,\mathbf{a}_2-z_{13}\,\mathbf{a}_3 & = & -x_{13}a\,\hat{\mathbf{x}}-y_{13}a\,\hat{\mathbf{y}}-z_{13}c\,\hat{\mathbf{z}} & (16k)\quad \text{O VI}\\
\mathbf{B}_{150} & = & \left(\frac{1}{2}+x_{13}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{13}\right)\mathbf{a}_2-z_{13}\,\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{13}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{13}\right)a\,\hat{\mathbf{y}}-z_{13}c\,\hat{\mathbf{z}} & (16k)\quad \text{O VI}\\
\mathbf{B}_{151} & = & \left(\frac{1}{2}+y_{13}\right)\mathbf{a}_1-x_{13}\,\mathbf{a}_2-z_{13}\,\mathbf{a}_3 & = & \left(\frac{1}{2}+y_{13}\right)a\,\hat{\mathbf{x}}-x_{13}a\,\hat{\mathbf{y}}-z_{13}c\,\hat{\mathbf{z}} & (16k)\quad \text{O VI}\\
\mathbf{B}_{152} & = & -y_{13}\,\mathbf{a}_1+\left(\frac{1}{2}+x_{13}\right)\mathbf{a}_2-z_{13}\,\mathbf{a}_3 & = & -y_{13}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_{13}\right)a\,\hat{\mathbf{y}}-z_{13}c\,\hat{\mathbf{z}} & (16k)\quad \text{O VI}\\
\mathbf{B}_{153} & = & \left(\frac{1}{2}+x_{13}\right)\mathbf{a}_1-y_{13}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{13}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{13}\right)a\,\hat{\mathbf{x}}-y_{13}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{13}\right)c\,\hat{\mathbf{z}} & (16k)\quad \text{O VI}\\
\mathbf{B}_{154} & = & -x_{13}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{13}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{13}\right)\mathbf{a}_3 & = & -x_{13}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{13}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{13}\right)c\,\hat{\mathbf{z}} & (16k)\quad \text{O VI}\\
\mathbf{B}_{155} & = & -y_{13}\,\mathbf{a}_1-x_{13}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{13}\right)\mathbf{a}_3 & = & -y_{13}a\,\hat{\mathbf{x}}-x_{13}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{13}\right)c\,\hat{\mathbf{z}} & (16k)\quad \text{O VI}\\
\mathbf{B}_{156} & = & \left(\frac{1}{2}+y_{13}\right)\mathbf{a}_1+\left(\frac{1}{2}+x_{13}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{13}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+y_{13}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_{13}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{13}\right)c\,\hat{\mathbf{z}} & (16k)\quad \text{O VI}\\
\mathbf{B}_{157} & = & x_{14}\,\mathbf{a}_1+y_{14}\,\mathbf{a}_2+z_{14}\,\mathbf{a}_3 & = & x_{14}a\,\hat{\mathbf{x}}+y_{14}a\,\hat{\mathbf{y}}+z_{14}c\,\hat{\mathbf{z}} & (16k)\quad \text{O VII}
\end{array}
$$

$$\mathbf{B}_{158} = \left(\tfrac{1}{2}-x_{14}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{14}\right)\mathbf{a}_2+z_{14}\,\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{14}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{14}\right)a\,\hat{\mathbf{y}}+z_{14}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O VII}$$

$$\mathbf{B}_{159} = \left(\tfrac{1}{2}-y_{14}\right)\mathbf{a}_1+x_{14}\,\mathbf{a}_2+z_{14}\,\mathbf{a}_3 = \left(\tfrac{1}{2}-y_{14}\right)a\,\hat{\mathbf{x}}+x_{14}a\,\hat{\mathbf{y}}+z_{14}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O VII}$$

$$\mathbf{B}_{160} = y_{14}\,\mathbf{a}_1+\left(\tfrac{1}{2}-x_{14}\right)\mathbf{a}_2+z_{14}\,\mathbf{a}_3 = y_{14}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_{14}\right)a\,\hat{\mathbf{y}}+z_{14}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O VII}$$

$$\mathbf{B}_{161} = \left(\tfrac{1}{2}-x_{14}\right)\mathbf{a}_1+y_{14}\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_{14}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{14}\right)a\,\hat{\mathbf{x}}+y_{14}a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{14}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O VII}$$

$$\mathbf{B}_{162} = x_{14}\,\mathbf{a}_1+\left(\tfrac{1}{2}-y_{14}\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_{14}\right)\mathbf{a}_3 = x_{14}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{14}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{14}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O VII}$$

$$\mathbf{B}_{163} = y_{14}\,\mathbf{a}_1+x_{14}\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_{14}\right)\mathbf{a}_3 = y_{14}a\,\hat{\mathbf{x}}+x_{14}a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{14}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O VII}$$

$$\mathbf{B}_{164} = \left(\tfrac{1}{2}-y_{14}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-x_{14}\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_{14}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-y_{14}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_{14}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{14}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O VII}$$

$$\mathbf{B}_{165} = -x_{14}\,\mathbf{a}_1-y_{14}\,\mathbf{a}_2-z_{14}\,\mathbf{a}_3 = -x_{14}a\,\hat{\mathbf{x}}-y_{14}a\,\hat{\mathbf{y}}-z_{14}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O VII}$$

$$\mathbf{B}_{166} = \left(\tfrac{1}{2}+x_{14}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{14}\right)\mathbf{a}_2-z_{14}\,\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{14}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{14}\right)a\,\hat{\mathbf{y}}-z_{14}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O VII}$$

$$\mathbf{B}_{167} = \left(\tfrac{1}{2}+y_{14}\right)\mathbf{a}_1-x_{14}\,\mathbf{a}_2-z_{14}\,\mathbf{a}_3 = \left(\tfrac{1}{2}+y_{14}\right)a\,\hat{\mathbf{x}}-x_{14}a\,\hat{\mathbf{y}}-z_{14}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O VII}$$

$$\mathbf{B}_{168} = -y_{14}\,\mathbf{a}_1+\left(\tfrac{1}{2}+x_{14}\right)\mathbf{a}_2-z_{14}\,\mathbf{a}_3 = -y_{14}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_{14}\right)a\,\hat{\mathbf{y}}-z_{14}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O VII}$$

$$\mathbf{B}_{169} = \left(\tfrac{1}{2}+x_{14}\right)\mathbf{a}_1-y_{14}\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_{14}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{14}\right)a\,\hat{\mathbf{x}}-y_{14}a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{14}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O VII}$$

$$\mathbf{B}_{170} = -x_{14}\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_{14}\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_{14}\right)\mathbf{a}_3 = -x_{14}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{14}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{14}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O VII}$$

$$\mathbf{B}_{171} = -y_{14}\,\mathbf{a}_1-x_{14}\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_{14}\right)\mathbf{a}_3 = -y_{14}a\,\hat{\mathbf{x}}-x_{14}a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{14}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O VII}$$

$$\mathbf{B}_{172} = \left(\tfrac{1}{2}+y_{14}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_{14}\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_{14}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+y_{14}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_{14}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{14}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O VII}$$

$$\mathbf{B}_{173} = x_{15}\,\mathbf{a}_1+y_{15}\,\mathbf{a}_2+z_{15}\,\mathbf{a}_3 = x_{15}a\,\hat{\mathbf{x}}+y_{15}a\,\hat{\mathbf{y}}+z_{15}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O VIII}$$

$$\mathbf{B}_{174} = \left(\tfrac{1}{2}-x_{15}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{15}\right)\mathbf{a}_2+z_{15}\,\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{15}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{15}\right)a\,\hat{\mathbf{y}}+z_{15}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O VIII}$$

$$\mathbf{B}_{175} = \left(\tfrac{1}{2}-y_{15}\right)\mathbf{a}_1+x_{15}\,\mathbf{a}_2+z_{15}\,\mathbf{a}_3 = \left(\tfrac{1}{2}-y_{15}\right)a\,\hat{\mathbf{x}}+x_{15}a\,\hat{\mathbf{y}}+z_{15}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O VIII}$$

$$\mathbf{B}_{176} = y_{15}\,\mathbf{a}_1+\left(\tfrac{1}{2}-x_{15}\right)\mathbf{a}_2+z_{15}\,\mathbf{a}_3 = y_{15}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_{15}\right)a\,\hat{\mathbf{y}}+z_{15}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O VIII}$$

$$\mathbf{B}_{177} = \left(\tfrac{1}{2}-x_{15}\right)\mathbf{a}_1+y_{15}\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_{15}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{15}\right)a\,\hat{\mathbf{x}}+y_{15}a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{15}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O VIII}$$

$$\mathbf{B}_{178} = x_{15}\,\mathbf{a}_1+\left(\tfrac{1}{2}-y_{15}\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_{15}\right)\mathbf{a}_3 = x_{15}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{15}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{15}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O VIII}$$

$$\mathbf{B}_{179} = y_{15}\,\mathbf{a}_1+x_{15}\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_{15}\right)\mathbf{a}_3 = y_{15}a\,\hat{\mathbf{x}}+x_{15}a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{15}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O VIII}$$

$$\mathbf{B}_{180} = \left(\tfrac{1}{2}-y_{15}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-x_{15}\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_{15}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-y_{15}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_{15}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{15}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O VIII}$$

$$\mathbf{B}_{181} = -x_{15}\,\mathbf{a}_1-y_{15}\,\mathbf{a}_2-z_{15}\,\mathbf{a}_3 = -x_{15}a\,\hat{\mathbf{x}}-y_{15}a\,\hat{\mathbf{y}}-z_{15}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O VIII}$$

$$\mathbf{B}_{182} = \left(\tfrac{1}{2}+x_{15}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{15}\right)\mathbf{a}_2-z_{15}\,\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{15}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{15}\right)a\,\hat{\mathbf{y}}-z_{15}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O VIII}$$

$$\mathbf{B}_{183} = \left(\tfrac{1}{2}+y_{15}\right)\mathbf{a}_1-x_{15}\,\mathbf{a}_2-z_{15}\,\mathbf{a}_3 = \left(\tfrac{1}{2}+y_{15}\right)a\,\hat{\mathbf{x}}-x_{15}a\,\hat{\mathbf{y}}-z_{15}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O VIII}$$

$$\mathbf{B}_{184} = -y_{15}\,\mathbf{a}_1+\left(\tfrac{1}{2}+x_{15}\right)\mathbf{a}_2-z_{15}\,\mathbf{a}_3 = -y_{15}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_{15}\right)a\,\hat{\mathbf{y}}-z_{15}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O VIII}$$

$$\mathbf{B}_{185} = \left(\tfrac{1}{2}+x_{15}\right)\mathbf{a}_1-y_{15}\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_{15}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{15}\right)a\,\hat{\mathbf{x}}-y_{15}a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{15}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O VIII}$$

$$\mathbf{B}_{186} = -x_{15}\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_{15}\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_{15}\right)\mathbf{a}_3 = -x_{15}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{15}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{15}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O VIII}$$

$$\mathbf{B}_{187} = -y_{15}\,\mathbf{a}_1-x_{15}\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_{15}\right)\mathbf{a}_3 = -y_{15}a\,\hat{\mathbf{x}}-x_{15}a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{15}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O VIII}$$

$$\mathbf{B}_{188} = \left(\tfrac{1}{2}+y_{15}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_{15}\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_{15}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+y_{15}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_{15}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{15}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O VIII}$$

$$\mathbf{B}_{189} = x_{16}\,\mathbf{a}_1+y_{16}\,\mathbf{a}_2+z_{16}\,\mathbf{a}_3 = x_{16}a\,\hat{\mathbf{x}}+y_{16}a\,\hat{\mathbf{y}}+z_{16}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O IX}$$

$$\mathbf{B}_{190} = \left(\tfrac{1}{2} - x_{16}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{16}\right)\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{16}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{16}\right)a\,\hat{\mathbf{y}} + z_{16}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O IX}$$

$$\mathbf{B}_{191} = \left(\tfrac{1}{2} - y_{16}\right)\mathbf{a}_1 + x_{16}\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 = \left(\tfrac{1}{2} - y_{16}\right)a\,\hat{\mathbf{x}} + x_{16}a\,\hat{\mathbf{y}} + z_{16}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O IX}$$

$$\mathbf{B}_{192} = y_{16}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_{16}\right)\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 = y_{16}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_{16}\right)a\,\hat{\mathbf{y}} + z_{16}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O IX}$$

$$\mathbf{B}_{193} = \left(\tfrac{1}{2} - x_{16}\right)\mathbf{a}_1 + y_{16}\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{16}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{16}\right)a\,\hat{\mathbf{x}} + y_{16}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{16}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O IX}$$

$$\mathbf{B}_{194} = x_{16}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{16}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{16}\right)\mathbf{a}_3 = x_{16}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{16}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{16}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O IX}$$

$$\mathbf{B}_{195} = y_{16}\,\mathbf{a}_1 + x_{16}\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{16}\right)\mathbf{a}_3 = y_{16}a\,\hat{\mathbf{x}} + x_{16}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{16}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O IX}$$

$$\mathbf{B}_{196} = \left(\tfrac{1}{2} - y_{16}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_{16}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{16}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - y_{16}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_{16}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{16}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O IX}$$

$$\mathbf{B}_{197} = -x_{16}\,\mathbf{a}_1 - y_{16}\,\mathbf{a}_2 - z_{16}\,\mathbf{a}_3 = -x_{16}a\,\hat{\mathbf{x}} - y_{16}a\,\hat{\mathbf{y}} - z_{16}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O IX}$$

$$\mathbf{B}_{198} = \left(\tfrac{1}{2} + x_{16}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{16}\right)\mathbf{a}_2 - z_{16}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{16}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{16}\right)a\,\hat{\mathbf{y}} - z_{16}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O IX}$$

$$\mathbf{B}_{199} = \left(\tfrac{1}{2} + y_{16}\right)\mathbf{a}_1 - x_{16}\,\mathbf{a}_2 - z_{16}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + y_{16}\right)a\,\hat{\mathbf{x}} - x_{16}a\,\hat{\mathbf{y}} - z_{16}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O IX}$$

$$\mathbf{B}_{200} = -y_{16}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_{16}\right)\mathbf{a}_2 - z_{16}\,\mathbf{a}_3 = -y_{16}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_{16}\right)a\,\hat{\mathbf{y}} - z_{16}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O IX}$$

$$\mathbf{B}_{201} = \left(\tfrac{1}{2} + x_{16}\right)\mathbf{a}_1 - y_{16}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{16}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{16}\right)a\,\hat{\mathbf{x}} - y_{16}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{16}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O IX}$$

$$\mathbf{B}_{202} = -x_{16}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{16}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{16}\right)\mathbf{a}_3 = -x_{16}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{16}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{16}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O IX}$$

$$\mathbf{B}_{203} = -y_{16}\,\mathbf{a}_1 - x_{16}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{16}\right)\mathbf{a}_3 = -y_{16}a\,\hat{\mathbf{x}} - x_{16}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{16}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O IX}$$

$$\mathbf{B}_{204} = \left(\tfrac{1}{2} + y_{16}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_{16}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{16}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + y_{16}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_{16}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{16}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{O IX}$$

$$\mathbf{B}_{205} = x_{17}\,\mathbf{a}_1 + y_{17}\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 = x_{17}a\,\hat{\mathbf{x}} + y_{17}a\,\hat{\mathbf{y}} + z_{17}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{OH}$$

$$\mathbf{B}_{206} = \left(\tfrac{1}{2} - x_{17}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{17}\right)\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{17}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{17}\right)a\,\hat{\mathbf{y}} + z_{17}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{OH}$$

$$\mathbf{B}_{207} = \left(\tfrac{1}{2} - y_{17}\right)\mathbf{a}_1 + x_{17}\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 = \left(\tfrac{1}{2} - y_{17}\right)a\,\hat{\mathbf{x}} + x_{17}a\,\hat{\mathbf{y}} + z_{17}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{OH}$$

$$\mathbf{B}_{208} = y_{17}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_{17}\right)\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 = y_{17}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_{17}\right)a\,\hat{\mathbf{y}} + z_{17}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{OH}$$

$$\mathbf{B}_{209} = \left(\tfrac{1}{2} - x_{17}\right)\mathbf{a}_1 + y_{17}\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{17}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{17}\right)a\,\hat{\mathbf{x}} + y_{17}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{17}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{OH}$$

$$\mathbf{B}_{210} = x_{17}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{17}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{17}\right)\mathbf{a}_3 = x_{17}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{17}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{17}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{OH}$$

$$\mathbf{B}_{211} = y_{17}\,\mathbf{a}_1 + x_{17}\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{17}\right)\mathbf{a}_3 = y_{17}a\,\hat{\mathbf{x}} + x_{17}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{17}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{OH}$$

$$\mathbf{B}_{212} = \left(\tfrac{1}{2} - y_{17}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_{17}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{17}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - y_{17}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_{17}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{17}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{OH}$$

$$\mathbf{B}_{213} = -x_{17}\,\mathbf{a}_1 - y_{17}\,\mathbf{a}_2 - z_{17}\,\mathbf{a}_3 = -x_{17}a\,\hat{\mathbf{x}} - y_{17}a\,\hat{\mathbf{y}} - z_{17}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{OH}$$

$$\mathbf{B}_{214} = \left(\tfrac{1}{2} + x_{17}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{17}\right)\mathbf{a}_2 - z_{17}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{17}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{17}\right)a\,\hat{\mathbf{y}} - z_{17}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{OH}$$

$$\mathbf{B}_{215} = \left(\tfrac{1}{2} + y_{17}\right)\mathbf{a}_1 - x_{17}\,\mathbf{a}_2 - z_{17}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + y_{17}\right)a\,\hat{\mathbf{x}} - x_{17}a\,\hat{\mathbf{y}} - z_{17}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{OH}$$

$$\mathbf{B}_{216} = -y_{17}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_{17}\right)\mathbf{a}_2 - z_{17}\,\mathbf{a}_3 = -y_{17}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_{17}\right)a\,\hat{\mathbf{y}} - z_{17}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{OH}$$

$$\mathbf{B}_{217} = \left(\tfrac{1}{2} + x_{17}\right)\mathbf{a}_1 - y_{17}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{17}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{17}\right)a\,\hat{\mathbf{x}} - y_{17}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{17}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{OH}$$

$$\mathbf{B}_{218} = -x_{17}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{17}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{17}\right)\mathbf{a}_3 = -x_{17}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{17}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{17}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{OH}$$

$$\mathbf{B}_{219} = -y_{17}\,\mathbf{a}_1 - x_{17}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{17}\right)\mathbf{a}_3 = -y_{17}a\,\hat{\mathbf{x}} - x_{17}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{17}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{OH}$$

$$\mathbf{B}_{220} = \left(\tfrac{1}{2} + y_{17}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_{17}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{17}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + y_{17}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_{17}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{17}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{OH}$$

$$\mathbf{B}_{221} = x_{18}\,\mathbf{a}_1 + y_{18}\,\mathbf{a}_2 + z_{18}\,\mathbf{a}_3 = x_{18}a\,\hat{\mathbf{x}} + y_{18}a\,\hat{\mathbf{y}} + z_{18}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{Si II}$$

$$\mathbf{B}_{222} = \left(\tfrac{1}{2} - x_{18}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{18}\right)\mathbf{a}_2 + z_{18}\,\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{18}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{18}\right)a\,\hat{\mathbf{y}} + z_{18}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{Si II}$$

$$\mathbf{B}_{223} = \left(\tfrac{1}{2} - y_{18}\right)\mathbf{a}_1 + x_{18}\,\mathbf{a}_2 + z_{18}\,\mathbf{a}_3 = \left(\tfrac{1}{2} - y_{18}\right)a\,\hat{\mathbf{x}} + x_{18}a\,\hat{\mathbf{y}} + z_{18}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{Si II}$$

$$\mathbf{B}_{224} = y_{18}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_{18}\right)\mathbf{a}_2 + z_{18}\,\mathbf{a}_3 = y_{18}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_{18}\right)a\,\hat{\mathbf{y}} + z_{18}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{Si II}$$

$$\mathbf{B}_{225} = \left(\tfrac{1}{2} - x_{18}\right)\mathbf{a}_1 + y_{18}\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{18}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{18}\right)a\,\hat{\mathbf{x}} + y_{18}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{18}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{Si II}$$

$$\mathbf{B}_{226} = x_{18}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{18}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{18}\right)\mathbf{a}_3 = x_{18}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{18}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{18}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{Si II}$$

$$\mathbf{B}_{227} = y_{18}\,\mathbf{a}_1 + x_{18}\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{18}\right)\mathbf{a}_3 = y_{18}a\,\hat{\mathbf{x}} + x_{18}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{18}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{Si II}$$

$$\mathbf{B}_{228} = \left(\tfrac{1}{2} - y_{18}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_{18}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{18}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - y_{18}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_{18}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{18}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{Si II}$$

$$\mathbf{B}_{229} = -x_{18}\,\mathbf{a}_1 - y_{18}\,\mathbf{a}_2 - z_{18}\,\mathbf{a}_3 = -x_{18}a\,\hat{\mathbf{x}} - y_{18}a\,\hat{\mathbf{y}} - z_{18}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{Si II}$$

$$\mathbf{B}_{230} = \left(\tfrac{1}{2} + x_{18}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{18}\right)\mathbf{a}_2 - z_{18}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{18}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{18}\right)a\,\hat{\mathbf{y}} - z_{18}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{Si II}$$

$$\mathbf{B}_{231} = \left(\tfrac{1}{2} + y_{18}\right)\mathbf{a}_1 - x_{18}\,\mathbf{a}_2 - z_{18}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + y_{18}\right)a\,\hat{\mathbf{x}} - x_{18}a\,\hat{\mathbf{y}} - z_{18}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{Si II}$$

$$\mathbf{B}_{232} = -y_{18}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_{18}\right)\mathbf{a}_2 - z_{18}\,\mathbf{a}_3 = -y_{18}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_{18}\right)a\,\hat{\mathbf{y}} - z_{18}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{Si II}$$

$$\mathbf{B}_{233} = \left(\tfrac{1}{2} + x_{18}\right)\mathbf{a}_1 - y_{18}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{18}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{18}\right)a\,\hat{\mathbf{x}} - y_{18}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{18}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{Si II}$$

$$\mathbf{B}_{234} = -x_{18}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{18}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{18}\right)\mathbf{a}_3 = -x_{18}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{18}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{18}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{Si II}$$

$$\mathbf{B}_{235} = -y_{18}\,\mathbf{a}_1 - x_{18}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{18}\right)\mathbf{a}_3 = -y_{18}a\,\hat{\mathbf{x}} - x_{18}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{18}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{Si II}$$

$$\mathbf{B}_{236} = \left(\tfrac{1}{2} + y_{18}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_{18}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{18}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + y_{18}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_{18}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{18}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{Si II}$$

$$\mathbf{B}_{237} = x_{19}\,\mathbf{a}_1 + y_{19}\,\mathbf{a}_2 + z_{19}\,\mathbf{a}_3 = x_{19}a\,\hat{\mathbf{x}} + y_{19}a\,\hat{\mathbf{y}} + z_{19}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{Si III}$$

$$\mathbf{B}_{238} = \left(\tfrac{1}{2} - x_{19}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{19}\right)\mathbf{a}_2 + z_{19}\,\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{19}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{19}\right)a\,\hat{\mathbf{y}} + z_{19}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{Si III}$$

$$\mathbf{B}_{239} = \left(\tfrac{1}{2} - y_{19}\right)\mathbf{a}_1 + x_{19}\,\mathbf{a}_2 + z_{19}\,\mathbf{a}_3 = \left(\tfrac{1}{2} - y_{19}\right)a\,\hat{\mathbf{x}} + x_{19}a\,\hat{\mathbf{y}} + z_{19}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{Si III}$$

$$\mathbf{B}_{240} = y_{19}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_{19}\right)\mathbf{a}_2 + z_{19}\,\mathbf{a}_3 = y_{19}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_{19}\right)a\,\hat{\mathbf{y}} + z_{19}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{Si III}$$

$$\mathbf{B}_{241} = \left(\tfrac{1}{2} - x_{19}\right)\mathbf{a}_1 + y_{19}\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{19}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{19}\right)a\,\hat{\mathbf{x}} + y_{19}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{19}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{Si III}$$

$$\mathbf{B}_{242} = x_{19}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{19}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{19}\right)\mathbf{a}_3 = x_{19}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{19}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{19}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{Si III}$$

$$\mathbf{B}_{243} = y_{19}\,\mathbf{a}_1 + x_{19}\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{19}\right)\mathbf{a}_3 = y_{19}a\,\hat{\mathbf{x}} + x_{19}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{19}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{Si III}$$

$$\mathbf{B}_{244} = \left(\tfrac{1}{2} - y_{19}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_{19}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{19}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - y_{19}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_{19}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{19}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{Si III}$$

$$\mathbf{B}_{245} = -x_{19}\,\mathbf{a}_1 - y_{19}\,\mathbf{a}_2 - z_{19}\,\mathbf{a}_3 = -x_{19}a\,\hat{\mathbf{x}} - y_{19}a\,\hat{\mathbf{y}} - z_{19}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{Si III}$$

$$\mathbf{B}_{246} = \left(\tfrac{1}{2} + x_{19}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{19}\right)\mathbf{a}_2 - z_{19}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{19}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{19}\right)a\,\hat{\mathbf{y}} - z_{19}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{Si III}$$

$$\mathbf{B}_{247} = \left(\tfrac{1}{2} + y_{19}\right)\mathbf{a}_1 - x_{19}\,\mathbf{a}_2 - z_{19}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + y_{19}\right)a\,\hat{\mathbf{x}} - x_{19}a\,\hat{\mathbf{y}} - z_{19}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{Si III}$$

$$\mathbf{B}_{248} = -y_{19}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_{19}\right)\mathbf{a}_2 - z_{19}\,\mathbf{a}_3 = -y_{19}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_{19}\right)a\,\hat{\mathbf{y}} - z_{19}c\,\hat{\mathbf{z}} \quad (16k) \quad \text{Si III}$$

$$\mathbf{B}_{249} = \left(\tfrac{1}{2} + x_{19}\right)\mathbf{a}_1 - y_{19}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{19}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{19}\right)a\,\hat{\mathbf{x}} - y_{19}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{19}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{Si III}$$

$$\mathbf{B}_{250} = -x_{19}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{19}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{19}\right)\mathbf{a}_3 = -x_{19}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{19}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{19}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{Si III}$$

$$\mathbf{B}_{251} = -y_{19}\,\mathbf{a}_1 - x_{19}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{19}\right)\mathbf{a}_3 = -y_{19}a\,\hat{\mathbf{x}} - x_{19}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{19}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{Si III}$$

$$\mathbf{B}_{252} = \left(\tfrac{1}{2} + y_{19}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_{19}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{19}\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + y_{19}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_{19}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{19}\right)c\,\hat{\mathbf{z}} \quad (16k) \quad \text{Si III}$$

References:

- B. E. Warren and D. I. Modell, *The Structure of Vesuvianite* $Ca_{10}Al_4(Mg,Fe)_2Si_9O_{34}(OH)_4$, Zeitschrift für Kristallographie - Crystalline Materials **78**, 422–432 (1931), doi:10.1524/zkri.1931.78.1.422.
- F. M. Allen and C. W. Burnham, *A comprehensive structure-model for vesuvianite; symmetry variations and crystal growth*, Can. Mineral. **30**, 1–18 (1992).
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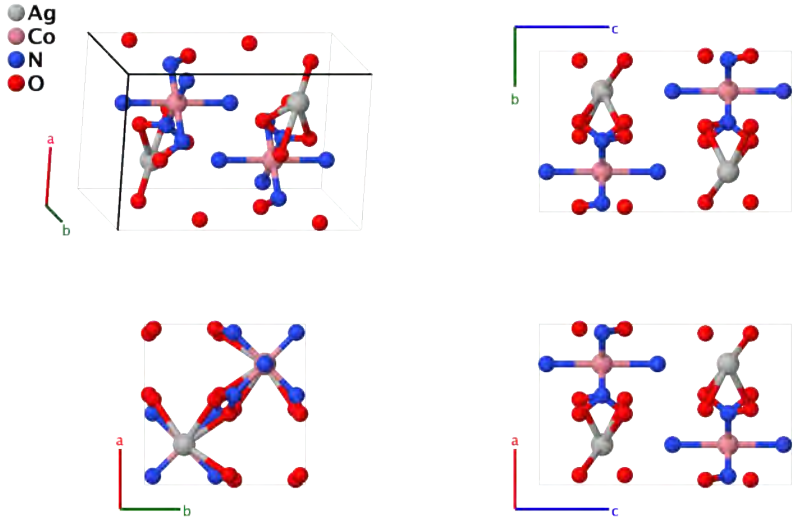
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1692
- POSCAR: pp. 1693

$Ag[Co(NH_3)_2(NO_2)_4]$ ($J1_9$) Structure: ABC4D2E8_tP32_126_a_b_h_e_k

http://aflow.org/prototype-encyclopedia/ABC4D2E8_tP32_126_a_b_h_e_k

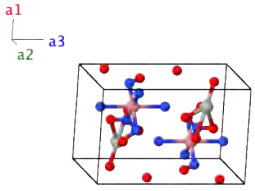


Prototype	:	$AgCoN_4(NH_3)_2O_8$
AFLOW prototype label	:	ABC4D2E8_tP32_126_a_b_h_e_k
***Strukturbericht* designation**	:	$J1_9$
Pearson symbol	:	tP32
Space group number	:	126
Space group symbol	:	$P4/nnc$
AFLOW prototype command	:	`aflow --proto=ABC4D2E8_tP32_126_a_b_h_e_k --params=`$a, c/a, z_3, x_4, x_5, y_5, z_5$

- The positions of the hydrogen atoms in the ammonia molecule are not determined, so we only provide the positions of the nitrogen atoms (labeled as NH_3).
- (Wells, 1936) gives the lattice coordinates in setting 1 of space group #126. We used FINDSYM to change this to the standard setting 2.

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$ =	$\frac{1}{4}\mathbf{a}_1 + \frac{1}{4}\mathbf{a}_2 + \frac{1}{4}\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($2a$)	Ag

$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(2*a*)	Ag
$\mathbf{B}_3$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(2*b*)	Co
$\mathbf{B}_4$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(2*b*)	Co
$\mathbf{B}_5$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	(4*e*)	NH_3
$\mathbf{B}_6$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(4*e*)	NH_3
$\mathbf{B}_7$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	(4*e*)	NH_3
$\mathbf{B}_8$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(4*e*)	NH_3
$\mathbf{B}_9$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(8*h*)	N
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_4\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - x_4\right)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(8*h*)	N
$\mathbf{B}_{11}$	=	$\left(\frac{1}{2} - x_4\right)\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(8*h*)	N
$\mathbf{B}_{12}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2} - x_4\right)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(8*h*)	N
$\mathbf{B}_{13}$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(8*h*)	N
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} + x_4\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + x_4\right)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(8*h*)	N
$\mathbf{B}_{15}$	=	$\left(\frac{1}{2} + x_4\right)\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(8*h*)	N
$\mathbf{B}_{16}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + x_4\right)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(8*h*)	N
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}} + y_5a\,\hat{\mathbf{y}} + z_5c\,\hat{\mathbf{z}}$	(16*k*)	O
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} - x_5\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right)a\,\hat{\mathbf{y}} + z_5c\,\hat{\mathbf{z}}$	(16*k*)	O
$\mathbf{B}_{19}$	=	$\left(\frac{1}{2} - y_5\right)\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_5\right)a\,\hat{\mathbf{x}} + x_5a\,\hat{\mathbf{y}} + z_5c\,\hat{\mathbf{z}}$	(16*k*)	O
$\mathbf{B}_{20}$	=	$y_5\,\mathbf{a}_1 + \left(\frac{1}{2} - x_5\right)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$y_5a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{y}} + z_5c\,\hat{\mathbf{z}}$	(16*k*)	O
$\mathbf{B}_{21}$	=	$\left(\frac{1}{2} - x_5\right)\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + y_5a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	(16*k*)	O
$\mathbf{B}_{22}$	=	$x_5\,\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	(16*k*)	O
$\mathbf{B}_{23}$	=	$y_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\,\mathbf{a}_3$	=	$y_5a\,\hat{\mathbf{x}} + x_5a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	(16*k*)	O
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} - y_5\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - x_5\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	(16*k*)	O
$\mathbf{B}_{25}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}} - y_5a\,\hat{\mathbf{y}} - z_5c\,\hat{\mathbf{z}}$	(16*k*)	O
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} + x_5\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right)a\,\hat{\mathbf{y}} - z_5c\,\hat{\mathbf{z}}$	(16*k*)	O
$\mathbf{B}_{27}$	=	$\left(\frac{1}{2} + y_5\right)\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_5\right)a\,\hat{\mathbf{x}} - x_5a\,\hat{\mathbf{y}} - z_5c\,\hat{\mathbf{z}}$	(16*k*)	O
$\mathbf{B}_{28}$	=	$-y_5\,\mathbf{a}_1 + \left(\frac{1}{2} + x_5\right)\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-y_5a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{y}} - z_5c\,\hat{\mathbf{z}}$	(16*k*)	O
$\mathbf{B}_{29}$	=	$\left(\frac{1}{2} + x_5\right)\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} - y_5a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	(16*k*)	O
$\mathbf{B}_{30}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	(16*k*)	O
$\mathbf{B}_{31}$	=	$-y_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\,\mathbf{a}_3$	=	$-y_5a\,\hat{\mathbf{x}} - x_5a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	(16*k*)	O
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2} + y_5\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + x_5\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	(16*k*)	O

References:

- A. F. Wells, *The Crystal Structure of Silver Diamminotetranitro-cobaltiate Ag[Co(NH$_3$)$_2$(NO$_2$)$_4$]*, Zeitschrift für Kristallographie - Crystalline Materials **95**, 74–82 (1936), doi:10.1524/zkri.1936.95.1.74.

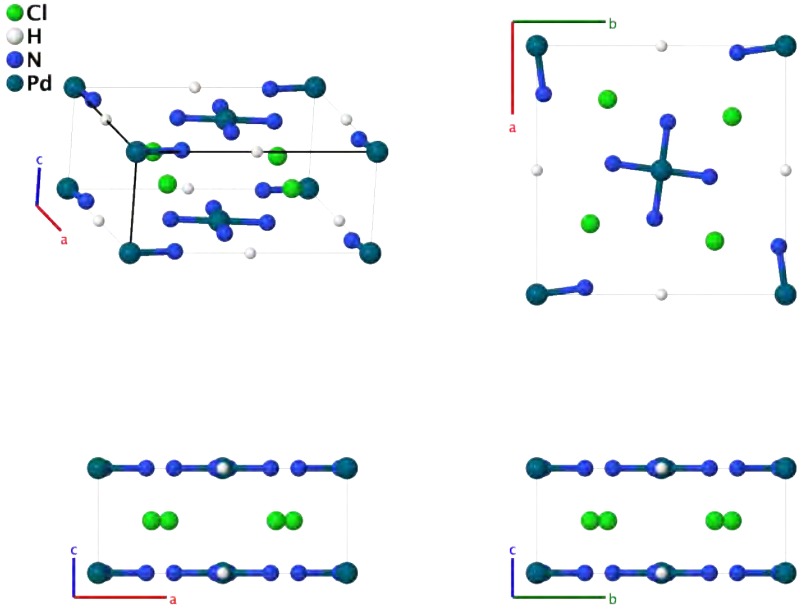
Found in:

- C. Gottfried, ed., *Strukturbericht Band IV 1936* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1938).

Geometry files:
- CIF: pp. 1694
- POSCAR: pp. 1694

$Pd(NH_3)_4Cl_2{\cdot}H_2O$ ($H4_9$) Structure: A2BC4D_tP16_127_h_d_i_a

http://aflow.org/prototype-encyclopedia/A2BC4D_tP16_127_h_d_i_a

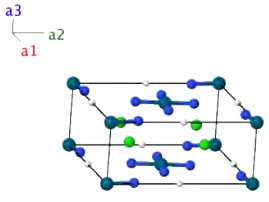


Prototype	:	$Cl_2(H_2O)(NH_3)_4Pd$
AFLOW prototype label	:	A2BC4D_tP16_127_h_d_i_a
***Strukturbericht* designation**	:	$H4_9$
Pearson symbol	:	tP16
Space group number	:	127
Space group symbol	:	$P4/mbm$
AFLOW prototype command	:	`aflow --proto=A2BC4D_tP16_127_h_d_i_a --params=`$a, c/a, x_3, x_4, y_4$

- The positions of the hydrogen atoms have not been measured, assuming they are actually fixed. We group them together with the central atoms of their molecules.

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(2*a*)	Pd
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$	(2*a*)	Pd
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{y}}$	(2*d*)	H_2O
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	(2*d*)	H_2O
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*h*)	Cl
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*h*)	Cl
$\mathbf{B}_7$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*h*)	Cl
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*h*)	Cl
$\mathbf{B}_9$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 a\,\hat{\mathbf{y}}$	(8*i*)	NH_3
$\mathbf{B}_{10}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 a\,\hat{\mathbf{y}}$	(8*i*)	NH_3
$\mathbf{B}_{11}$	=	$-y_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2$	=	$-y_4 a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}}$	(8*i*)	NH_3
$\mathbf{B}_{12}$	=	$y_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2$	=	$y_4 a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}}$	(8*i*)	NH_3
$\mathbf{B}_{13}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)a\,\hat{\mathbf{y}}$	(8*i*)	NH_3
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)a\,\hat{\mathbf{y}}$	(8*i*)	NH_3
$\mathbf{B}_{15}$	=	$\left(\frac{1}{2} + y_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_4\right)\mathbf{a}_2$	=	$\left(\frac{1}{2} + y_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{y}}$	(8*i*)	NH_3
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} - y_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_2$	=	$\left(\frac{1}{2} - y_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{y}}$	(8*i*)	NH_3

References:

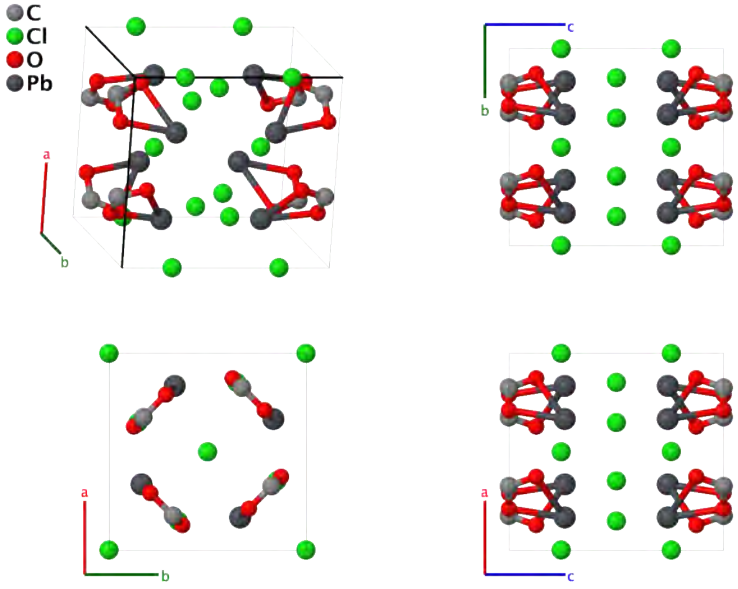
- B. N. Dickinson, *The Crystal Structure of Tetramminopalladous Chloride $Pd(NH_3)_4Cl_2{\cdot}H_2O$*, Zeitschrift für Kristallographie - Crystalline Materials **88**, 281–297 (1934), doi:10.1524/zkri.1934.88.1.281.

Geometry files:

- CIF: pp. 1695
- POSCAR: pp. 1695

Phosgenite [$Pb_2Cl_2(CO_3)$] Structure: AB2C3D2_tP32_127_g_eh_gk_k

http://aflow.org/prototype-encyclopedia/AB2C3D2_tP32_127_g_eh_gk_k

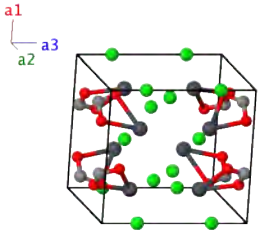


Prototype	:	$CCl_2O_3Pb_2$
AFLOW prototype label	:	AB2C3D2_tP32_127_g_eh_gk_k
***Strukturbericht* designation**	:	None
Pearson symbol	:	tP32
Space group number	:	127
Space group symbol	:	$P4/mbm$
AFLOW prototype command	:	`aflow --proto=AB2C3D2_tP32_127_g_eh_gk_k --params=`$a, c/a, z_1, x_2, x_3, x_4, x_5, z_5, x_6, z_6$

- (Onotaro, 1934) made the first determination of the structure of phosgenite, which was given *Strukturbericht* designation $G7_5$ by (Gottfried). (Giuseppetti, 1974) and others found that the actual structure has a unit cell twice that of $G7_5$. This newer structure is presented here.

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$z_1\,\mathbf{a}_3$	=	$z_1 c\,\hat{\mathbf{z}}$	($4e$)	Cl I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	($4e$)	Cl I
$\mathbf{B}_3$	=	$-z_1\,\mathbf{a}_3$	=	$-z_1 c\,\hat{\mathbf{z}}$	($4e$)	Cl I
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	($4e$)	Cl I
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_2$	=	$x_2 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{y}}$	($4g$)	C
$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_2$	=	$-x_2 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{y}}$	($4g$)	C
$\mathbf{B}_7$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + x_2\,\mathbf{a}_2$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + x_2 a\,\hat{\mathbf{y}}$	($4g$)	C
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 - x_2\,\mathbf{a}_2$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} - x_2 a\,\hat{\mathbf{y}}$	($4g$)	C
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2$	=	$x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}}$	($4g$)	O I
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2$	=	$-x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}}$	($4g$)	O I
$\mathbf{B}_{11}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + x_3\,\mathbf{a}_2$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}}$	($4g$)	O I
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 - x_3\,\mathbf{a}_2$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}}$	($4g$)	O I
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + x_4\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4h$)	Cl II
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4h$)	Cl II
$\mathbf{B}_{15}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4h$)	Cl II
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4h$)	Cl II
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + \left(\frac{1}{2} + x_5\right)\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	($8k$)	O II
$\mathbf{B}_{18}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2} - x_5\right)\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	($8k$)	O II
$\mathbf{B}_{19}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	($8k$)	O II
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 - x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	($8k$)	O II
$\mathbf{B}_{21}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	($8k$)	O II
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 - x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	($8k$)	O II
$\mathbf{B}_{23}$	=	$x_5\,\mathbf{a}_1 + \left(\frac{1}{2} + x_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	($8k$)	O II
$\mathbf{B}_{24}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2} - x_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	($8k$)	O II
$\mathbf{B}_{25}$	=	$x_6\,\mathbf{a}_1 + \left(\frac{1}{2} + x_6\right)\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	($8k$)	Pb
$\mathbf{B}_{26}$	=	$-x_6\,\mathbf{a}_1 + \left(\frac{1}{2} - x_6\right)\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	($8k$)	Pb
$\mathbf{B}_{27}$	=	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 + x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	($8k$)	Pb
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 - x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	($8k$)	Pb
$\mathbf{B}_{29}$	=	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 + x_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	($8k$)	Pb
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 - x_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	($8k$)	Pb
$\mathbf{B}_{31}$	=	$x_6\,\mathbf{a}_1 + \left(\frac{1}{2} + x_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	($8k$)	Pb
$\mathbf{B}_{32}$	=	$-x_6\,\mathbf{a}_1 + \left(\frac{1}{2} - x_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	($8k$)	Pb

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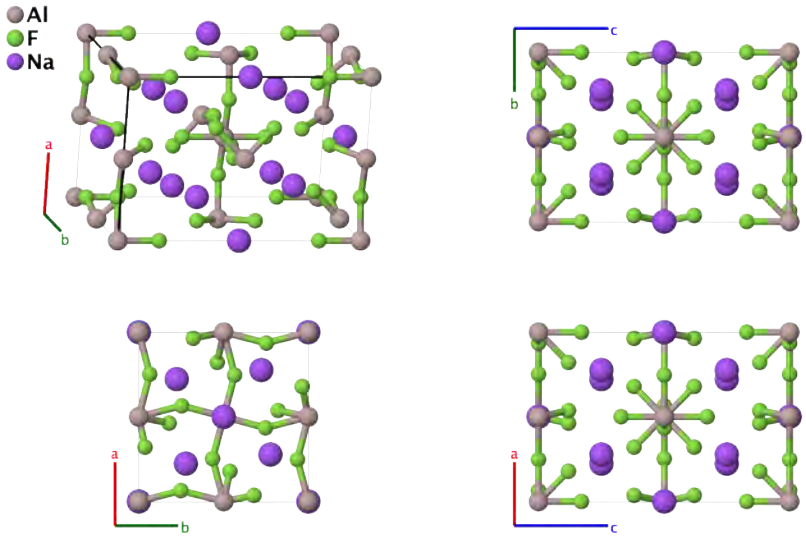
- C. Gottfried and F. Schossberger, eds., *Strukturbericht Band III 1933-1935* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:
- CIF: pp. 1695
- POSCAR: pp. 1696

Chiolite ($Na_5Al_3F_{14}$, $K7_5$) Structure: A3B14C5_tP44_128_ac_ehi_bg

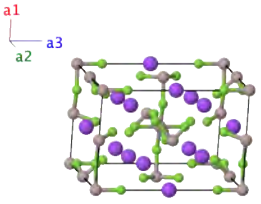
http://aflow.org/prototype-encyclopedia/A3B14C5_tP44_128_ac_ehi_bg



Prototype	:	$Al_3F_{14}Na_5$
AFLOW prototype label	:	A3B14C5_tP44_128_ac_ehi_bg
***Strukturbericht* designation**	:	$K7_5$
Pearson symbol	:	tP44
Space group number	:	128
Space group symbol	:	$P4/mnc$
AFLOW prototype command	:	`aflow --proto=A3B14C5_tP44_128_ac_ehi_bg --params=`$a, c/a, z_4, x_5, x_6, y_6, x_7, y_7, z_7$

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($2a$)	Al I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($2a$)	Al I
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	($2b$)	Na I
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$	($2b$)	Na I

$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{y}}$	(4*c*)	Al II
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	(4*c*)	Al II
$\mathbf{B}_7$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*c*)	Al II
$\mathbf{B}_8$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*c*)	Al II
$\mathbf{B}_9$	=	$z_4\,\mathbf{a}_3$	=	$z_4 c\,\hat{\mathbf{z}}$	(4*e*)	F I
$\mathbf{B}_{10}$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(4*e*)	F I
$\mathbf{B}_{11}$	=	$-z_4\,\mathbf{a}_3$	=	$-z_4 c\,\hat{\mathbf{z}}$	(4*e*)	F I
$\mathbf{B}_{12}$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(4*e*)	F I
$\mathbf{B}_{13}$	=	$x_5\,\mathbf{a}_1 + \left(\frac{1}{2} + x_5\right)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(8*g*)	Na II
$\mathbf{B}_{14}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2} - x_5\right)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(8*g*)	Na II
$\mathbf{B}_{15}$	=	$\left(\frac{1}{2} - x_5\right)\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(8*g*)	Na II
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_5\right)\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(8*g*)	Na II
$\mathbf{B}_{17}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2} - x_5\right)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(8*g*)	Na II
$\mathbf{B}_{18}$	=	$x_5\,\mathbf{a}_1 + \left(\frac{1}{2} + x_5\right)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(8*g*)	Na II
$\mathbf{B}_{19}$	=	$\left(\frac{1}{2} + x_5\right)\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(8*g*)	Na II
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} - x_5\right)\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(8*g*)	Na II
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2$	=	$x_6 a\,\hat{\mathbf{x}} + y_6 a\,\hat{\mathbf{y}}$	(8*h*)	F II
$\mathbf{B}_{22}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2$	=	$-x_6 a\,\hat{\mathbf{x}} - y_6 a\,\hat{\mathbf{y}}$	(8*h*)	F II
$\mathbf{B}_{23}$	=	$-y_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2$	=	$-y_6 a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}}$	(8*h*)	F II
$\mathbf{B}_{24}$	=	$y_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2$	=	$y_6 a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}}$	(8*h*)	F II
$\mathbf{B}_{25}$	=	$\left(\frac{1}{2} - x_6\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right)a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(8*h*)	F II
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} + x_6\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right)a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(8*h*)	F II
$\mathbf{B}_{27}$	=	$\left(\frac{1}{2} + y_6\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + x_6\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(8*h*)	F II
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} - y_6\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - x_6\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(8*h*)	F II
$\mathbf{B}_{29}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + y_7 a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(16*i*)	F III
$\mathbf{B}_{30}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} - y_7 a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(16*i*)	F III
$\mathbf{B}_{31}$	=	$-y_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$-y_7 a\,\hat{\mathbf{x}} + x_7 a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(16*i*)	F III
$\mathbf{B}_{32}$	=	$y_7\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$y_7 a\,\hat{\mathbf{x}} - x_7 a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(16*i*)	F III
$\mathbf{B}_{33}$	=	$\left(\frac{1}{2} - x_7\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right)c\,\hat{\mathbf{z}}$	(16*i*)	F III
$\mathbf{B}_{34}$	=	$\left(\frac{1}{2} + x_7\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - y_7\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_7\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right)c\,\hat{\mathbf{z}}$	(16*i*)	F III
$\mathbf{B}_{35}$	=	$\left(\frac{1}{2} + y_7\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + x_7\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_7\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right)c\,\hat{\mathbf{z}}$	(16*i*)	F III
$\mathbf{B}_{36}$	=	$\left(\frac{1}{2} - y_7\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - x_7\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_7\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right)c\,\hat{\mathbf{z}}$	(16*i*)	F III
$\mathbf{B}_{37}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} - y_7 a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	(16*i*)	F III

$\mathbf{B}_{38}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} + y_7 a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	(16*i*)	F III
$\mathbf{B}_{39}$	=	$y_7\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$y_7 a\,\hat{\mathbf{x}} - x_7 a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	(16*i*)	F III
$\mathbf{B}_{40}$	=	$-y_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$-y_7 a\,\hat{\mathbf{x}} + x_7 a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	(16*i*)	F III
$\mathbf{B}_{41}$	=	$\left(\frac{1}{2} + x_7\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_7\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_7\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right)c\,\hat{\mathbf{z}}$	(16*i*)	F III
$\mathbf{B}_{42}$	=	$\left(\frac{1}{2} - x_7\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_7\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_7\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right)c\,\hat{\mathbf{z}}$	(16*i*)	F III
$\mathbf{B}_{43}$	=	$\left(\frac{1}{2} - y_7\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_7\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_7\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right)c\,\hat{\mathbf{z}}$	(16*i*)	F III
$\mathbf{B}_{44}$	=	$\left(\frac{1}{2} + y_7\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_7\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_7\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right)c\,\hat{\mathbf{z}}$	(16*i*)	F III

References:

- C. Jacoboni, A. Leble, and J. J. Rousseau, *Détermination précise de la structure de la chiolite $Na_5Al_3F_{14}$ et étude par R.P.E. de $Na_5Al_3F_{14}$:Cr^{3+}*, J. Solid State Chem. **36**, 297–304 (1981), doi:10.1016/0022-4596(81)90440-0.

Found in:

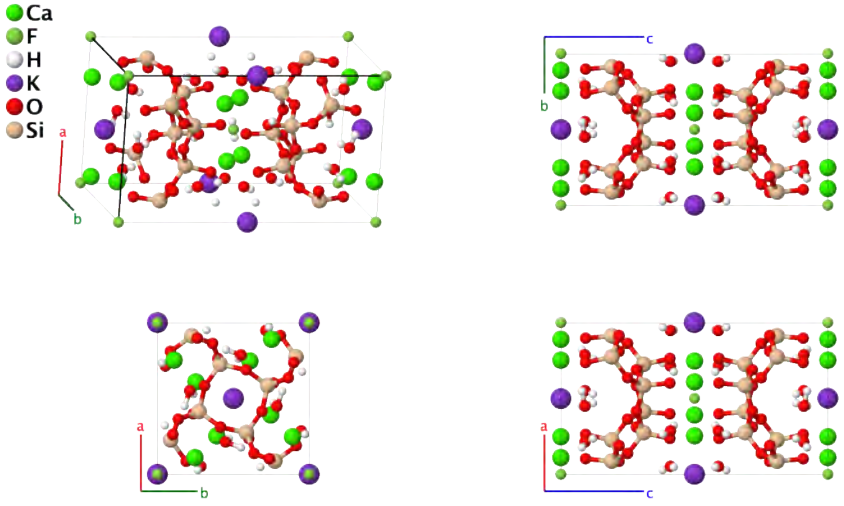
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1696
- POSCAR: pp. 1696

Apophyllite ($KCa_4Si_8O_{20}F{\cdot}8H_2O$, $S5_2$) Structure: A4BC16DE28F8_tP116_128_h_a_2i_b_g3i_i

http://aflow.org/prototype-encyclopedia/A4BC16DE28F8_tP116_128_h_a_2i_b_g3i_i

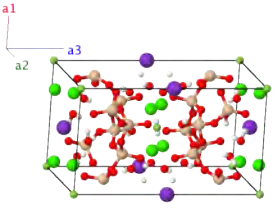


Prototype	:	$Ca_4FH_{16}KO_{28}Si_8$
AFLOW prototype label	:	A4BC16DE28F8_tP116_128_h_a_2i_b_g3i_i
***Strukturbericht* designation**	:	$S5_2$
Pearson symbol	:	tP116
Space group number	:	128
Space group symbol	:	$P4/mnc$
AFLOW prototype command	:	`aflow --proto=A4BC16DE28F8_tP116_128_h_a_2i_b_g3i_i --params=`$a, c/a, x_3, x_4, y_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}$

- Although we use the structure found by (Chao, 1971), we should note that there is some disagreement between Chao's x-ray diffraction data and the neutron diffraction data taken by (Prince, 1971): while both agree on the positions of the heavy atoms, Prince's work suggests that some of the hydrogens may form OH radicals rather than water molecules.
- In any case, the fluorine atoms are usually partially replaced by OH radicals. This sample, which is predominantly fluorine, is technically labeled apophyllite-(KF).

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(2*a*)	F
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(2*a*)	F
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	(2*b*)	K
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$	(2*b*)	K
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(8*g*)	O I
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(8*g*)	O I
$\mathbf{B}_7$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(8*g*)	O I
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(8*g*)	O I
$\mathbf{B}_9$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(8*g*)	O I
$\mathbf{B}_{10}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(8*g*)	O I
$\mathbf{B}_{11}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(8*g*)	O I
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(8*g*)	O I
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2$	=	$x_4 a\,\hat{\mathbf{x}} + y_4 a\,\hat{\mathbf{y}}$	(8*h*)	Ca
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2$	=	$-x_4 a\,\hat{\mathbf{x}} - y_4 a\,\hat{\mathbf{y}}$	(8*h*)	Ca
$\mathbf{B}_{15}$	=	$-y_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2$	=	$-y_4 a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}}$	(8*h*)	Ca
$\mathbf{B}_{16}$	=	$y_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2$	=	$y_4 a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}}$	(8*h*)	Ca
$\mathbf{B}_{17}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(8*h*)	Ca
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(8*h*)	Ca
$\mathbf{B}_{19}$	=	$\left(\frac{1}{2} + y_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_4\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(8*h*)	Ca
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} - y_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(8*h*)	Ca
$\mathbf{B}_{21}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(16*i*)	H I
$\mathbf{B}_{22}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(16*i*)	H I
$\mathbf{B}_{23}$	=	$-y_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$-y_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(16*i*)	H I
$\mathbf{B}_{24}$	=	$y_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$y_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(16*i*)	H I
$\mathbf{B}_{25}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	(16*i*)	H I
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_5\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	(16*i*)	H I
$\mathbf{B}_{27}$	=	$\left(\frac{1}{2} + y_5\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_5\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	(16*i*)	H I
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} - y_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_5\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	(16*i*)	H I
$\mathbf{B}_{29}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - y_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(16*i*)	H I
$\mathbf{B}_{30}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(16*i*)	H I
$\mathbf{B}_{31}$	=	$y_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$y_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(16*i*)	H I
$\mathbf{B}_{32}$	=	$-y_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-y_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(16*i*)	H I

$$\mathbf{B}_{33} = \left(\tfrac{1}{2}+x_5\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_5\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_5\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_5\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_5\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{H I}$$

$$\mathbf{B}_{34} = \left(\tfrac{1}{2}-x_5\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_5\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_5\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_5\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_5\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{H I}$$

$$\mathbf{B}_{35} = \left(\tfrac{1}{2}-y_5\right)\mathbf{a}_1+\left(\tfrac{1}{2}-x_5\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-y_5\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_5\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_5\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{H I}$$

$$\mathbf{B}_{36} = \left(\tfrac{1}{2}+y_5\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_5\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+y_5\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_5\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_5\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{H I}$$

$$\mathbf{B}_{37} = x_6\,\mathbf{a}_1+y_6\,\mathbf{a}_2+z_6\,\mathbf{a}_3 = x_6a\,\hat{\mathbf{x}}+y_6a\,\hat{\mathbf{y}}+z_6c\,\hat{\mathbf{z}} \quad (16i) \quad \text{H II}$$

$$\mathbf{B}_{38} = -x_6\,\mathbf{a}_1-y_6\,\mathbf{a}_2+z_6\,\mathbf{a}_3 = -x_6a\,\hat{\mathbf{x}}-y_6a\,\hat{\mathbf{y}}+z_6c\,\hat{\mathbf{z}} \quad (16i) \quad \text{H II}$$

$$\mathbf{B}_{39} = -y_6\,\mathbf{a}_1+x_6\,\mathbf{a}_2+z_6\,\mathbf{a}_3 = -y_6a\,\hat{\mathbf{x}}+x_6a\,\hat{\mathbf{y}}+z_6c\,\hat{\mathbf{z}} \quad (16i) \quad \text{H II}$$

$$\mathbf{B}_{40} = y_6\,\mathbf{a}_1-x_6\,\mathbf{a}_2+z_6\,\mathbf{a}_3 = y_6a\,\hat{\mathbf{x}}-x_6a\,\hat{\mathbf{y}}+z_6c\,\hat{\mathbf{z}} \quad (16i) \quad \text{H II}$$

$$\mathbf{B}_{41} = \left(\tfrac{1}{2}-x_6\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_6\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_6\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_6\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{H II}$$

$$\mathbf{B}_{42} = \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_6\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_6\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_6\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{H II}$$

$$\mathbf{B}_{43} = \left(\tfrac{1}{2}+y_6\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_6\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+y_6\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_6\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{H II}$$

$$\mathbf{B}_{44} = \left(\tfrac{1}{2}-y_6\right)\mathbf{a}_1+\left(\tfrac{1}{2}-x_6\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-y_6\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_6\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{H II}$$

$$\mathbf{B}_{45} = -x_6\,\mathbf{a}_1-y_6\,\mathbf{a}_2-z_6\,\mathbf{a}_3 = -x_6a\,\hat{\mathbf{x}}-y_6a\,\hat{\mathbf{y}}-z_6c\,\hat{\mathbf{z}} \quad (16i) \quad \text{H II}$$

$$\mathbf{B}_{46} = x_6\,\mathbf{a}_1+y_6\,\mathbf{a}_2-z_6\,\mathbf{a}_3 = x_6a\,\hat{\mathbf{x}}+y_6a\,\hat{\mathbf{y}}-z_6c\,\hat{\mathbf{z}} \quad (16i) \quad \text{H II}$$

$$\mathbf{B}_{47} = y_6\,\mathbf{a}_1-x_6\,\mathbf{a}_2-z_6\,\mathbf{a}_3 = y_6a\,\hat{\mathbf{x}}-x_6a\,\hat{\mathbf{y}}-z_6c\,\hat{\mathbf{z}} \quad (16i) \quad \text{H II}$$

$$\mathbf{B}_{48} = -y_6\,\mathbf{a}_1+x_6\,\mathbf{a}_2-z_6\,\mathbf{a}_3 = -y_6a\,\hat{\mathbf{x}}+x_6a\,\hat{\mathbf{y}}-z_6c\,\hat{\mathbf{z}} \quad (16i) \quad \text{H II}$$

$$\mathbf{B}_{49} = \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_6\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_6\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_6\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{H II}$$

$$\mathbf{B}_{50} = \left(\tfrac{1}{2}-x_6\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_6\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_6\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_6\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{H II}$$

$$\mathbf{B}_{51} = \left(\tfrac{1}{2}-y_6\right)\mathbf{a}_1+\left(\tfrac{1}{2}-x_6\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-y_6\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_6\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{H II}$$

$$\mathbf{B}_{52} = \left(\tfrac{1}{2}+y_6\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_6\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+y_6\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_6\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{H II}$$

$$\mathbf{B}_{53} = x_7\,\mathbf{a}_1+y_7\,\mathbf{a}_2+z_7\,\mathbf{a}_3 = x_7a\,\hat{\mathbf{x}}+y_7a\,\hat{\mathbf{y}}+z_7c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O II}$$

$$\mathbf{B}_{54} = -x_7\,\mathbf{a}_1-y_7\,\mathbf{a}_2+z_7\,\mathbf{a}_3 = -x_7a\,\hat{\mathbf{x}}-y_7a\,\hat{\mathbf{y}}+z_7c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O II}$$

$$\mathbf{B}_{55} = -y_7\,\mathbf{a}_1+x_7\,\mathbf{a}_2+z_7\,\mathbf{a}_3 = -y_7a\,\hat{\mathbf{x}}+x_7a\,\hat{\mathbf{y}}+z_7c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O II}$$

$$\mathbf{B}_{56} = y_7\,\mathbf{a}_1-x_7\,\mathbf{a}_2+z_7\,\mathbf{a}_3 = y_7a\,\hat{\mathbf{x}}-x_7a\,\hat{\mathbf{y}}+z_7c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O II}$$

$$\mathbf{B}_{57} = \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_7\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_7\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O II}$$

$$\mathbf{B}_{58} = \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_7\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O II}$$

$$\mathbf{B}_{59} = \left(\tfrac{1}{2}+y_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+y_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_7\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O II}$$

$$\mathbf{B}_{60} = \left(\tfrac{1}{2}-y_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-y_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_7\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O II}$$

$$\mathbf{B}_{61} = -x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3 = -x_7a\,\hat{\mathbf{x}} - y_7a\,\hat{\mathbf{y}} - z_7c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O II}$$

$$\mathbf{B}_{62} = x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3 = x_7a\,\hat{\mathbf{x}} + y_7a\,\hat{\mathbf{y}} - z_7c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O II}$$

$$\mathbf{B}_{63} = y_7\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3 = y_7a\,\hat{\mathbf{x}} - x_7a\,\hat{\mathbf{y}} - z_7c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O II}$$

$$\mathbf{B}_{64} = -y_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3 = -y_7a\,\hat{\mathbf{x}} + x_7a\,\hat{\mathbf{y}} - z_7c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O II}$$

$$\mathbf{B}_{65} = \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_7\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O II}$$

$$\mathbf{B}_{66} = \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_7\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O II}$$

$$\mathbf{B}_{67} = \left(\tfrac{1}{2}-y_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-y_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_7\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O II}$$

$$\mathbf{B}_{68} = \left(\tfrac{1}{2}+y_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+y_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_7\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O II}$$

$$\mathbf{B}_{69} = x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 = x_8a\,\hat{\mathbf{x}} + y_8a\,\hat{\mathbf{y}} + z_8c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O III}$$

$$\mathbf{B}_{70} = -x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 = -x_8a\,\hat{\mathbf{x}} - y_8a\,\hat{\mathbf{y}} + z_8c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O III}$$

$$\mathbf{B}_{71} = -y_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 = -y_8a\,\hat{\mathbf{x}} + x_8a\,\hat{\mathbf{y}} + z_8c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O III}$$

$$\mathbf{B}_{72} = y_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 = y_8a\,\hat{\mathbf{x}} - x_8a\,\hat{\mathbf{y}} + z_8c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O III}$$

$$\mathbf{B}_{73} = \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_8\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_8\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O III}$$

$$\mathbf{B}_{74} = \left(\tfrac{1}{2}+x_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_8\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_8\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O III}$$

$$\mathbf{B}_{75} = \left(\tfrac{1}{2}+y_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+x_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+y_8\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_8\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O III}$$

$$\mathbf{B}_{76} = \left(\tfrac{1}{2}-y_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-y_8\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_8\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O III}$$

$$\mathbf{B}_{77} = -x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3 = -x_8a\,\hat{\mathbf{x}} - y_8a\,\hat{\mathbf{y}} - z_8c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O III}$$

$$\mathbf{B}_{78} = x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3 = x_8a\,\hat{\mathbf{x}} + y_8a\,\hat{\mathbf{y}} - z_8c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O III}$$

$$\mathbf{B}_{79} = y_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3 = y_8a\,\hat{\mathbf{x}} - x_8a\,\hat{\mathbf{y}} - z_8c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O III}$$

$$\mathbf{B}_{80} = -y_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3 = -y_8a\,\hat{\mathbf{x}} + x_8a\,\hat{\mathbf{y}} - z_8c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O III}$$

$$\mathbf{B}_{81} = \left(\tfrac{1}{2}+x_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_8\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_8\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O III}$$

$$\mathbf{B}_{82} = \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_8\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_8\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O III}$$

$$\mathbf{B}_{83} = \left(\tfrac{1}{2}-y_8\right)\mathbf{a}_1+\left(\tfrac{1}{2}-x_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-y_8\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_8\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O III}$$

$$\mathbf{B}_{84} = \left(\tfrac{1}{2}+y_8\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+y_8\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_8\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O III}$$

$$\mathbf{B}_{85} = x_9\,\mathbf{a}_1+y_9\,\mathbf{a}_2+z_9\,\mathbf{a}_3 = x_9a\,\hat{\mathbf{x}}+y_9a\,\hat{\mathbf{y}}+z_9c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O IV}$$

$$\mathbf{B}_{86} = -x_9\,\mathbf{a}_1-y_9\,\mathbf{a}_2+z_9\,\mathbf{a}_3 = -x_9a\,\hat{\mathbf{x}}-y_9a\,\hat{\mathbf{y}}+z_9c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O IV}$$

$$\mathbf{B}_{87} = -y_9\,\mathbf{a}_1+x_9\,\mathbf{a}_2+z_9\,\mathbf{a}_3 = -y_9a\,\hat{\mathbf{x}}+x_9a\,\hat{\mathbf{y}}+z_9c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O IV}$$

$$\mathbf{B}_{88} = y_9\,\mathbf{a}_1-x_9\,\mathbf{a}_2+z_9\,\mathbf{a}_3 = y_9a\,\hat{\mathbf{x}}-x_9a\,\hat{\mathbf{y}}+z_9c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O IV}$$

$$\mathbf{B}_{89} = \left(\tfrac{1}{2}-x_9\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_9\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_9\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_9\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_9\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O IV}$$

$$\mathbf{B}_{90} = \left(\tfrac{1}{2}+x_9\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_9\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_9\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_9\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_9\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O IV}$$

$$\mathbf{B}_{91} = \left(\tfrac{1}{2}+y_9\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_9\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+y_9\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_9\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_9\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O IV}$$

$$\mathbf{B}_{92} = \left(\tfrac{1}{2}-y_9\right)\mathbf{a}_1+\left(\tfrac{1}{2}-x_9\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-y_9\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_9\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_9\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O IV}$$

$$\mathbf{B}_{93} = -x_9\,\mathbf{a}_1-y_9\,\mathbf{a}_2-z_9\,\mathbf{a}_3 = -x_9a\,\hat{\mathbf{x}}-y_9a\,\hat{\mathbf{y}}-z_9c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O IV}$$

$$\mathbf{B}_{94} = x_9\,\mathbf{a}_1+y_9\,\mathbf{a}_2-z_9\,\mathbf{a}_3 = x_9a\,\hat{\mathbf{x}}+y_9a\,\hat{\mathbf{y}}-z_9c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O IV}$$

$$\mathbf{B}_{95} = y_9\,\mathbf{a}_1-x_9\,\mathbf{a}_2-z_9\,\mathbf{a}_3 = y_9a\,\hat{\mathbf{x}}-x_9a\,\hat{\mathbf{y}}-z_9c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O IV}$$

$$\mathbf{B}_{96} = -y_9\,\mathbf{a}_1+x_9\,\mathbf{a}_2-z_9\,\mathbf{a}_3 = -y_9a\,\hat{\mathbf{x}}+x_9a\,\hat{\mathbf{y}}-z_9c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O IV}$$

$$\mathbf{B}_{97} = \left(\tfrac{1}{2}+x_9\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_9\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_9\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_9\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_9\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O IV}$$

$$\mathbf{B}_{98} = \left(\tfrac{1}{2}-x_9\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_9\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_9\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_9\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_9\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O IV}$$

$$\mathbf{B}_{99} = \left(\tfrac{1}{2}-y_9\right)\mathbf{a}_1+\left(\tfrac{1}{2}-x_9\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-y_9\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_9\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_9\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O IV}$$

$$\mathbf{B}_{100} = \left(\tfrac{1}{2}+y_9\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_9\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+y_9\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_9\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_9\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{O IV}$$

$$\mathbf{B}_{101} = x_{10}\,\mathbf{a}_1+y_{10}\,\mathbf{a}_2+z_{10}\,\mathbf{a}_3 = x_{10}a\,\hat{\mathbf{x}}+y_{10}a\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}} \quad (16i) \quad \text{Si}$$

$$\mathbf{B}_{102} = -x_{10}\,\mathbf{a}_1-y_{10}\,\mathbf{a}_2+z_{10}\,\mathbf{a}_3 = -x_{10}a\,\hat{\mathbf{x}}-y_{10}a\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}} \quad (16i) \quad \text{Si}$$

$$\mathbf{B}_{103} = -y_{10}\,\mathbf{a}_1+x_{10}\,\mathbf{a}_2+z_{10}\,\mathbf{a}_3 = -y_{10}a\,\hat{\mathbf{x}}+x_{10}a\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}} \quad (16i) \quad \text{Si}$$

$$\mathbf{B}_{104} = y_{10}\,\mathbf{a}_1-x_{10}\,\mathbf{a}_2+z_{10}\,\mathbf{a}_3 = y_{10}a\,\hat{\mathbf{x}}-x_{10}a\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}} \quad (16i) \quad \text{Si}$$

$$\mathbf{B}_{105} = \left(\tfrac{1}{2}-x_{10}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{10}\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{10}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{10}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{10}\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{Si}$$

$$\mathbf{B}_{106} = \left(\tfrac{1}{2}+x_{10}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{10}\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{10}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{10}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{10}\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{Si}$$

$$\mathbf{B}_{107} = \left(\tfrac{1}{2}+y_{10}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_{10}\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+y_{10}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_{10}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{10}\right)c\,\hat{\mathbf{z}} \quad (16i) \quad \text{Si}$$

$\mathbf{B}_{108}$	=	$\left(\frac{1}{2}-y_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}-x_{10}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{10}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{10}\right)c\,\hat{\mathbf{z}}$	(16*i*)	Si
$\mathbf{B}_{109}$	=	$-x_{10}\,\mathbf{a}_1-y_{10}\,\mathbf{a}_2-z_{10}\,\mathbf{a}_3$	=	$-x_{10}a\,\hat{\mathbf{x}}-y_{10}a\,\hat{\mathbf{y}}-z_{10}c\,\hat{\mathbf{z}}$	(16*i*)	Si
$\mathbf{B}_{110}$	=	$x_{10}\,\mathbf{a}_1+y_{10}\,\mathbf{a}_2-z_{10}\,\mathbf{a}_3$	=	$x_{10}a\,\hat{\mathbf{x}}+y_{10}a\,\hat{\mathbf{y}}-z_{10}c\,\hat{\mathbf{z}}$	(16*i*)	Si
$\mathbf{B}_{111}$	=	$y_{10}\,\mathbf{a}_1-x_{10}\,\mathbf{a}_2-z_{10}\,\mathbf{a}_3$	=	$y_{10}a\,\hat{\mathbf{x}}-x_{10}a\,\hat{\mathbf{y}}-z_{10}c\,\hat{\mathbf{z}}$	(16*i*)	Si
$\mathbf{B}_{112}$	=	$-y_{10}\,\mathbf{a}_1+x_{10}\,\mathbf{a}_2-z_{10}\,\mathbf{a}_3$	=	$-y_{10}a\,\hat{\mathbf{x}}+x_{10}a\,\hat{\mathbf{y}}-z_{10}c\,\hat{\mathbf{z}}$	(16*i*)	Si
$\mathbf{B}_{113}$	=	$\left(\frac{1}{2}+x_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{10}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{10}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{10}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{10}\right)c\,\hat{\mathbf{z}}$	(16*i*)	Si
$\mathbf{B}_{114}$	=	$\left(\frac{1}{2}-x_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{10}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{10}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{10}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{10}\right)c\,\hat{\mathbf{z}}$	(16*i*)	Si
$\mathbf{B}_{115}$	=	$\left(\frac{1}{2}-y_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}-x_{10}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{10}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{10}\right)c\,\hat{\mathbf{z}}$	(16*i*)	Si
$\mathbf{B}_{116}$	=	$\left(\frac{1}{2}+y_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}+x_{10}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{10}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{10}\right)c\,\hat{\mathbf{z}}$	(16*i*)	Si

References:

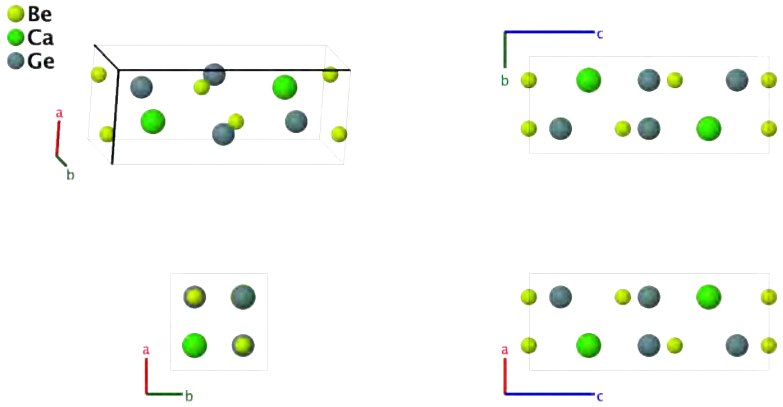
- G. Y. Chao, *The refinement of the crystal structure of apophyllite: II. Determination of the hydrogen positions by X-ray diffraction*, Am. Mineral. **56**, 1234–1242 (1971).
- E. Prince, *The refinement of the crystal structure of apophyllite: III. Determination of the hydrogen positions by neutron diffraction*, Am. Mineral. **56**, 1243–1251 (1971).

Geometry files:

- CIF: pp. 1696
- POSCAR: pp. 1697

$CaBe_2Ge_2$ Structure: A2BC2_tP10_129_ac_c_bc

http://aflow.org/prototype-encyclopedia/A2BC2_tP10_129_ac_c_bc



Prototype	:	Be_2CaGe_2
AFLOW prototype label	:	A2BC2_tP10_129_ac_c_bc
***Strukturbericht* designation**	:	None
Pearson symbol	:	tP10
Space group number	:	129
Space group symbol	:	$P4/nmm$
AFLOW prototype command	:	`aflow --proto=A2BC2_tP10_129_ac_c_bc --params=`$a, c/a, z_3, z_4, z_5$

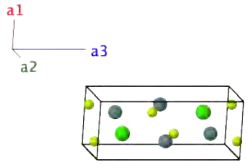
Other compounds with this structure

- $BaMg_2Pb_2$, $BaPd_2Sb_2$, $BaZn_2Sn_2$, $EuAu_2Al_2$, $EuPd_2Sb_2$, $EuPt_2Ge_2$, $LaPt_2Bi_2$, $LaPt_2Ge_2$, $LiPd_2Bi_2$, $SrPd_2Sb_2$, $SrPt_2As_2$, and $ThIr_2Si_2$

- This is a ternary form of the $D1_3$ ($BaAl_4$) structure. The atomic positions are approximately the same as in the conventional cell of $BaAl_4$, but the distribution of the atoms on those sites and the resulting relaxation leads to a different structure.
- Space group $P4/nmm$ #129 has two settings, but both have the same origin of the z-axis, so either setting will do here. We chose our standard setting 2 here.

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}}$	$(2a)$	Be I

$\mathbf{B}_2$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}}$	(2*a*)	Be I
$\mathbf{B}_3$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(2*b*)	Ge I
$\mathbf{B}_4$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(2*b*)	Ge I
$\mathbf{B}_5$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	(2*c*)	Be II
$\mathbf{B}_6$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	(2*c*)	Be II
$\mathbf{B}_7$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	(2*c*)	Ca
$\mathbf{B}_8$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	(2*c*)	Ca
$\mathbf{B}_9$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + z_5c\,\hat{\mathbf{z}}$	(2*c*)	Ge II
$\mathbf{B}_{10}$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} - z_5c\,\hat{\mathbf{z}}$	(2*c*)	Ge II

References:

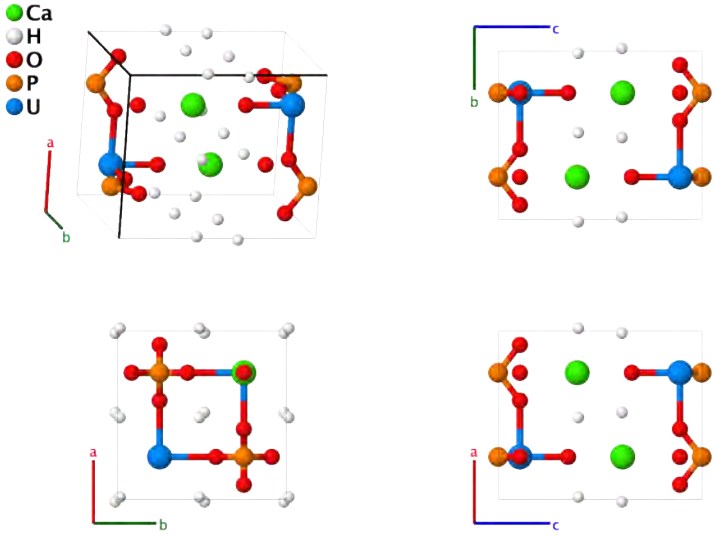
- B. Eisenmann, N. May, W. Müller, and H. Schäfer, *Eine neue strukturelle Variante des $BaAl_4$-Typs: Der $CaBe_2Ge_2$-Typ*, Z. Naturforsch. B **27**, 1155–1157 (1972), doi:10.1515/znb-1972-1008.

Geometry files:

- CIF: pp. 1698
- POSCAR: pp. 1698

Meta-autunite (I) [$Ca(UO_2)_2(PO_4)_2{\cdot}6H_2O$, $H5_{10}$] Structure: AB4C6DE_tP26_129_c_j_2ci_a_c

http://aflow.org/prototype-encyclopedia/AB4C6DE_tP26_129_c_j_2ci_a_c

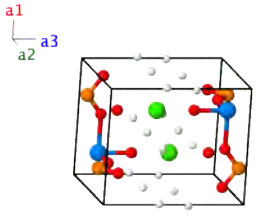


Prototype	:	$Ca(H_2O)_6O_{12}P_2U_2$
AFLOW prototype label	:	AB4C6DE_tP26_129_c_j_2ci_a_c
***Strukturbericht* designation**	:	$H5_{10}$
Pearson symbol	:	tP26
Space group number	:	129
Space group symbol	:	$P4/nmm$
AFLOW prototype command	:	`aflow --proto=AB4C6DE_tP26_129_c_j_2ci_a_c --params=`$a, c/a, z_2, z_3, z_4, z_5, y_6, z_6, x_7, z_7$

- Autunite $Ca(UO_2)_2(PO_4)_2{\cdot}nH_2O$, is found in three varieties: naturally occurring autunite, with $n \gtrsim 10$, and meta-autunite (I), which is partially dehydrated, $6 \gtrsim n \gtrsim 10$. Further dehydration in the laboratory produces meta-autunite (II).
- (Beintema, 1938) proposed a structure for meta-autunite (I), which (Herrmann, 1941) designated $H5_{10}$. He did not locate the calcium and oxygen atoms nor the water molecules. This structure was improved by (Makarov, 1960), and we include it here as our prototype for $H5_{10}$. The Ca-I site is 50% occupied, while the H_2O site is 75% occupied.
- The AFLOW label models the structure as if the sites were fully occupied.

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}}$	(2*a*)	P
$\mathbf{B}_2$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}}$	(2*a*)	P
$\mathbf{B}_3$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + z_2c\,\hat{\mathbf{z}}$	(2*c*)	Ca
$\mathbf{B}_4$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} - z_2c\,\hat{\mathbf{z}}$	(2*c*)	Ca
$\mathbf{B}_5$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	(2*c*)	O I
$\mathbf{B}_6$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	(2*c*)	O I
$\mathbf{B}_7$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	(2*c*)	O II
$\mathbf{B}_8$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	(2*c*)	O II
$\mathbf{B}_9$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + z_5c\,\hat{\mathbf{z}}$	(2*c*)	U
$\mathbf{B}_{10}$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} - z_5c\,\hat{\mathbf{z}}$	(2*c*)	U
$\mathbf{B}_{11}$	=	$\frac{1}{4}\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + y_6a\,\hat{\mathbf{y}} + z_6c\,\hat{\mathbf{z}}$	(8*i*)	O III
$\mathbf{B}_{12}$	=	$\frac{1}{4}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right)a\,\hat{\mathbf{y}} + z_6c\,\hat{\mathbf{z}}$	(8*i*)	O III
$\mathbf{B}_{13}$	=	$\left(\frac{1}{2} - y_6\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_6\right)a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + z_6c\,\hat{\mathbf{z}}$	(8*i*)	O III
$\mathbf{B}_{14}$	=	$y_6\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$y_6a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + z_6c\,\hat{\mathbf{z}}$	(8*i*)	O III
$\mathbf{B}_{15}$	=	$\frac{3}{4}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right)a\,\hat{\mathbf{y}} - z_6c\,\hat{\mathbf{z}}$	(8*i*)	O III
$\mathbf{B}_{16}$	=	$\frac{3}{4}\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} - y_6a\,\hat{\mathbf{y}} - z_6c\,\hat{\mathbf{z}}$	(8*i*)	O III
$\mathbf{B}_{17}$	=	$\left(\frac{1}{2} + y_6\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_6\right)a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} - z_6c\,\hat{\mathbf{z}}$	(8*i*)	O III
$\mathbf{B}_{18}$	=	$-y_6\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$-y_6a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} - z_6c\,\hat{\mathbf{z}}$	(8*i*)	O III
$\mathbf{B}_{19}$	=	$x_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$x_7a\,\hat{\mathbf{x}} + x_7a\,\hat{\mathbf{y}} + z_7c\,\hat{\mathbf{z}}$	(8*j*)	H_2O
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} - x_7\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_7\right)\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_7\right)a\,\hat{\mathbf{y}} + z_7c\,\hat{\mathbf{z}}$	(8*j*)	H_2O
$\mathbf{B}_{21}$	=	$\left(\frac{1}{2} - x_7\right)\mathbf{a}_1 + x_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_7\right)a\,\hat{\mathbf{x}} + x_7a\,\hat{\mathbf{y}} + z_7c\,\hat{\mathbf{z}}$	(8*j*)	H_2O
$\mathbf{B}_{22}$	=	$x_7\,\mathbf{a}_1 + \left(\frac{1}{2} - x_7\right)\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$x_7a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_7\right)a\,\hat{\mathbf{y}} + z_7c\,\hat{\mathbf{z}}$	(8*j*)	H_2O
$\mathbf{B}_{23}$	=	$-x_7\,\mathbf{a}_1 + \left(\frac{1}{2} + x_7\right)\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$-x_7a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_7\right)a\,\hat{\mathbf{y}} - z_7c\,\hat{\mathbf{z}}$	(8*j*)	H_2O
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} + x_7\right)\mathbf{a}_1 - x_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_7\right)a\,\hat{\mathbf{x}} - x_7a\,\hat{\mathbf{y}} - z_7c\,\hat{\mathbf{z}}$	(8*j*)	H_2O
$\mathbf{B}_{25}$	=	$\left(\frac{1}{2} + x_7\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_7\right)\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_7\right)a\,\hat{\mathbf{y}} - z_7c\,\hat{\mathbf{z}}$	(8*j*)	H_2O
$\mathbf{B}_{26}$	=	$-x_7\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$-x_7a\,\hat{\mathbf{x}} - x_7a\,\hat{\mathbf{y}} - z_7c\,\hat{\mathbf{z}}$	(8*j*)	H_2O

References:

- Y. S. Makarov and V. I. Ivanov, *The crystal structure of meta-autunite, $Ca(UO_2)_2(PO_4)_2$*6H_2O*, Doklady Akademii Nauk SSSR **132**, 601–603 (1960).
- J. Beintema, *On the composition and the crystallography of autunite and the meta-autunites*, Rec. Trav. Chim. Pays-Bas **57**, 155–175 (1938), doi:10.1002/recl.19380570206.
- K. Herrmann, ed., *Strukturbericht Band VI 1938* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1941).

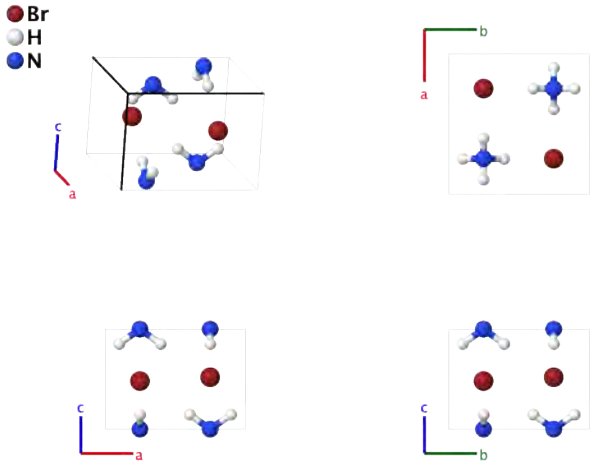
Found in:

- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:
- CIF: pp. 1698
- POSCAR: pp. 1698

NH_4Br (*B*25) Structure: AB4C_tP12_129_c_i_a

http://aflow.org/prototype-encyclopedia/AB4C_tP12_129_c_i_a



Prototype	:	BrH_4N
AFLOW prototype label	:	AB4C_tP12_129_c_i_a
***Strukturbericht* designation**	:	*B*25
Pearson symbol	:	tP12
Space group number	:	129
Space group symbol	:	$P4/nmm$
AFLOW prototype command	:	`aflow --proto=AB4C_tP12_129_c_i_a --params=`$a, c/a, z_2, y_3, z_3$

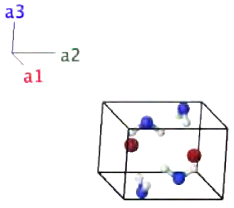
Other compounds with this structure

- NH_4I

- Data was taken at 100 °C using deuterium.
- The atomic positions were given for setting 1 of space group #129. We have shifted this to setting 2, placing the origin at the inversion site of the structure.

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$\mathbf{B}_1$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}}$	(2*a*)	N
$\mathbf{B}_2$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}}$	(2*a*)	N
$\mathbf{B}_3$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + z_2c\,\hat{\mathbf{z}}$	(2*c*)	Br
$\mathbf{B}_4$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} - z_2c\,\hat{\mathbf{z}}$	(2*c*)	Br
$\mathbf{B}_5$	=	$\frac{1}{4}\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + y_3a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	(8*i*)	H
$\mathbf{B}_6$	=	$\frac{1}{4}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	(8*i*)	H
$\mathbf{B}_7$	=	$\left(\frac{1}{2} - y_3\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	(8*i*)	H
$\mathbf{B}_8$	=	$y_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$y_3a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	(8*i*)	H
$\mathbf{B}_9$	=	$\frac{3}{4}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	(8*i*)	H
$\mathbf{B}_{10}$	=	$\frac{3}{4}\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} - y_3a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	(8*i*)	H
$\mathbf{B}_{11}$	=	$\left(\frac{1}{2} + y_3\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	(8*i*)	H
$\mathbf{B}_{12}$	=	$-y_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-y_3a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	(8*i*)	H

References:

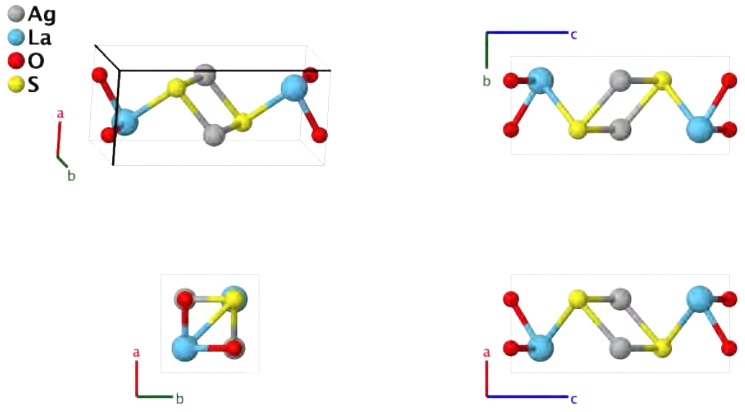
- H. A. Levy and S. W. Peterson, *Neutron Diffraction Determination of the Crystal Structure of Ammonium Bromide in Four Phases*, J. Am. Chem. Soc. **75**, 1536–1542 (1953), doi:10.1021/ja01103a006.

Geometry files:

- CIF: pp. 1699
- POSCAR: pp. 1699

LaOAgS Structure: ABCD_tP8_129_b_c_a_c

http://aflow.org/prototype-encyclopedia/ABCD_tP8_129_b_c_a_c



Prototype	:	AgLaOS
AFLOW prototype label	:	ABCD_tP8_129_b_c_a_c
***Strukturbericht* designation**	:	None
Pearson symbol	:	tP8
Space group number	:	129
Space group symbol	:	$P4/nmm$
AFLOW prototype command	:	`aflow --proto=ABCD_tP8_129_b_c_a_c --params=`$a, c/a, z_3, z_4$

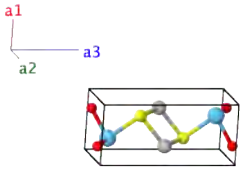
Other compounds with this structure

- LaAgSeO and LaAgTeO

- This structure has the same space group and Wyckoff positions as AsCuSiZr, but there is a substantial difference in the c/a ratio leading to different bonding between the layers.

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{3}{4}\mathbf{a}_1 + \frac{1}{4}\mathbf{a}_2$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}}$	$(2a)$	O
$\mathbf{B}_2$	=	$\frac{1}{4}\mathbf{a}_1 + \frac{3}{4}\mathbf{a}_2$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}}$	$(2a)$	O
$\mathbf{B}_3$	=	$\frac{3}{4}\mathbf{a}_1 + \frac{1}{4}\mathbf{a}_2 + \frac{1}{2}\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(2b)$	Ag
$\mathbf{B}_4$	=	$\frac{1}{4}\mathbf{a}_1 + \frac{3}{4}\mathbf{a}_2 + \frac{1}{2}\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(2b)$	Ag

$\mathbf{B}_5$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(2*c*)	La
$\mathbf{B}_6$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(2*c*)	La
$\mathbf{B}_7$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(2*c*)	S
$\mathbf{B}_8$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(2*c*)	S

References:

- M. Palazzi and S. Jaulmes, *Structure du Conducteur Ionique (LaO)AgS*, Acta Crystallogr. Sect. B Struct. Sci. **37**, 1337–1339 (1981), doi:10.1107/S0567740881005876.

Found in:

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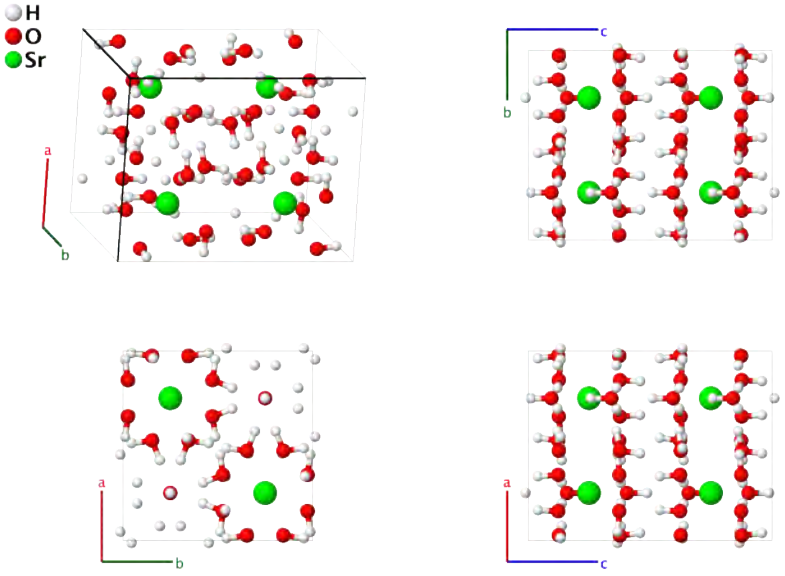
Geometry files:

- CIF: pp. 1699
- POSCAR: pp. 1699

$Sr(OH)_2(H_2O)_8$ Structure: A18B10C_tP116_130_2c4g_2c2g_a

http://aflow.org/prototype-encyclopedia/A18B10C_tP116_130_2c4g_2c2g_a

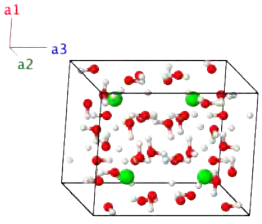


Prototype	:	$H_{18}O_{10}Sr$
AFLOW prototype label	:	A18B10C_tP116_130_2c4g_2c2g_a
***Strukturbericht* designation**	:	None
Pearson symbol	:	tP116
Space group number	:	130
Space group symbol	:	$P4/ncc$
AFLOW prototype command	:	`aflow --proto=A18B10C_tP116_130_2c4g_2c2g_a --params=`$a, c/a, z_2, z_3, z_4, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}$

- This determination of the crystal structure of $Sr(OH)_2(H_2O)_8$ improves upon the structure found by (Natta, 1928), which was given the *Strukturbericht* designation $E6_1$ by (Hermann, 1937). The new structure quadruples the size of the unit cell, locates the hydrogen atoms, and changes the space group.
- We use the data taken by (Ricci, 2005) at 20 K.

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*a*)	Sr
$\mathbf{B}_2$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*a*)	Sr
$\mathbf{B}_3$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(4*a*)	Sr
$\mathbf{B}_4$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(4*a*)	Sr
$\mathbf{B}_5$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + z_2c\,\hat{\mathbf{z}}$	(4*c*)	H I
$\mathbf{B}_6$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	(4*c*)	H I
$\mathbf{B}_7$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} - z_2c\,\hat{\mathbf{z}}$	(4*c*)	H I
$\mathbf{B}_8$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	(4*c*)	H I
$\mathbf{B}_9$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	(4*c*)	H II
$\mathbf{B}_{10}$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(4*c*)	H II
$\mathbf{B}_{11}$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	(4*c*)	H II
$\mathbf{B}_{12}$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(4*c*)	H II
$\mathbf{B}_{13}$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_{14}$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_{15}$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_{16}$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(4*c*)	O I
$\mathbf{B}_{17}$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + z_5c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_{18}$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_{19}$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} - z_5c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_{20}$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	(4*c*)	O II
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6a\,\hat{\mathbf{x}} + y_6a\,\hat{\mathbf{y}} + z_6c\,\hat{\mathbf{z}}$	(16*g*)	H III
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right)a\,\hat{\mathbf{y}} + z_6c\,\hat{\mathbf{z}}$	(16*g*)	H III
$\mathbf{B}_{23}$	=	$\left(\frac{1}{2} - y_6\right)\mathbf{a}_1 + x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_6\right)a\,\hat{\mathbf{x}} + x_6a\,\hat{\mathbf{y}} + z_6c\,\hat{\mathbf{z}}$	(16*g*)	H III
$\mathbf{B}_{24}$	=	$y_6\,\mathbf{a}_1 + \left(\frac{1}{2} - x_6\right)\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$y_6a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{y}} + z_6c\,\hat{\mathbf{z}}$	(16*g*)	H III
$\mathbf{B}_{25}$	=	$-x_6\,\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$-x_6a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\,\hat{\mathbf{z}}$	(16*g*)	H III
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{x}} - y_6a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\,\hat{\mathbf{z}}$	(16*g*)	H III
$\mathbf{B}_{27}$	=	$\left(\frac{1}{2} + y_6\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_6\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\,\hat{\mathbf{z}}$	(16*g*)	H III
$\mathbf{B}_{28}$	=	$-y_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$-y_6a\,\hat{\mathbf{x}} - x_6a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\,\hat{\mathbf{z}}$	(16*g*)	H III
$\mathbf{B}_{29}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$-x_6a\,\hat{\mathbf{x}} - y_6a\,\hat{\mathbf{y}} - z_6c\,\hat{\mathbf{z}}$	(16*g*)	H III
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right)a\,\hat{\mathbf{y}} - z_6c\,\hat{\mathbf{z}}$	(16*g*)	H III
$\mathbf{B}_{31}$	=	$\left(\frac{1}{2} + y_6\right)\mathbf{a}_1 - x_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_6\right)a\,\hat{\mathbf{x}} - x_6a\,\hat{\mathbf{y}} - z_6c\,\hat{\mathbf{z}}$	(16*g*)	H III
$\mathbf{B}_{32}$	=	$-y_6\,\mathbf{a}_1 + \left(\frac{1}{2} + x_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$-y_6a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{y}} - z_6c\,\hat{\mathbf{z}}$	(16*g*)	H III
$\mathbf{B}_{33}$	=	$x_6\,\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$x_6a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\,\hat{\mathbf{z}}$	(16*g*)	H III
$\mathbf{B}_{34}$	=	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 + y_6\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{x}} + y_6a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\,\hat{\mathbf{z}}$	(16*g*)	H III

$\mathbf{B}_{35}$	=	$\left(\frac{1}{2}-y_6\right)\mathbf{a}_1+\left(\frac{1}{2}-x_6\right)\mathbf{a}_2+\left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	(16*g*)	H III
$\mathbf{B}_{36}$	=	$y_6\,\mathbf{a}_1+x_6\,\mathbf{a}_2+\left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	=	$y_6a\,\hat{\mathbf{x}}+x_6a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	(16*g*)	H III
$\mathbf{B}_{37}$	=	$x_7\,\mathbf{a}_1+y_7\,\mathbf{a}_2+z_7\,\mathbf{a}_3$	=	$x_7a\,\hat{\mathbf{x}}+y_7a\,\hat{\mathbf{y}}+z_7c\,\hat{\mathbf{z}}$	(16*g*)	H IV
$\mathbf{B}_{38}$	=	$\left(\frac{1}{2}-x_7\right)\mathbf{a}_1+\left(\frac{1}{2}-y_7\right)\mathbf{a}_2+z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_7\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_7\right)a\,\hat{\mathbf{y}}+z_7c\,\hat{\mathbf{z}}$	(16*g*)	H IV
$\mathbf{B}_{39}$	=	$\left(\frac{1}{2}-y_7\right)\mathbf{a}_1+x_7\,\mathbf{a}_2+z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_7\right)a\,\hat{\mathbf{x}}+x_7a\,\hat{\mathbf{y}}+z_7c\,\hat{\mathbf{z}}$	(16*g*)	H IV
$\mathbf{B}_{40}$	=	$y_7\,\mathbf{a}_1+\left(\frac{1}{2}-x_7\right)\mathbf{a}_2+z_7\,\mathbf{a}_3$	=	$y_7a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_7\right)a\,\hat{\mathbf{y}}+z_7c\,\hat{\mathbf{z}}$	(16*g*)	H IV
$\mathbf{B}_{41}$	=	$-x_7\,\mathbf{a}_1+\left(\frac{1}{2}+y_7\right)\mathbf{a}_2+\left(\frac{1}{2}-z_7\right)\mathbf{a}_3$	=	$-x_7a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_7\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_7\right)c\,\hat{\mathbf{z}}$	(16*g*)	H IV
$\mathbf{B}_{42}$	=	$\left(\frac{1}{2}+x_7\right)\mathbf{a}_1-y_7\,\mathbf{a}_2+\left(\frac{1}{2}-z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{x}}-y_7a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_7\right)c\,\hat{\mathbf{z}}$	(16*g*)	H IV
$\mathbf{B}_{43}$	=	$\left(\frac{1}{2}+y_7\right)\mathbf{a}_1+\left(\frac{1}{2}+x_7\right)\mathbf{a}_2+\left(\frac{1}{2}-z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_7\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_7\right)c\,\hat{\mathbf{z}}$	(16*g*)	H IV
$\mathbf{B}_{44}$	=	$-y_7\,\mathbf{a}_1-x_7\,\mathbf{a}_2+\left(\frac{1}{2}-z_7\right)\mathbf{a}_3$	=	$-y_7a\,\hat{\mathbf{x}}-x_7a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_7\right)c\,\hat{\mathbf{z}}$	(16*g*)	H IV
$\mathbf{B}_{45}$	=	$-x_7\,\mathbf{a}_1-y_7\,\mathbf{a}_2-z_7\,\mathbf{a}_3$	=	$-x_7a\,\hat{\mathbf{x}}-y_7a\,\hat{\mathbf{y}}-z_7c\,\hat{\mathbf{z}}$	(16*g*)	H IV
$\mathbf{B}_{46}$	=	$\left(\frac{1}{2}+x_7\right)\mathbf{a}_1+\left(\frac{1}{2}+y_7\right)\mathbf{a}_2-z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_7\right)a\,\hat{\mathbf{y}}-z_7c\,\hat{\mathbf{z}}$	(16*g*)	H IV
$\mathbf{B}_{47}$	=	$\left(\frac{1}{2}+y_7\right)\mathbf{a}_1-x_7\,\mathbf{a}_2-z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_7\right)a\,\hat{\mathbf{x}}-x_7a\,\hat{\mathbf{y}}-z_7c\,\hat{\mathbf{z}}$	(16*g*)	H IV
$\mathbf{B}_{48}$	=	$-y_7\,\mathbf{a}_1+\left(\frac{1}{2}+x_7\right)\mathbf{a}_2-z_7\,\mathbf{a}_3$	=	$-y_7a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{y}}-z_7c\,\hat{\mathbf{z}}$	(16*g*)	H IV
$\mathbf{B}_{49}$	=	$x_7\,\mathbf{a}_1+\left(\frac{1}{2}-y_7\right)\mathbf{a}_2+\left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	=	$x_7a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_7\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	(16*g*)	H IV
$\mathbf{B}_{50}$	=	$\left(\frac{1}{2}-x_7\right)\mathbf{a}_1+y_7\,\mathbf{a}_2+\left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_7\right)a\,\hat{\mathbf{x}}+y_7a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	(16*g*)	H IV
$\mathbf{B}_{51}$	=	$\left(\frac{1}{2}-y_7\right)\mathbf{a}_1+\left(\frac{1}{2}-x_7\right)\mathbf{a}_2+\left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_7\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_7\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	(16*g*)	H IV
$\mathbf{B}_{52}$	=	$y_7\,\mathbf{a}_1+x_7\,\mathbf{a}_2+\left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	=	$y_7a\,\hat{\mathbf{x}}+x_7a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	(16*g*)	H IV
$\mathbf{B}_{53}$	=	$x_8\,\mathbf{a}_1+y_8\,\mathbf{a}_2+z_8\,\mathbf{a}_3$	=	$x_8a\,\hat{\mathbf{x}}+y_8a\,\hat{\mathbf{y}}+z_8c\,\hat{\mathbf{z}}$	(16*g*)	H V
$\mathbf{B}_{54}$	=	$\left(\frac{1}{2}-x_8\right)\mathbf{a}_1+\left(\frac{1}{2}-y_8\right)\mathbf{a}_2+z_8\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_8\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_8\right)a\,\hat{\mathbf{y}}+z_8c\,\hat{\mathbf{z}}$	(16*g*)	H V
$\mathbf{B}_{55}$	=	$\left(\frac{1}{2}-y_8\right)\mathbf{a}_1+x_8\,\mathbf{a}_2+z_8\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_8\right)a\,\hat{\mathbf{x}}+x_8a\,\hat{\mathbf{y}}+z_8c\,\hat{\mathbf{z}}$	(16*g*)	H V
$\mathbf{B}_{56}$	=	$y_8\,\mathbf{a}_1+\left(\frac{1}{2}-x_8\right)\mathbf{a}_2+z_8\,\mathbf{a}_3$	=	$y_8a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_8\right)a\,\hat{\mathbf{y}}+z_8c\,\hat{\mathbf{z}}$	(16*g*)	H V
$\mathbf{B}_{57}$	=	$-x_8\,\mathbf{a}_1+\left(\frac{1}{2}+y_8\right)\mathbf{a}_2+\left(\frac{1}{2}-z_8\right)\mathbf{a}_3$	=	$-x_8a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_8\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_8\right)c\,\hat{\mathbf{z}}$	(16*g*)	H V
$\mathbf{B}_{58}$	=	$\left(\frac{1}{2}+x_8\right)\mathbf{a}_1-y_8\,\mathbf{a}_2+\left(\frac{1}{2}-z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_8\right)a\,\hat{\mathbf{x}}-y_8a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_8\right)c\,\hat{\mathbf{z}}$	(16*g*)	H V
$\mathbf{B}_{59}$	=	$\left(\frac{1}{2}+y_8\right)\mathbf{a}_1+\left(\frac{1}{2}+x_8\right)\mathbf{a}_2+\left(\frac{1}{2}-z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_8\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_8\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_8\right)c\,\hat{\mathbf{z}}$	(16*g*)	H V
$\mathbf{B}_{60}$	=	$-y_8\,\mathbf{a}_1-x_8\,\mathbf{a}_2+\left(\frac{1}{2}-z_8\right)\mathbf{a}_3$	=	$-y_8a\,\hat{\mathbf{x}}-x_8a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_8\right)c\,\hat{\mathbf{z}}$	(16*g*)	H V
$\mathbf{B}_{61}$	=	$-x_8\,\mathbf{a}_1-y_8\,\mathbf{a}_2-z_8\,\mathbf{a}_3$	=	$-x_8a\,\hat{\mathbf{x}}-y_8a\,\hat{\mathbf{y}}-z_8c\,\hat{\mathbf{z}}$	(16*g*)	H V
$\mathbf{B}_{62}$	=	$\left(\frac{1}{2}+x_8\right)\mathbf{a}_1+\left(\frac{1}{2}+y_8\right)\mathbf{a}_2-z_8\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_8\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_8\right)a\,\hat{\mathbf{y}}-z_8c\,\hat{\mathbf{z}}$	(16*g*)	H V
$\mathbf{B}_{63}$	=	$\left(\frac{1}{2}+y_8\right)\mathbf{a}_1-x_8\,\mathbf{a}_2-z_8\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_8\right)a\,\hat{\mathbf{x}}-x_8a\,\hat{\mathbf{y}}-z_8c\,\hat{\mathbf{z}}$	(16*g*)	H V
$\mathbf{B}_{64}$	=	$-y_8\,\mathbf{a}_1+\left(\frac{1}{2}+x_8\right)\mathbf{a}_2-z_8\,\mathbf{a}_3$	=	$-y_8a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_8\right)a\,\hat{\mathbf{y}}-z_8c\,\hat{\mathbf{z}}$	(16*g*)	H V
$\mathbf{B}_{65}$	=	$x_8\,\mathbf{a}_1+\left(\frac{1}{2}-y_8\right)\mathbf{a}_2+\left(\frac{1}{2}+z_8\right)\mathbf{a}_3$	=	$x_8a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_8\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_8\right)c\,\hat{\mathbf{z}}$	(16*g*)	H V
$\mathbf{B}_{66}$	=	$\left(\frac{1}{2}-x_8\right)\mathbf{a}_1+y_8\,\mathbf{a}_2+\left(\frac{1}{2}+z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_8\right)a\,\hat{\mathbf{x}}+y_8a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_8\right)c\,\hat{\mathbf{z}}$	(16*g*)	H V

$$
\begin{array}{lcccl}
\mathbf{B}_{67} & = & \left(\frac{1}{2}-y_8\right)\mathbf{a}_1+\left(\frac{1}{2}-x_8\right)\mathbf{a}_2+\left(\frac{1}{2}+z_8\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-y_8\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_8\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_8\right)c\,\hat{\mathbf{z}} \quad (16g) \quad \text{H V} \\
\mathbf{B}_{68} & = & y_8\,\mathbf{a}_1+x_8\,\mathbf{a}_2+\left(\frac{1}{2}+z_8\right)\mathbf{a}_3 & = & y_8a\,\hat{\mathbf{x}}+x_8a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_8\right)c\,\hat{\mathbf{z}} \quad (16g) \quad \text{H V} \\
\mathbf{B}_{69} & = & x_9\,\mathbf{a}_1+y_9\,\mathbf{a}_2+z_9\,\mathbf{a}_3 & = & x_9a\,\hat{\mathbf{x}}+y_9a\,\hat{\mathbf{y}}+z_9c\,\hat{\mathbf{z}} \quad (16g) \quad \text{H VI} \\
\mathbf{B}_{70} & = & \left(\frac{1}{2}-x_9\right)\mathbf{a}_1+\left(\frac{1}{2}-y_9\right)\mathbf{a}_2+z_9\,\mathbf{a}_3 & = & \left(\frac{1}{2}-x_9\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_9\right)a\,\hat{\mathbf{y}}+z_9c\,\hat{\mathbf{z}} \quad (16g) \quad \text{H VI} \\
\mathbf{B}_{71} & = & \left(\frac{1}{2}-y_9\right)\mathbf{a}_1+x_9\,\mathbf{a}_2+z_9\,\mathbf{a}_3 & = & \left(\frac{1}{2}-y_9\right)a\,\hat{\mathbf{x}}+x_9a\,\hat{\mathbf{y}}+z_9c\,\hat{\mathbf{z}} \quad (16g) \quad \text{H VI} \\
\mathbf{B}_{72} & = & y_9\,\mathbf{a}_1+\left(\frac{1}{2}-x_9\right)\mathbf{a}_2+z_9\,\mathbf{a}_3 & = & y_9a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_9\right)a\,\hat{\mathbf{y}}+z_9c\,\hat{\mathbf{z}} \quad (16g) \quad \text{H VI} \\
\mathbf{B}_{73} & = & -x_9\,\mathbf{a}_1+\left(\frac{1}{2}+y_9\right)\mathbf{a}_2+\left(\frac{1}{2}-z_9\right)\mathbf{a}_3 & = & -x_9a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_9\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_9\right)c\,\hat{\mathbf{z}} \quad (16g) \quad \text{H VI} \\
\mathbf{B}_{74} & = & \left(\frac{1}{2}+x_9\right)\mathbf{a}_1-y_9\,\mathbf{a}_2+\left(\frac{1}{2}-z_9\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_9\right)a\,\hat{\mathbf{x}}-y_9a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_9\right)c\,\hat{\mathbf{z}} \quad (16g) \quad \text{H VI} \\
\mathbf{B}_{75} & = & \left(\frac{1}{2}+y_9\right)\mathbf{a}_1+\left(\frac{1}{2}+x_9\right)\mathbf{a}_2+\left(\frac{1}{2}-z_9\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+y_9\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_9\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_9\right)c\,\hat{\mathbf{z}} \quad (16g) \quad \text{H VI} \\
\mathbf{B}_{76} & = & -y_9\,\mathbf{a}_1-x_9\,\mathbf{a}_2+\left(\frac{1}{2}-z_9\right)\mathbf{a}_3 & = & -y_9a\,\hat{\mathbf{x}}-x_9a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_9\right)c\,\hat{\mathbf{z}} \quad (16g) \quad \text{H VI} \\
\mathbf{B}_{77} & = & -x_9\,\mathbf{a}_1-y_9\,\mathbf{a}_2-z_9\,\mathbf{a}_3 & = & -x_9a\,\hat{\mathbf{x}}-y_9a\,\hat{\mathbf{y}}-z_9c\,\hat{\mathbf{z}} \quad (16g) \quad \text{H VI} \\
\mathbf{B}_{78} & = & \left(\frac{1}{2}+x_9\right)\mathbf{a}_1+\left(\frac{1}{2}+y_9\right)\mathbf{a}_2-z_9\,\mathbf{a}_3 & = & \left(\frac{1}{2}+x_9\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_9\right)a\,\hat{\mathbf{y}}-z_9c\,\hat{\mathbf{z}} \quad (16g) \quad \text{H VI} \\
\mathbf{B}_{79} & = & \left(\frac{1}{2}+y_9\right)\mathbf{a}_1-x_9\,\mathbf{a}_2-z_9\,\mathbf{a}_3 & = & \left(\frac{1}{2}+y_9\right)a\,\hat{\mathbf{x}}-x_9a\,\hat{\mathbf{y}}-z_9c\,\hat{\mathbf{z}} \quad (16g) \quad \text{H VI} \\
\mathbf{B}_{80} & = & -y_9\,\mathbf{a}_1+\left(\frac{1}{2}+x_9\right)\mathbf{a}_2-z_9\,\mathbf{a}_3 & = & -y_9a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_9\right)a\,\hat{\mathbf{y}}-z_9c\,\hat{\mathbf{z}} \quad (16g) \quad \text{H VI} \\
\mathbf{B}_{81} & = & x_9\,\mathbf{a}_1+\left(\frac{1}{2}-y_9\right)\mathbf{a}_2+\left(\frac{1}{2}+z_9\right)\mathbf{a}_3 & = & x_9a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_9\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_9\right)c\,\hat{\mathbf{z}} \quad (16g) \quad \text{H VI} \\
\mathbf{B}_{82} & = & \left(\frac{1}{2}-x_9\right)\mathbf{a}_1+y_9\,\mathbf{a}_2+\left(\frac{1}{2}+z_9\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_9\right)a\,\hat{\mathbf{x}}+y_9a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_9\right)c\,\hat{\mathbf{z}} \quad (16g) \quad \text{H VI} \\
\mathbf{B}_{83} & = & \left(\frac{1}{2}-y_9\right)\mathbf{a}_1+\left(\frac{1}{2}-x_9\right)\mathbf{a}_2+\left(\frac{1}{2}+z_9\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-y_9\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_9\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_9\right)c\,\hat{\mathbf{z}} \quad (16g) \quad \text{H VI} \\
\mathbf{B}_{84} & = & y_9\,\mathbf{a}_1+x_9\,\mathbf{a}_2+\left(\frac{1}{2}+z_9\right)\mathbf{a}_3 & = & y_9a\,\hat{\mathbf{x}}+x_9a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_9\right)c\,\hat{\mathbf{z}} \quad (16g) \quad \text{H VI} \\
\mathbf{B}_{85} & = & x_{10}\,\mathbf{a}_1+y_{10}\,\mathbf{a}_2+z_{10}\,\mathbf{a}_3 & = & x_{10}a\,\hat{\mathbf{x}}+y_{10}a\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}} \quad (16g) \quad \text{O III} \\
\mathbf{B}_{86} & = & \left(\frac{1}{2}-x_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{10}\right)\mathbf{a}_2+z_{10}\,\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{10}\right)a\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}} \quad (16g) \quad \text{O III} \\
\mathbf{B}_{87} & = & \left(\frac{1}{2}-y_{10}\right)\mathbf{a}_1+x_{10}\,\mathbf{a}_2+z_{10}\,\mathbf{a}_3 & = & \left(\frac{1}{2}-y_{10}\right)a\,\hat{\mathbf{x}}+x_{10}a\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}} \quad (16g) \quad \text{O III} \\
\mathbf{B}_{88} & = & y_{10}\,\mathbf{a}_1+\left(\frac{1}{2}-x_{10}\right)\mathbf{a}_2+z_{10}\,\mathbf{a}_3 & = & y_{10}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}} \quad (16g) \quad \text{O III} \\
\mathbf{B}_{89} & = & -x_{10}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{10}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{10}\right)\mathbf{a}_3 & = & -x_{10}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{10}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{10}\right)c\,\hat{\mathbf{z}} \quad (16g) \quad \text{O III} \\
\mathbf{B}_{90} & = & \left(\frac{1}{2}+x_{10}\right)\mathbf{a}_1-y_{10}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{10}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{x}}-y_{10}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{10}\right)c\,\hat{\mathbf{z}} \quad (16g) \quad \text{O III} \\
\mathbf{B}_{91} & = & \left(\frac{1}{2}+y_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}+x_{10}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{10}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+y_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{10}\right)c\,\hat{\mathbf{z}} \quad (16g) \quad \text{O III} \\
\mathbf{B}_{92} & = & -y_{10}\,\mathbf{a}_1-x_{10}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{10}\right)\mathbf{a}_3 & = & -y_{10}a\,\hat{\mathbf{x}}-x_{10}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{10}\right)c\,\hat{\mathbf{z}} \quad (16g) \quad \text{O III} \\
\mathbf{B}_{93} & = & -x_{10}\,\mathbf{a}_1-y_{10}\,\mathbf{a}_2-z_{10}\,\mathbf{a}_3 & = & -x_{10}a\,\hat{\mathbf{x}}-y_{10}a\,\hat{\mathbf{y}}-z_{10}c\,\hat{\mathbf{z}} \quad (16g) \quad \text{O III} \\
\mathbf{B}_{94} & = & \left(\frac{1}{2}+x_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{10}\right)\mathbf{a}_2-z_{10}\,\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{10}\right)a\,\hat{\mathbf{y}}-z_{10}c\,\hat{\mathbf{z}} \quad (16g) \quad \text{O III} \\
\mathbf{B}_{95} & = & \left(\frac{1}{2}+y_{10}\right)\mathbf{a}_1-x_{10}\,\mathbf{a}_2-z_{10}\,\mathbf{a}_3 & = & \left(\frac{1}{2}+y_{10}\right)a\,\hat{\mathbf{x}}-x_{10}a\,\hat{\mathbf{y}}-z_{10}c\,\hat{\mathbf{z}} \quad (16g) \quad \text{O III} \\
\mathbf{B}_{96} & = & -y_{10}\,\mathbf{a}_1+\left(\frac{1}{2}+x_{10}\right)\mathbf{a}_2-z_{10}\,\mathbf{a}_3 & = & -y_{10}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{y}}-z_{10}c\,\hat{\mathbf{z}} \quad (16g) \quad \text{O III} \\
\mathbf{B}_{97} & = & x_{10}\,\mathbf{a}_1+\left(\frac{1}{2}-y_{10}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{10}\right)\mathbf{a}_3 & = & x_{10}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{10}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{10}\right)c\,\hat{\mathbf{z}} \quad (16g) \quad \text{O III} \\
\mathbf{B}_{98} & = & \left(\frac{1}{2}-x_{10}\right)\mathbf{a}_1+y_{10}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{10}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{x}}+y_{10}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{10}\right)c\,\hat{\mathbf{z}} \quad (16g) \quad \text{O III}
\end{array}
$$

$$\mathbf{B}_{99} = \left(\tfrac{1}{2}-y_{10}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-x_{10}\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-y_{10}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_{10}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{10}\right)c\,\hat{\mathbf{z}} \quad (16g) \quad \text{O III}$$

$$\mathbf{B}_{100} = y_{10}\,\mathbf{a}_1+x_{10}\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_{10}\right)\mathbf{a}_3 = y_{10}a\,\hat{\mathbf{x}}+x_{10}a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{10}\right)c\,\hat{\mathbf{z}} \quad (16g) \quad \text{O III}$$

$$\mathbf{B}_{101} = x_{11}\,\mathbf{a}_1+y_{11}\,\mathbf{a}_2+z_{11}\,\mathbf{a}_3 = x_{11}a\,\hat{\mathbf{x}}+y_{11}a\,\hat{\mathbf{y}}+z_{11}c\,\hat{\mathbf{z}} \quad (16g) \quad \text{O IV}$$

$$\mathbf{B}_{102} = \left(\tfrac{1}{2}-x_{11}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{11}\right)\mathbf{a}_2+z_{11}\,\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{11}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{11}\right)a\,\hat{\mathbf{y}}+z_{11}c\,\hat{\mathbf{z}} \quad (16g) \quad \text{O IV}$$

$$\mathbf{B}_{103} = \left(\tfrac{1}{2}-y_{11}\right)\mathbf{a}_1+x_{11}\,\mathbf{a}_2+z_{11}\,\mathbf{a}_3 = \left(\tfrac{1}{2}-y_{11}\right)a\,\hat{\mathbf{x}}+x_{11}a\,\hat{\mathbf{y}}+z_{11}c\,\hat{\mathbf{z}} \quad (16g) \quad \text{O IV}$$

$$\mathbf{B}_{104} = y_{11}\,\mathbf{a}_1+\left(\tfrac{1}{2}-x_{11}\right)\mathbf{a}_2+z_{11}\,\mathbf{a}_3 = y_{11}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_{11}\right)a\,\hat{\mathbf{y}}+z_{11}c\,\hat{\mathbf{z}} \quad (16g) \quad \text{O IV}$$

$$\mathbf{B}_{105} = -x_{11}\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_{11}\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_{11}\right)\mathbf{a}_3 = -x_{11}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{11}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{11}\right)c\,\hat{\mathbf{z}} \quad (16g) \quad \text{O IV}$$

$$\mathbf{B}_{106} = \left(\tfrac{1}{2}+x_{11}\right)\mathbf{a}_1-y_{11}\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{11}\right)a\,\hat{\mathbf{x}}-y_{11}a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{11}\right)c\,\hat{\mathbf{z}} \quad (16g) \quad \text{O IV}$$

$$\mathbf{B}_{107} = \left(\tfrac{1}{2}+y_{11}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_{11}\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+y_{11}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_{11}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{11}\right)c\,\hat{\mathbf{z}} \quad (16g) \quad \text{O IV}$$

$$\mathbf{B}_{108} = -y_{11}\,\mathbf{a}_1-x_{11}\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_{11}\right)\mathbf{a}_3 = -y_{11}a\,\hat{\mathbf{x}}-x_{11}a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{11}\right)c\,\hat{\mathbf{z}} \quad (16g) \quad \text{O IV}$$

$$\mathbf{B}_{109} = -x_{11}\,\mathbf{a}_1-y_{11}\,\mathbf{a}_2-z_{11}\,\mathbf{a}_3 = -x_{11}a\,\hat{\mathbf{x}}-y_{11}a\,\hat{\mathbf{y}}-z_{11}c\,\hat{\mathbf{z}} \quad (16g) \quad \text{O IV}$$

$$\mathbf{B}_{110} = \left(\tfrac{1}{2}+x_{11}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{11}\right)\mathbf{a}_2-z_{11}\,\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{11}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{11}\right)a\,\hat{\mathbf{y}}-z_{11}c\,\hat{\mathbf{z}} \quad (16g) \quad \text{O IV}$$

$$\mathbf{B}_{111} = \left(\tfrac{1}{2}+y_{11}\right)\mathbf{a}_1-x_{11}\,\mathbf{a}_2-z_{11}\,\mathbf{a}_3 = \left(\tfrac{1}{2}+y_{11}\right)a\,\hat{\mathbf{x}}-x_{11}a\,\hat{\mathbf{y}}-z_{11}c\,\hat{\mathbf{z}} \quad (16g) \quad \text{O IV}$$

$$\mathbf{B}_{112} = -y_{11}\,\mathbf{a}_1+\left(\tfrac{1}{2}+x_{11}\right)\mathbf{a}_2-z_{11}\,\mathbf{a}_3 = -y_{11}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_{11}\right)a\,\hat{\mathbf{y}}-z_{11}c\,\hat{\mathbf{z}} \quad (16g) \quad \text{O IV}$$

$$\mathbf{B}_{113} = x_{11}\,\mathbf{a}_1+\left(\tfrac{1}{2}-y_{11}\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_{11}\right)\mathbf{a}_3 = x_{11}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{11}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{11}\right)c\,\hat{\mathbf{z}} \quad (16g) \quad \text{O IV}$$

$$\mathbf{B}_{114} = \left(\tfrac{1}{2}-x_{11}\right)\mathbf{a}_1+y_{11}\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{11}\right)a\,\hat{\mathbf{x}}+y_{11}a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{11}\right)c\,\hat{\mathbf{z}} \quad (16g) \quad \text{O IV}$$

$$\mathbf{B}_{115} = \left(\tfrac{1}{2}-y_{11}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-x_{11}\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-y_{11}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_{11}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{11}\right)c\,\hat{\mathbf{z}} \quad (16g) \quad \text{O IV}$$

$$\mathbf{B}_{116} = y_{11}\,\mathbf{a}_1+x_{11}\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_{11}\right)\mathbf{a}_3 = y_{11}a\,\hat{\mathbf{x}}+x_{11}a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{11}\right)c\,\hat{\mathbf{z}} \quad (16g) \quad \text{O IV}$$

References:

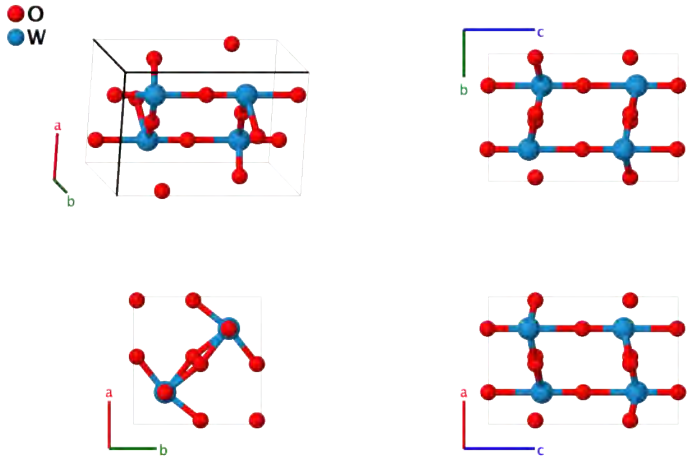
- J. S. Ricci, R. C. Stevens, R. K. McMullan, and W. T. Klooster, *Structure of strontium hydroxide octahydrate, $Sr(OH)_2{\cdot}8H_2O$, at 20, 100 and 200 K from neutron diffraction*, Acta Crystallogr. Sect. B Struct. Sci. **61**, 381–386 (2005), doi:10.1107/S0108768105013480.
- G. Natta, *Constitution of hydroxides and of hydrates. III. Octahydrated strontium hydroxide*, Gazz. Chim. Ital. **58**, 870–872 (1928).
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1700
- POSCAR: pp. 1700

α-WO_3 Structure: A3B_tP16_130_cf_c

http://aflow.org/prototype-encyclopedia/A3B_tP16_130_cf_c



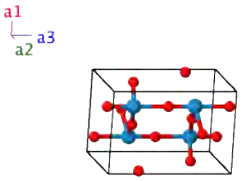
Prototype	:	O_3W
AFLOW prototype label	:	A3B_tP16_130_cf_c
***Strukturbericht* designation**	:	None
Pearson symbol	:	tP16
Space group number	:	130
Space group symbol	:	$P4/ncc$
AFLOW prototype command	:	`aflow --proto=A3B_tP16_130_cf_c --params=`$a, c/a, z_1, z_2, x_3$

- All stable phases of WO_3 are distortions of the cubic α-ReO_3 ($D0_9$) phase. Based on (Woodward, 1997 and Vogt, 1999), the known stable phases and their approximate temperature ranges are:
  - α-WO_3 (1010-1170 K) (Vogt, 1999), this structure
  - β-WO_3 (600-1170 K) (Vogt, 1999)
  - γ-WO_3 (290-600 K) (Vogt, 1999)
  - δ-WO_3 (230-290 K) (Diehl, 1978)
  - ϵ-WO_3 (below 23 K) (Woodward, 1997)

  Woodward's caption says "The transition temperatures display large hysteresis effects and universal agreement is not found in the literature."
- In addition, several other structures have been proposed and/or found:
  - The original $D0_{10}$ structure (Bräkken, 1931), (Hermann, 1937) superseded by δ-WO_3
  - Original β-WO_3 (Salje, 1977)
  - Hexagonal WO_3 (Gerand, 1979) (metastable)

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	$(4c)$	O I
$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right)c\,\hat{\mathbf{z}}$	$(4c)$	O I
$\mathbf{B}_3$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_1\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	$(4c)$	O I
$\mathbf{B}_4$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\,\hat{\mathbf{z}}$	$(4c)$	O I
$\mathbf{B}_5$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(4c)$	W
$\mathbf{B}_6$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	$(4c)$	W
$\mathbf{B}_7$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	$(4c)$	W
$\mathbf{B}_8$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	$(4c)$	W
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{11}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{12}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{13}$	=	$-x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{15}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(8f)$	O II
$\mathbf{B}_{16}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(8f)$	O II

References:

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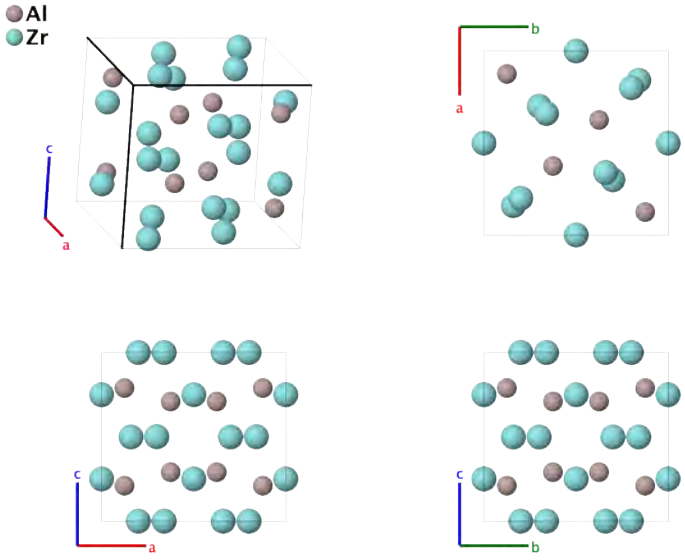
Geometry files:
- CIF: pp. 1701
- POSCAR: pp. 1701

Zr_3Al_2 Structure: A2B3_tP20_136_j_dfg

http://aflow.org/prototype-encyclopedia/A2B3_tP20_136_j_dfg



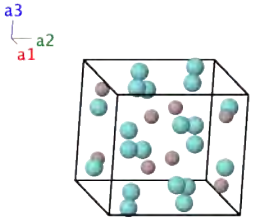
Prototype	:	Al_2Zr_3
AFLOW prototype label	:	A2B3_tP20_136_j_dfg
***Strukturbericht* designation**	:	None
Pearson symbol	:	tP20
Space group number	:	136
Space group symbol	:	$P4_2/mnm$
AFLOW prototype command	:	`aflow --proto=A2B3_tP20_136_j_dfg --params=`$a, c/a, x_2, x_3, x_4, z_4$

Other compounds with this structure

- Hf_3Al_2 and Y_3Al_2

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(4d)$	Zr I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(4d)$	Zr I

$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($4d$)	Zr I
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	($4d$)	Zr I
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2$	=	$x_2a\,\hat{\mathbf{x}} + x_2a\,\hat{\mathbf{y}}$	($4f$)	Zr II
$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2$	=	$-x_2a\,\hat{\mathbf{x}} - x_2a\,\hat{\mathbf{y}}$	($4f$)	Zr II
$\mathbf{B}_7$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4f$)	Zr II
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4f$)	Zr II
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2$	=	$x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}}$	($4g$)	Zr III
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2$	=	$-x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}}$	($4g$)	Zr III
$\mathbf{B}_{11}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4g$)	Zr III
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4g$)	Zr III
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	($8j$)	Al
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	($8j$)	Al
$\mathbf{B}_{15}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	($8j$)	Al
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	($8j$)	Al
$\mathbf{B}_{17}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	($8j$)	Al
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	($8j$)	Al
$\mathbf{B}_{19}$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	($8j$)	Al
$\mathbf{B}_{20}$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	($8j$)	Al

References:

- C. G. Wilson and F. J. Spooner, *The Crystal Structure of Zr_3Al_2*, Acta Cryst. **13**, 358–359 (1960), doi:10.1107/S0365110X60000844.

Found in:

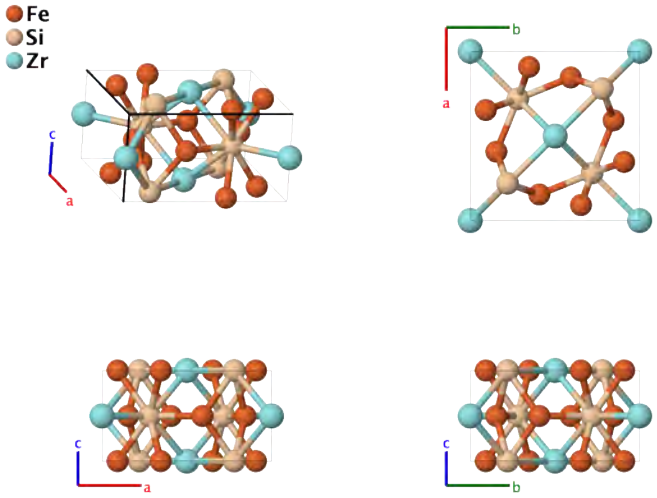
- L.-E. Edshammar, *Crystal Structure Investigations on the Zr-Al and Hf-Al Systems*, Acta Chem. Scand. **16**, 20–30 (1962), doi:10.3891/acta.chem.scand.16-0020.

Geometry files:

- CIF: pp. 1701
- POSCAR: pp. 1702

$ZrFe_4Si_2$ Structure: A4B2C_tP14_136_i_g_b

http://aflow.org/prototype-encyclopedia/A4B2C_tP14_136_i_g_b



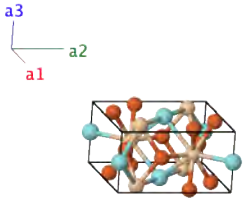
Prototype	:	Fe_4Si_2Zr
AFLOW prototype label	:	A4B2C_tP14_136_i_g_b
***Strukturbericht* designation**	:	None
Pearson symbol	:	tP14
Space group number	:	136
Space group symbol	:	$P4_2/mnm$
AFLOW prototype command	:	`aflow --proto=A4B2C_tP14_136_i_g_b --params=`$a, c/a, x_2, x_3, y_3$

Other compounds with this structure

- $BaCd_4Pt_2$, $DyFe_4Ge_2$, $DyNi_4As_2$, $ErFe_4Ge_2$, $ErNi_4P_2$, $GdNi_4As_2$, $GdRe_4Si_2$, $HoFe_4Ge_2$, $LuFe_4Ge_2$, $LuNi_4As_2$, $LuRe_4Si_2$, $ScFe_4P_2$, $ScFe_4Si_2$, $ScNi_4As_2$, $SmRe_4Si_2$, $SrCd_4Pt_2$, $TbRe_4Si_2$, $TmFe_4Ge_2$, $TmRe_4Si_2$, UMn_4P_2, YFe_4Ge_2, YNi_4As_2, YNi_4P_2, YRe_4Si_2, $YbNi_4P_2$, $ZrFe_4P_2$, and $ZrNi_4As_2$

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	$(2b)$	Zr

$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$	($2b$)	Zr
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2$	=	$x_2 a\,\hat{\mathbf{x}} - x_2 a\,\hat{\mathbf{y}}$	($4g$)	Si
$\mathbf{B}_4$	=	$-x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2$	=	$-x_2 a\,\hat{\mathbf{x}} + x_2 a\,\hat{\mathbf{y}}$	($4g$)	Si
$\mathbf{B}_5$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4g$)	Si
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4g$)	Si
$\mathbf{B}_7$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 a\,\hat{\mathbf{y}}$	($8i$)	Fe
$\mathbf{B}_8$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 a\,\hat{\mathbf{y}}$	($8i$)	Fe
$\mathbf{B}_9$	=	$\left(\frac{1}{2} - y_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($8i$)	Fe
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} + y_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($8i$)	Fe
$\mathbf{B}_{11}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($8i$)	Fe
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($8i$)	Fe
$\mathbf{B}_{13}$	=	$y_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2$	=	$y_3 a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}}$	($8i$)	Fe
$\mathbf{B}_{14}$	=	$-y_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2$	=	$-y_3 a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}}$	($8i$)	Fe

References:

- Y. P. Yarmolyuk, L. A. Lysenko, and E. I. Gladyshevsky, *Crystal Structure of $ZrFe_4Si_2$ – A New Structure Type of Ternary Silicides*, Dopov. Akad. Nauk Ukr. RSR, Ser. A **37**, 281–284 (1975). In Russian.

Found in:

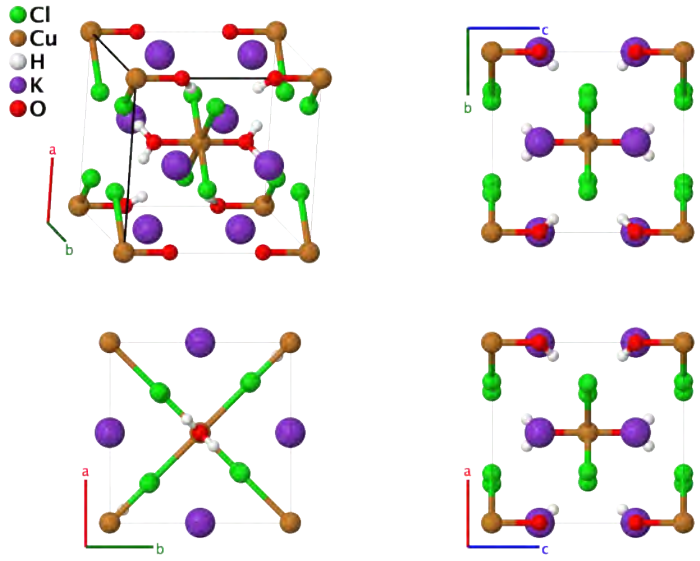
- P. Schobinger-Papamantellos, J. Rodríguez-Carvajal, G. André, N. P. Duong, K. H. J. Buschow, and P. Tolédano, *Simultaneous structural and magnetic transitions in YFe_4Ge_2 studied by neutron diffraction and magnetic measurements*, J. Magn. Magn. Mater. **236**, 14–27 (2001), doi:10.1016/S0304-8853(01)00442-5.

Geometry files:

- CIF: pp. 1702
- POSCAR: pp. 1702

$K_2CuCl_4{\cdot}2H_2O$ ($H4_1$) Structure: A4BC4D2E2_tP26_136_fg_a_j_d_e

http://aflow.org/prototype-encyclopedia/A4BC4D2E2_tP26_136_fg_a_j_d_e



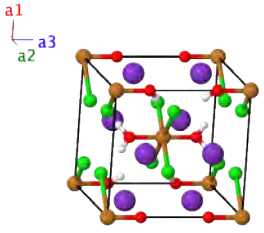
Prototype	:	$Cl_4CuH_4K_2O_2$
AFLOW prototype label	:	A4BC4D2E2_tP26_136_fg_a_j_d_e
***Strukturbericht* designation**	:	$H4_1$
Pearson symbol	:	tP26
Space group number	:	136
Space group symbol	:	$P4_2/mnm$
AFLOW prototype command	:	`aflow --proto=A4BC4D2E2_tP26_136_fg_a_j_d_e --params=`$a, c/a, z_3, x_4, x_5, x_6, z_6$

Other compounds with this structure

- $Rb_2CuCl_4 \cdot 2H_2O$, $(NH_4)_2CuCl_4 \cdot 2H_2O$, and $(NH_4)_2CuBr_4 \cdot 2H_2O$

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$ =	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	Cu

$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(2*a*)	Cu
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*d*)	K
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(4*d*)	K
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*d*)	K
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(4*d*)	K
$\mathbf{B}_7$	=	$z_3\,\mathbf{a}_3$	=	$z_3 c\,\hat{\mathbf{z}}$	(4*e*)	O
$\mathbf{B}_8$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(4*e*)	O
$\mathbf{B}_9$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(4*e*)	O
$\mathbf{B}_{10}$	=	$-z_3\,\mathbf{a}_3$	=	$-z_3 c\,\hat{\mathbf{z}}$	(4*e*)	O
$\mathbf{B}_{11}$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2$	=	$x_4 a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}}$	(4*f*)	Cl I
$\mathbf{B}_{12}$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2$	=	$-x_4 a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}}$	(4*f*)	Cl I
$\mathbf{B}_{13}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_4\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*f*)	Cl I
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*f*)	Cl I
$\mathbf{B}_{15}$	=	$x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2$	=	$x_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}}$	(4*g*)	Cl II
$\mathbf{B}_{16}$	=	$-x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2$	=	$-x_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}}$	(4*g*)	Cl II
$\mathbf{B}_{17}$	=	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_5\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*g*)	Cl II
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_5\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*g*)	Cl II
$\mathbf{B}_{19}$	=	$x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(8*j*)	H
$\mathbf{B}_{20}$	=	$-x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(8*j*)	H
$\mathbf{B}_{21}$	=	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_6\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\,\hat{\mathbf{z}}$	(8*j*)	H
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_6\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\,\hat{\mathbf{z}}$	(8*j*)	H
$\mathbf{B}_{23}$	=	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_6\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\,\hat{\mathbf{z}}$	(8*j*)	H
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_6\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\,\hat{\mathbf{z}}$	(8*j*)	H
$\mathbf{B}_{25}$	=	$x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(8*j*)	H
$\mathbf{B}_{26}$	=	$-x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(8*j*)	H

References:

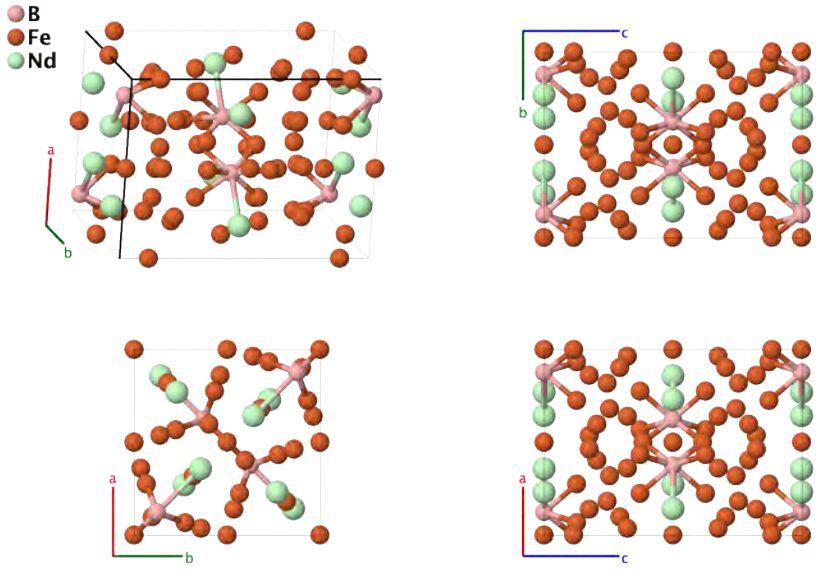
- R. Chidambaram, Q. O. Navarro, A. Garcia, K. Linggoatmodjo, L. Shi-Chien, and I.-H. Suh, *Neutron diffraction refinement of the crystal structure of potassium copper chloride dihydrate, $K_2CuCl_4 \cdot 2H_2O$*, Acta Crystallogr. Sect. B Struct. Sci. **26**, 827–830 (1970), doi:10.1107/S0567740870003187.

Geometry files:

- CIF: pp. 1702
- POSCAR: pp. 1703

$Nd_2Fe_{14}B$ Structure: AB14C2_tP68_136_f_ce2j2k_fg

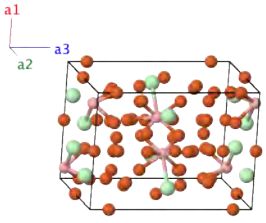
http://aflow.org/prototype-encyclopedia/AB14C2_tP68_136_f_ce2j2k_fg



Prototype	:	$BFe_{14}Nd_2$
AFLOW prototype label	:	AB14C2_tP68_136_f_ce2j2k_fg
***Strukturbericht* designation**	:	None
Pearson symbol	:	tP68
Space group number	:	136
Space group symbol	:	$P4_2/mnm$
AFLOW prototype command	:	`aflow --proto=AB14C2_tP68_136_f_ce2j2k_fg --params=`$a, c/a, z_2, x_3, x_4, x_5, x_6, z_6, x_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9$

Simple Tetragonal primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{y}}$	($4c$)	Fe I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4c$)	Fe I
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	($4c$)	Fe I
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	($4c$)	Fe I
$\mathbf{B}_5$	=	$z_2\,\mathbf{a}_3$	=	$z_2 c\,\hat{\mathbf{z}}$	($4e$)	Fe II

$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	$(4e)$	Fe II
$\mathbf{B}_7$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	$(4e)$	Fe II
$\mathbf{B}_8$	=	$-z_2\,\mathbf{a}_3$	=	$-z_2c\,\hat{\mathbf{z}}$	$(4e)$	Fe II
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2$	=	$x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}}$	$(4f)$	B
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2$	=	$-x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}}$	$(4f)$	B
$\mathbf{B}_{11}$	=	$\left(\frac{1}{2} - x_3\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(4f)$	B
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(4f)$	B
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2$	=	$x_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}}$	$(4f)$	Nd I
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2$	=	$-x_4a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}}$	$(4f)$	Nd I
$\mathbf{B}_{15}$	=	$\left(\frac{1}{2} - x_4\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + x_4\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(4f)$	Nd I
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_4\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - x_4\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(4f)$	Nd I
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2$	=	$x_5a\,\hat{\mathbf{x}} - x_5a\,\hat{\mathbf{y}}$	$(4g)$	Nd II
$\mathbf{B}_{18}$	=	$-x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2$	=	$-x_5a\,\hat{\mathbf{x}} + x_5a\,\hat{\mathbf{y}}$	$(4g)$	Nd II
$\mathbf{B}_{19}$	=	$\left(\frac{1}{2} + x_5\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + x_5\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(4g)$	Nd II
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} - x_5\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - x_5\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(4g)$	Nd II
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$x_6a\,\hat{\mathbf{x}} + x_6a\,\hat{\mathbf{y}} + z_6c\,\hat{\mathbf{z}}$	$(8j)$	Fe III
$\mathbf{B}_{22}$	=	$-x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$-x_6a\,\hat{\mathbf{x}} - x_6a\,\hat{\mathbf{y}} + z_6c\,\hat{\mathbf{z}}$	$(8j)$	Fe III
$\mathbf{B}_{23}$	=	$\left(\frac{1}{2} - x_6\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + x_6\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\,\hat{\mathbf{z}}$	$(8j)$	Fe III
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} + x_6\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - x_6\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\,\hat{\mathbf{z}}$	$(8j)$	Fe III
$\mathbf{B}_{25}$	=	$\left(\frac{1}{2} - x_6\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + x_6\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\,\hat{\mathbf{z}}$	$(8j)$	Fe III
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} + x_6\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - x_6\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\,\hat{\mathbf{z}}$	$(8j)$	Fe III
$\mathbf{B}_{27}$	=	$x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$x_6a\,\hat{\mathbf{x}} + x_6a\,\hat{\mathbf{y}} - z_6c\,\hat{\mathbf{z}}$	$(8j)$	Fe III
$\mathbf{B}_{28}$	=	$-x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$-x_6a\,\hat{\mathbf{x}} - x_6a\,\hat{\mathbf{y}} - z_6c\,\hat{\mathbf{z}}$	$(8j)$	Fe III
$\mathbf{B}_{29}$	=	$x_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$x_7a\,\hat{\mathbf{x}} + x_7a\,\hat{\mathbf{y}} + z_7c\,\hat{\mathbf{z}}$	$(8j)$	Fe IV
$\mathbf{B}_{30}$	=	$-x_7\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$-x_7a\,\hat{\mathbf{x}} - x_7a\,\hat{\mathbf{y}} + z_7c\,\hat{\mathbf{z}}$	$(8j)$	Fe IV
$\mathbf{B}_{31}$	=	$\left(\frac{1}{2} - x_7\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + x_7\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_7\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right)c\,\hat{\mathbf{z}}$	$(8j)$	Fe IV
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2} + x_7\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - x_7\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_7\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right)c\,\hat{\mathbf{z}}$	$(8j)$	Fe IV
$\mathbf{B}_{33}$	=	$\left(\frac{1}{2} - x_7\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + x_7\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_7\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right)c\,\hat{\mathbf{z}}$	$(8j)$	Fe IV
$\mathbf{B}_{34}$	=	$\left(\frac{1}{2} + x_7\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - x_7\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_7\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right)c\,\hat{\mathbf{z}}$	$(8j)$	Fe IV
$\mathbf{B}_{35}$	=	$x_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$x_7a\,\hat{\mathbf{x}} + x_7a\,\hat{\mathbf{y}} - z_7c\,\hat{\mathbf{z}}$	$(8j)$	Fe IV

$\mathbf{B}_{36}$	=	$-x_7\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} - x_7 a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	$(8j)$	Fe IV
$\mathbf{B}_{37}$	=	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + y_8 a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	$(16k)$	Fe V
$\mathbf{B}_{38}$	=	$-x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} - y_8 a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	$(16k)$	Fe V
$\mathbf{B}_{39}$	=	$\left(\frac{1}{2} - y_8\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_8\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_8\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right)c\,\hat{\mathbf{z}}$	$(16k)$	Fe V
$\mathbf{B}_{40}$	=	$\left(\frac{1}{2} + y_8\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_8\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_8\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right)c\,\hat{\mathbf{z}}$	$(16k)$	Fe V
$\mathbf{B}_{41}$	=	$\left(\frac{1}{2} - x_8\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_8\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_8\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right)c\,\hat{\mathbf{z}}$	$(16k)$	Fe V
$\mathbf{B}_{42}$	=	$\left(\frac{1}{2} + x_8\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_8\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_8\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right)c\,\hat{\mathbf{z}}$	$(16k)$	Fe V
$\mathbf{B}_{43}$	=	$y_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$y_8 a\,\hat{\mathbf{x}} + x_8 a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	$(16k)$	Fe V
$\mathbf{B}_{44}$	=	$-y_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$-y_8 a\,\hat{\mathbf{x}} - x_8 a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	$(16k)$	Fe V
$\mathbf{B}_{45}$	=	$-x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} - y_8 a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	$(16k)$	Fe V
$\mathbf{B}_{46}$	=	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + y_8 a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	$(16k)$	Fe V
$\mathbf{B}_{47}$	=	$\left(\frac{1}{2} + y_8\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_8\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_8\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right)c\,\hat{\mathbf{z}}$	$(16k)$	Fe V
$\mathbf{B}_{48}$	=	$\left(\frac{1}{2} - y_8\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_8\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_8\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right)c\,\hat{\mathbf{z}}$	$(16k)$	Fe V
$\mathbf{B}_{49}$	=	$\left(\frac{1}{2} + x_8\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_8\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_8\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right)c\,\hat{\mathbf{z}}$	$(16k)$	Fe V
$\mathbf{B}_{50}$	=	$\left(\frac{1}{2} - x_8\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_8\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_8\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right)c\,\hat{\mathbf{z}}$	$(16k)$	Fe V
$\mathbf{B}_{51}$	=	$-y_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$-y_8 a\,\hat{\mathbf{x}} - x_8 a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	$(16k)$	Fe V
$\mathbf{B}_{52}$	=	$y_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$y_8 a\,\hat{\mathbf{x}} + x_8 a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	$(16k)$	Fe V
$\mathbf{B}_{53}$	=	$x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}} + y_9 a\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	$(16k)$	Fe VI
$\mathbf{B}_{54}$	=	$-x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{x}} - y_9 a\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	$(16k)$	Fe VI
$\mathbf{B}_{55}$	=	$\left(\frac{1}{2} - y_9\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_9\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_9\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_9\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right)c\,\hat{\mathbf{z}}$	$(16k)$	Fe VI
$\mathbf{B}_{56}$	=	$\left(\frac{1}{2} + y_9\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_9\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_9\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_9\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right)c\,\hat{\mathbf{z}}$	$(16k)$	Fe VI
$\mathbf{B}_{57}$	=	$\left(\frac{1}{2} - x_9\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_9\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_9\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_9\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_9\right)c\,\hat{\mathbf{z}}$	$(16k)$	Fe VI
$\mathbf{B}_{58}$	=	$\left(\frac{1}{2} + x_9\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_9\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_9\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_9\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_9\right)c\,\hat{\mathbf{z}}$	$(16k)$	Fe VI
$\mathbf{B}_{59}$	=	$y_9\,\mathbf{a}_1 + x_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$y_9 a\,\hat{\mathbf{x}} + x_9 a\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}}$	$(16k)$	Fe VI
$\mathbf{B}_{60}$	=	$-y_9\,\mathbf{a}_1 - x_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$-y_9 a\,\hat{\mathbf{x}} - x_9 a\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}}$	$(16k)$	Fe VI
$\mathbf{B}_{61}$	=	$-x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{x}} - y_9 a\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}}$	$(16k)$	Fe VI
$\mathbf{B}_{62}$	=	$x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}} + y_9 a\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}}$	$(16k)$	Fe VI

$\mathbf{B}_{63}$	=	$\left(\frac{1}{2}+y_9\right)\mathbf{a}_1+\left(\frac{1}{2}-x_9\right)\mathbf{a}_2+\left(\frac{1}{2}-z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_9\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_9\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_9\right)c\,\hat{\mathbf{z}}$	(16*k*)	Fe VI
$\mathbf{B}_{64}$	=	$\left(\frac{1}{2}-y_9\right)\mathbf{a}_1+\left(\frac{1}{2}+x_9\right)\mathbf{a}_2+\left(\frac{1}{2}-z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_9\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_9\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_9\right)c\,\hat{\mathbf{z}}$	(16*k*)	Fe VI
$\mathbf{B}_{65}$	=	$\left(\frac{1}{2}+x_9\right)\mathbf{a}_1+\left(\frac{1}{2}-y_9\right)\mathbf{a}_2+\left(\frac{1}{2}+z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_9\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_9\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_9\right)c\,\hat{\mathbf{z}}$	(16*k*)	Fe VI
$\mathbf{B}_{66}$	=	$\left(\frac{1}{2}-x_9\right)\mathbf{a}_1+\left(\frac{1}{2}+y_9\right)\mathbf{a}_2+\left(\frac{1}{2}+z_9\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_9\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_9\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_9\right)c\,\hat{\mathbf{z}}$	(16*k*)	Fe VI
$\mathbf{B}_{67}$	=	$-y_9\,\mathbf{a}_1-x_9\,\mathbf{a}_2+z_9\,\mathbf{a}_3$	=	$-y_9a\,\hat{\mathbf{x}}-x_9a\,\hat{\mathbf{y}}+z_9c\,\hat{\mathbf{z}}$	(16*k*)	Fe VI
$\mathbf{B}_{68}$	=	$y_9\,\mathbf{a}_1+x_9\,\mathbf{a}_2+z_9\,\mathbf{a}_3$	=	$y_9a\,\hat{\mathbf{x}}+x_9a\,\hat{\mathbf{y}}+z_9c\,\hat{\mathbf{z}}$	(16*k*)	Fe VI

References:

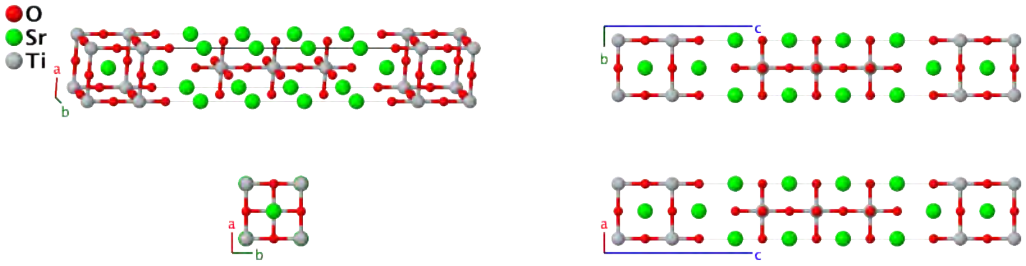
- D. Givord, H. S. Li, and J. M. Moreau, *Magnetic properties and crystal structure of $Nd_2Fe_{14}B$*, Solid State Commun. **50**, 497–499 (1984), doi:10.1016/0038-1098(84)90315-6.

Geometry files:

- CIF: pp. 1703
- POSCAR: pp. 1703

$Sr_4Ti_3O_{10}$ Structure: A10B4C3_tI34_139_c2eg_2e_ae

http://aflow.org/prototype-encyclopedia/A10B4C3_tI34_139_c2eg_2e_ae



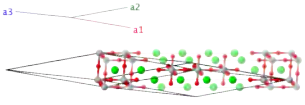
Prototype	:	$O_{10}Sr_4Ti_3$
AFLOW prototype label	:	A10B4C3_tI34_139_c2eg_2e_ae
***Strukturbericht* designation**	:	None
Pearson symbol	:	tI34
Space group number	:	139
Space group symbol	:	$I4/mmm$
AFLOW prototype command	:	`aflow --proto=A10B4C3_tI34_139_c2eg_2e_ae --params=`$a, c/a, z_3, z_4, z_5, z_6, z_7, z_8$

Other compounds with this structure

- $K_2La_2Ti_3O_{10}$, $Li_2Eu_2Ti_3O_{10}$, $Na_2Eu_2Ti_3O_{10}$, and $Na_2Sr_2Nb_2MnO_{10}$

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} - \frac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	Ti I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}}$	$(4c)$	O I
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	$(4c)$	O I
$\mathbf{B}_4$	=	$z_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_2$	=	$z_3c\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_5$	=	$-z_3\,\mathbf{a}_1 - z_3\,\mathbf{a}_2$	=	$-z_3c\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_6$	=	$z_4\,\mathbf{a}_1 + z_4\,\mathbf{a}_2$	=	$z_4c\,\hat{\mathbf{z}}$	$(4e)$	O III
$\mathbf{B}_7$	=	$-z_4\,\mathbf{a}_1 - z_4\,\mathbf{a}_2$	=	$-z_4c\,\hat{\mathbf{z}}$	$(4e)$	O III
$\mathbf{B}_8$	=	$z_5\,\mathbf{a}_1 + z_5\,\mathbf{a}_2$	=	$z_5c\,\hat{\mathbf{z}}$	$(4e)$	Sr I
$\mathbf{B}_9$	=	$-z_5\,\mathbf{a}_1 - z_5\,\mathbf{a}_2$	=	$-z_5c\,\hat{\mathbf{z}}$	$(4e)$	Sr I
$\mathbf{B}_{10}$	=	$z_6\,\mathbf{a}_1 + z_6\,\mathbf{a}_2$	=	$z_6c\,\hat{\mathbf{z}}$	$(4e)$	Sr II
$\mathbf{B}_{11}$	=	$-z_6\,\mathbf{a}_1 - z_6\,\mathbf{a}_2$	=	$-z_6c\,\hat{\mathbf{z}}$	$(4e)$	Sr II

$\mathbf{B}_{12}$	=	$z_7\,\mathbf{a}_1 + z_7\,\mathbf{a}_2$	=	$z_7 c\,\hat{\mathbf{z}}$	($4e$)	Ti II
$\mathbf{B}_{13}$	=	$-z_7\,\mathbf{a}_1 - z_7\,\mathbf{a}_2$	=	$-z_7 c\,\hat{\mathbf{z}}$	($4e$)	Ti II
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} + z_8\right)\mathbf{a}_1 + z_8\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	($8g$)	O IV
$\mathbf{B}_{15}$	=	$z_8\,\mathbf{a}_1 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + z_8 c\,\hat{\mathbf{z}}$	($8g$)	O IV
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} - z_8\right)\mathbf{a}_1 - z_8\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	($8g$)	O IV
$\mathbf{B}_{17}$	=	$-z_8\,\mathbf{a}_1 + \left(\frac{1}{2} - z_8\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - z_8 c\,\hat{\mathbf{z}}$	($8g$)	O IV

References:

- S. N. Ruddlesden and P. Popper, *The compound $Sr_3Ti_2O_7$ and its structure*, Acta Cryst. **11**, 54–55 (1958), doi:10.1107/S0365110X58000128.

Found in:

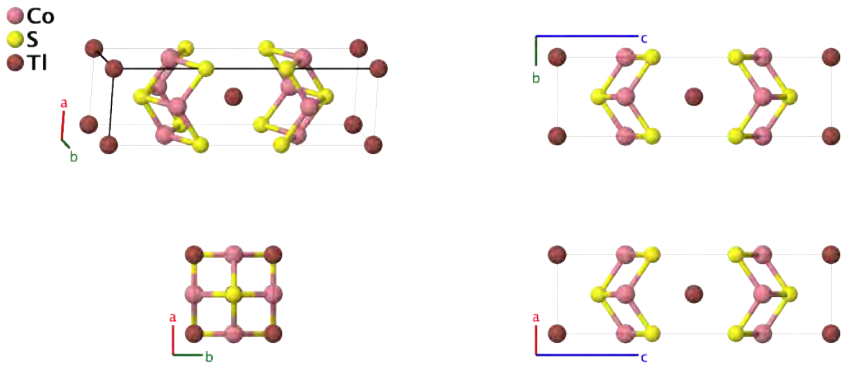
- Wikipedia, *Ruddlesden-Popper phase*, https://en.wikipedia.org/wiki/Ruddlesden-Popper_phase. $A_3B_2X_7$ series.

Geometry files:

- CIF: pp. 1704
- POSCAR: pp. 1704

$TlCo_2S_2$ Structure: A2B2C_tI10_139_d_e_a

http://aflow.org/prototype-encyclopedia/A2B2C_tI10_139_d_e_a.TlCo2S2



Prototype	:	Co_2S_2Tl
AFLOW prototype label	:	A2B2C_tI10_139_d_e_a
***Strukturbericht* designation**	:	None
Pearson symbol	:	tI10
Space group number	:	139
Space group symbol	:	$I4/mmm$
AFLOW prototype command	:	`aflow --proto=A2B2C_tI10_139_d_e_a --params=`$a, c/a, z_3$

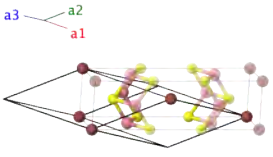
Other compounds with this structure

- $BaCr_2As_2$, $BaFe_2As_2$, BaP_2Zn_2, $BaSr_2As_2$, $BiCe_2O_2$, KCo_2Se_2, $TeCe_2O_2$, $TlCo_2Se_2$, $TlCu_2S_2$, $TlCu_2Se_2$, $TlCu_2Te_2$, $TlFe_2S_2$, $TlFe_2Se_2$, $TlNi_2S_2$, $TlNi_2Se_2$, and $Cl(Li_{0.25}Be_{0.75})_2O_2$

- This is a ternary form of the $D1_3$ ($BaAl_4$) structure. It differs from the $ThCr_2Si_2$ structure in that here $c/a > 3$.
- Other authors designate $TlCu_2Se_2$ or $Cl(Li_{0.25}Be_{0.75})_2O_2$ as the prototype for this structure, but (Klepp, 1978) give complete structural information for $TlCo_2S_2$, so we use that as the prototype.

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} - \frac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	Tl
$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(4d)$	Co
$\mathbf{B}_3$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(4d)$	Co

$\mathbf{B}_4$	=	$z_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_2$	=	$z_3 c\,\hat{\mathbf{z}}$	($4e$)	S
$\mathbf{B}_5$	=	$-z_3\,\mathbf{a}_1 - z_3\,\mathbf{a}_2$	=	$-z_3 c\,\hat{\mathbf{z}}$	($4e$)	S

References:

- K. Klepp and H. Boller, *Ternäre Thallium-Übergangsmetall-Chalkogenide mit $ThCr_2Si_2$-Struktur*, Monatshefte für Chemie - Chemical Monthly **109**, 1049–1057 (1978), doi:10.1007/BF00913007.

Found in:

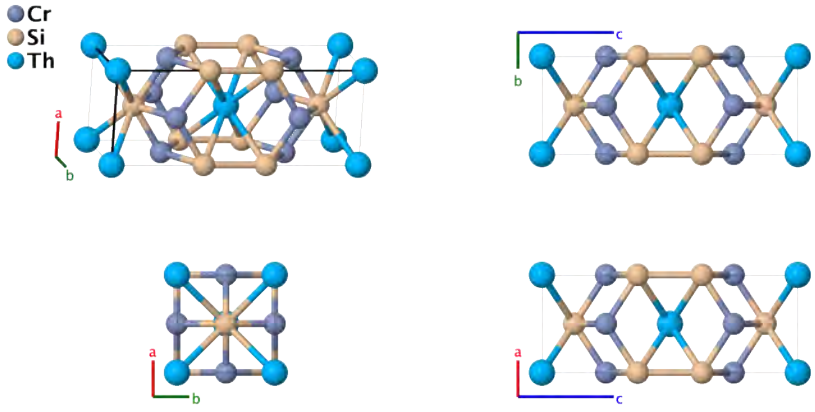
- R. Berger and C. F. Van Bruggen, *$TlCu_2Se_2$: A p-type metal with a layer structure*, J. Less-Common Met. **99**, 113–123 (1984), doi:10.1016/0022-5088(84)90340-0.

Geometry files:

- CIF: pp. 1704
- POSCAR: pp. 1705

$ThCr_2Si_2$ Structure: A2B2C_tI10_139_d_e_a

http://aflow.org/prototype-encyclopedia/A2B2C_tI10_139_d_e_a



Prototype	:	Cr_2Si_2Th
AFLOW prototype label	:	A2B2C_tI10_139_d_e_a
***Strukturbericht* designation**	:	None
Pearson symbol	:	tI10
Space group number	:	139
Space group symbol	:	$I4/mmm$
AFLOW prototype command	:	`aflow --proto=A2B2C_tI10_139_d_e_a --params=`$a, c/a, z_3$

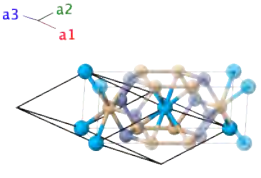
Other compounds with this structure

- $BaAl_2Ge_2$, $BaFe_2As_2$, $BaFe_2P_2$, $BaMn_2Ge_2$, $BaNi_2As_2$, $BaNi_2Ge_2$, $BaRh_2B_2$, $CaAl_2Zn_2$, $CaAu_2Si_2$, $CaFe_2As_2$, $CaMn_2Ge_2$, $CaNi_2Ge_2$, $CeAl_2Ga_2$, $CsCo_2As_2$, $CsCo_2P_2$, $CsFe_2As_2$, $CsFe_2P_2$, $CsIr_2As_2$, $CsIr_2P_2$, $CsRh_2As_2$, $CsRh_2P_2$, $CsRu_2As_2$, $CsRu_2P_2$, $DyCr_2Si_2$, $EuCo_2As_2$, $EuCo_2P_2$, $EuFe_2P_2$, $EuRh_2P_2$, KNi_2S_2, KNi_2Se_2, $LaCo_2P_2$, $LaRu_2P_2$, $LuFe_2B_2$, $LuRu_2Si_2$, $PuCr_2Si_2$, $SrCo_2P_2$, $SrFe_2As_2$, $SrFe_2P_2$, $SrNi_2P_2$, $SrRh_2P_2$, $SrRu_2P_2$, $ThCu_2Si_2$, $ThMn_2Ge_2$, $ThMn_2Si_2$, $ThNi_2Si_2$, UCr_2Si_2, and YNi_2Ge_2

- (Shatruk, 2019) refers to this as "the perovskite of intermetallics." The list of compounds above is by no means complete.
- This is a ternary form of the $D1_3$ ($BaAl_4$) structure.
- The structures generally identified as having prototype $ThCr_2Si_2$ actually divide into two distinct regions, based on c/a ratio. We assign structures with $c/a \leq 3$ to the $ThCr_2Si_2$ structure, and those with $c/a > 3$ to the $TlCo_2S_2$ structure.

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} - \frac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($2a$)	Th
$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($4d$)	Cr
$\mathbf{B}_3$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($4d$)	Cr
$\mathbf{B}_4$	=	$z_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_2$	=	$z_3 c\,\hat{\mathbf{z}}$	($4e$)	Si
$\mathbf{B}_5$	=	$-z_3\,\mathbf{a}_1 - z_3\,\mathbf{a}_2$	=	$-z_3 c\,\hat{\mathbf{z}}$	($4e$)	Si

References:

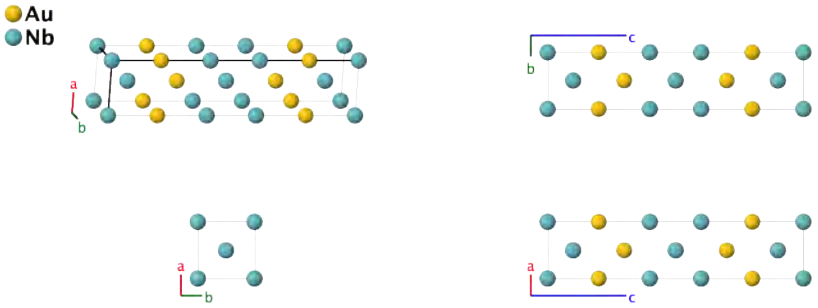
- Z. Ban and M. Sikirica, *The crystal structure of ternary silicides ThM_2Si_2 (M = Cr, Mn, Fe, Co, Ni and Cu)*, Acta Cryst. **18**, 594–599 (1965), doi:10.1107/S0365110X6500141X.
- M. Shatruk, *$ThCr_2Si_2$ structure type: The "perovskite" of intermetallics*, J. Solid State Chem. **272**, 198–209 (2019), doi:10.1016/j.jssc.2019.02.012.

Geometry files:

- CIF: pp. 1705
- POSCAR: pp. 1705

Au_2Nb_3 Structure: A2B3_tI10_139_e_ae

http://aflow.org/prototype-encyclopedia/A2B3_tI10_139_e_ae



Prototype	:	Au_2Nb_3
AFLOW prototype label	:	A2B3_tI10_139_e_ae
***Strukturbericht* designation**	:	None
Pearson symbol	:	tI10
Space group number	:	139
Space group symbol	:	$I4/mmm$
AFLOW prototype command	:	`aflow --proto=A2B3_tI10_139_e_ae --params=`$a, c/a, z_2, z_3$

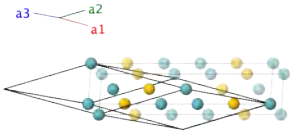
Other compounds with this structure

- Os_2Al_3

- (Schubert, 1960) listed the lattice constants in "kX" units. We used the conversion factor 1/1.00207789 to convert their units into Ångstroms (Arblaster, 1997).
- When $c/a = 5$, $z_2 = 1/5$ and $z_3 = 2/5$, the atoms are on the sites of a body-centered cubic lattice.

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{y}} + \tfrac{1}{2}\,c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2}\,a\,\hat{\mathbf{y}} + \tfrac{1}{2}\,c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{y}} - \tfrac{1}{2}\,c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	Nb I
$\mathbf{B}_2$	=	$z_2\,\mathbf{a}_1 + z_2\,\mathbf{a}_2$	=	$z_2 c\,\hat{\mathbf{z}}$	$(4e)$	Au
$\mathbf{B}_3$	=	$-z_2\,\mathbf{a}_1 - z_2\,\mathbf{a}_2$	=	$-z_2 c\,\hat{\mathbf{z}}$	$(4e)$	Au
$\mathbf{B}_4$	=	$z_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_2$	=	$z_3 c\,\hat{\mathbf{z}}$	$(4e)$	Nb II
$\mathbf{B}_5$	=	$-z_3\,\mathbf{a}_1 - z_3\,\mathbf{a}_2$	=	$-z_3 c\,\hat{\mathbf{z}}$	$(4e)$	Nb II

References:

- K. Schubert, T. R. Anantharaman, H. O. K. Ata, H. G. Meissner, M. Pötzschke, W. Rossteutscher, and E. Stolz, *Einige strukturelle Ergebnisse an metallischen Phasen (6)*, Naturwissenschaften **47**, 512 (1960), doi:10.1007/BF00641115.
- J. W. Arblaster, *Crystallographic Properties of Platinum*, Platin. Met. Ref. **41**, 12–21 (1997). http://www.technology.matthey.com/article/41/1/12-21/.

Found in:

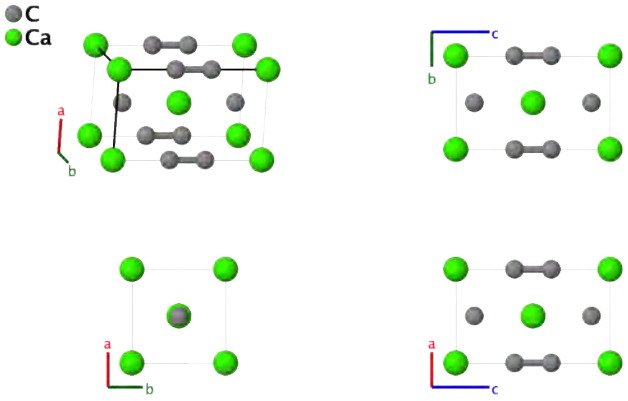
- L.-E. Edshammar, *The Crystal Structures of Os_2Al_3 and $OsAl_2$*, Acta Chem. Scand. **19**, 871–874 (1965), doi:10.3891/acta.chem.scand.19-0871.

Geometry files:

- CIF: pp. 1705
- POSCAR: pp. 1706

CaC_2-I ($C11_a$) Structure: A2B_tI6_139_e_a

http://aflow.org/prototype-encyclopedia/A2B_tI6_139_e_a

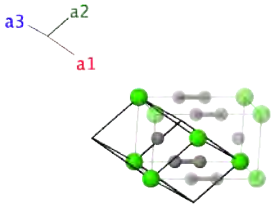


Prototype	:	C_2Ca
AFLOW prototype label	:	A2B_tI6_139_e_a
***Strukturbericht* designation**	:	$C11_a$
Pearson symbol	:	tI6
Space group number	:	139
Space group symbol	:	$I4/mmm$
AFLOW prototype command	:	`aflow --proto=A2B_tI6_139_e_a --params=`$a, c/a, z_2$

- (Ewald, 1931) designated both CaC_2 and $MoSi_2$ as *Strukturbericht* $C11$. (Smithells,1955) separated them into $C11_a$ (CaC_2) and $C11_b$ ($MoSi_2$, AB2_tI6_139_a_e).
- This is the stable room-temperature structure for CaC_2. There is also a meta-stable room temperature structure CaC_2-III (A2B_mC12_12_2i_i), and a low-temperature structure CaC_2, which has the ThC_2 (C_g) structure.
- (von Stackelberg, 1930) describes this structure in a face-centered-tetragonal setting. We follow (Smithells, 1955) and place it in the equivalent body-centered tetragonal setting.

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}a\,\hat{\mathbf{x}} - \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} - \tfrac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	Ca
$\mathbf{B}_2$	=	$z_2\,\mathbf{a}_1 + z_2\,\mathbf{a}_2$	=	$z_2 c\,\hat{\mathbf{z}}$	$(4e)$	C
$\mathbf{B}_3$	=	$-z_2\,\mathbf{a}_1 - z_2\,\mathbf{a}_2$	=	$-z_2 c\,\hat{\mathbf{z}}$	$(4e)$	C

References:

- M. von Stackelberg, *Die Krystallstruktur des CaC_2*, Naturwissenschaften **18**, 305–306 (1930), doi:10.1007/BF01495090.
- P. P. Ewald and C. Hermann, eds., *Strukturbericht 1913-1928* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1931).
- C. J. Smithells, *Metals Reference Book* (Butterworths Scientific, London, 1955), second edn.

Found in:

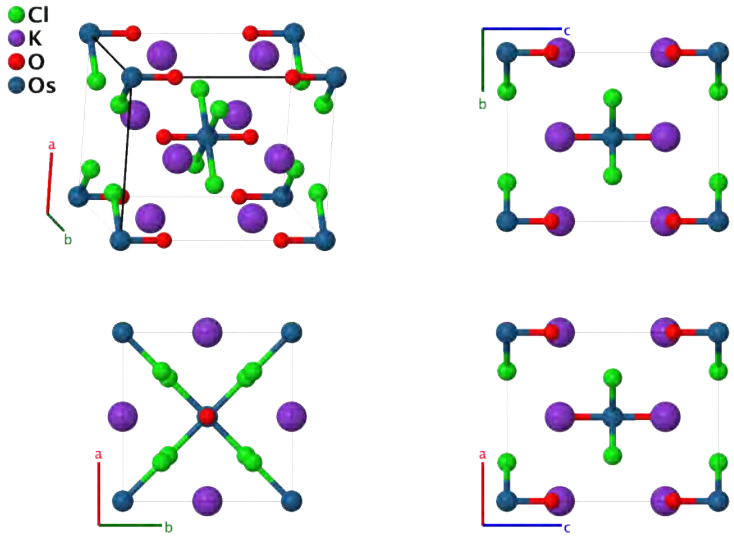
- N.-G. Vannerberg, *The Crystal Structure of Calcium Carbide III*, Acta Chem. Scand. **15**, 769–774 (1961), doi:10.3891/acta.chem.scand.15-0769.

Geometry files:

- CIF: pp. 1706
- POSCAR: pp. 1706

$K_2OsO_2Cl_4$ ($J1_5$) Structure: A4B2C2D_tI18_139_h_d_e_a

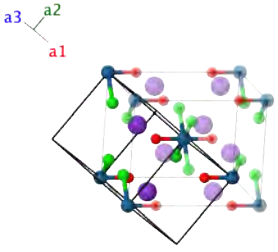
http://aflow.org/prototype-encyclopedia/A4B2C2D_tI18_139_h_d_e_a



Prototype	:	$Cl_4K_2O_2Os$
AFLOW prototype label	:	A4B2C2D_tI18_139_h_d_e_a
***Strukturbericht* designation**	:	$J1_5$
Pearson symbol	:	tI18
Space group number	:	139
Space group symbol	:	$I4/mmm$
AFLOW prototype command	:	`aflow --proto=A4B2C2D_tI18_139_h_d_e_a --params=`$a, c/a, z_3, x_4$

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}a\,\hat{\mathbf{x}} - \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} - \tfrac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	Os
$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(4d)$	K
$\mathbf{B}_3$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(4d)$	K
$\mathbf{B}_4$	=	$z_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_2$	=	$z_3 c\,\hat{\mathbf{z}}$	$(4e)$	O

$\mathbf{B}_5$	=	$-z_3\,\mathbf{a}_1 - z_3\,\mathbf{a}_2$	=	$-z_3 c\,\hat{\mathbf{z}}$	($4e$)	O
$\mathbf{B}_6$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + 2x_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}}$	($8h$)	Cl
$\mathbf{B}_7$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - 2x_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}}$	($8h$)	Cl
$\mathbf{B}_8$	=	$x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2$	=	$-x_4 a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}}$	($8h$)	Cl
$\mathbf{B}_9$	=	$-x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2$	=	$x_4 a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}}$	($8h$)	Cl

References:

- J. L. Hoard and J. D. Grenko, *The Crystal Structure of Potassium Osmyl Chloride, $K_2OsO_2Cl_4$*, Zeitschrift für Kristallographie - Crystalline Materials **87**, 100–109 (1934), doi:10.1524/zkri.1934.87.1.100.

Found in:

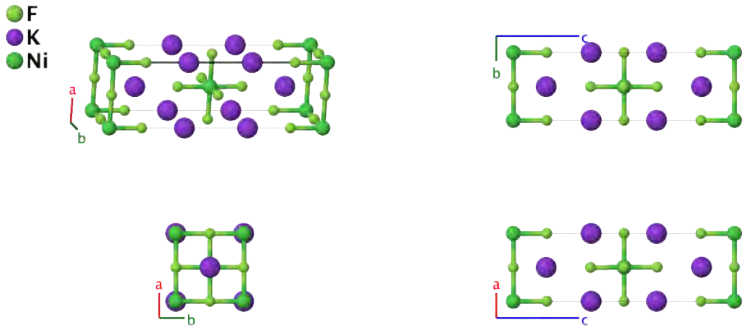
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1706
- POSCAR: pp. 1707

K_2NiF_4 Structure: A4B2C_tI14_139_ce_e_a

http://aflow.org/prototype-encyclopedia/A4B2C_tI14_139_ce_e_a



Prototype	:	F_4K_2Ni
AFLOW prototype label	:	A4B2C_tI14_139_ce_e_a
***Strukturbericht* designation**	:	None
Pearson symbol	:	tI14
Space group number	:	139
Space group symbol	:	$I4/mmm$
AFLOW prototype command	:	`aflow --proto=A4B2C_tI14_139_ce_e_a --params=`$a, c/a, z_3, z_4$

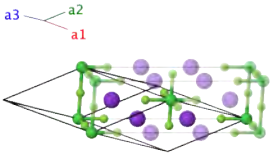
Other compounds with this structure

- Ca_2MnO_4, La_2PdO_4, Nd_2CuO_4, Sr_2MnO_4, Sr_2RuO_4, Sr_2TiO_4, Sr_2VO_4, $(Fe,La)_2SrO_4$, $(Sr,La)_2AlO_4$, and $(Sr,La)_2CoO_4$

- This is the parent compound of the simplest layered-perovskite Ruddlesden-Popper series (Wikipedia). The series also includes the parent of the high-T_c cuprates, $(La,Ba)_2CuO_4$, but we keep that separate as it represents a new class of materials.

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} - \frac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	Ni
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}}$	$(4c)$	F I
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	$(4c)$	F I
$\mathbf{B}_4$	=	$z_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_2$	=	$z_3 c\,\hat{\mathbf{z}}$	$(4e)$	F II
$\mathbf{B}_5$	=	$-z_3\,\mathbf{a}_1 - z_3\,\mathbf{a}_2$	=	$-z_3 c\,\hat{\mathbf{z}}$	$(4e)$	F II

$\mathbf{B}_6$	=	$z_4\,\mathbf{a}_1 + z_4\,\mathbf{a}_2$	=	$z_4 c\,\hat{\mathbf{z}}$	($4e$)	K
$\mathbf{B}_7$	=	$-z_4\,\mathbf{a}_1 - z_4\,\mathbf{a}_2$	=	$-z_4 c\,\hat{\mathbf{z}}$	($4e$)	K

References:

- S. N. Ruddlesden and P. Popper, *New compounds of the K_2NiF_4 type*, Acta Cryst. **10**, 538–539 (1957), doi:10.1107/S0365110X57001929.

Found in:

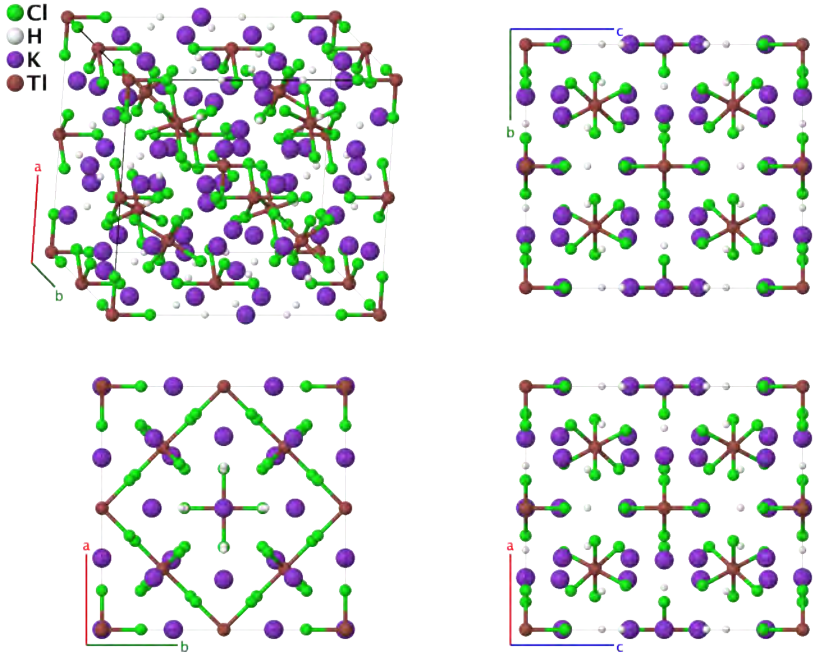
- Wikipedia, *Ruddlesden-Popper phase*, https://en.wikipedia.org/wiki/Ruddlesden-Popper_phase. $A_3B_2X_7$ series.

Geometry files:

- CIF: pp. 1707
- POSCAR: pp. 1707

$K_3TlCl_6{\cdot}2H_2O$ ($J3_1$) Structure: A6B2C3D_tI168_139_egikl2m_ejn_bh2n_acf

http://aflow.org/prototype-encyclopedia/A6B2C3D_tI168_139_egikl2m_ejn_bh2n_acf

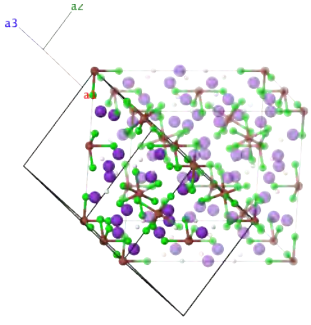


Prototype	:	$Cl_6(H_2O)_2K_3Tl$
AFLOW prototype label	:	A6B2C3D_tI168_139_egikl2m_ejn_bh2n_acf
***Strukturbericht* designation**	:	$J3_1$
Pearson symbol	:	tI168
Space group number	:	139
Space group symbol	:	$I4/mmm$
AFLOW prototype command	:	`aflow --proto=A6B2C3D_tI168_139_egikl2m_ejn_bh2n_acf --params=`$a, c/a, z_4, z_5, z_7, x_8, x_9, x_{10}, x_{11}, x_{12}, y_{12}, x_{13}, z_{13}, x_{14}, z_{14}, y_{15}, z_{15}, y_{16}, z_{16}, y_{17}, z_{17}$

- The positions of the hydrogen atoms in the water molecules were not determined, so we only provide the positions of the oxygen atoms (labeled as H_2O).

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} - \frac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(2*a*)	Tl I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	(2*b*)	K I
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}}$	(4*c*)	Tl II
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	(4*c*)	Tl II
$\mathbf{B}_5$	=	$z_4\,\mathbf{a}_1 + z_4\,\mathbf{a}_2$	=	$z_4 c\,\hat{\mathbf{z}}$	(4*e*)	Cl I
$\mathbf{B}_6$	=	$-z_4\,\mathbf{a}_1 - z_4\,\mathbf{a}_2$	=	$-z_4 c\,\hat{\mathbf{z}}$	(4*e*)	Cl I
$\mathbf{B}_7$	=	$z_5\,\mathbf{a}_1 + z_5\,\mathbf{a}_2$	=	$z_5 c\,\hat{\mathbf{z}}$	(4*e*)	H_2O I
$\mathbf{B}_8$	=	$-z_5\,\mathbf{a}_1 - z_5\,\mathbf{a}_2$	=	$-z_5 c\,\hat{\mathbf{z}}$	(4*e*)	H_2O I
$\mathbf{B}_9$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(8*f*)	Tl III
$\mathbf{B}_{10}$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} - \frac{1}{4}c\,\hat{\mathbf{z}}$	(8*f*)	Tl III
$\mathbf{B}_{11}$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(8*f*)	Tl III
$\mathbf{B}_{12}$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(8*f*)	Tl III
$\mathbf{B}_{13}$	=	$\left(\frac{1}{2} + z_7\right)\,\mathbf{a}_1 + z_7\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(8*g*)	Cl II
$\mathbf{B}_{14}$	=	$z_7\,\mathbf{a}_1 + \left(\frac{1}{2} + z_7\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + z_7 c\,\hat{\mathbf{z}}$	(8*g*)	Cl II
$\mathbf{B}_{15}$	=	$\left(\frac{1}{2} - z_7\right)\,\mathbf{a}_1 - z_7\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	(8*g*)	Cl II
$\mathbf{B}_{16}$	=	$-z_7\,\mathbf{a}_1 + \left(\frac{1}{2} - z_7\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - z_7 c\,\hat{\mathbf{z}}$	(8*g*)	Cl II
$\mathbf{B}_{17}$	=	$x_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 + 2x_8\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + x_8 a\,\hat{\mathbf{y}}$	(8*h*)	K II
$\mathbf{B}_{18}$	=	$-x_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 - 2x_8\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} - x_8 a\,\hat{\mathbf{y}}$	(8*h*)	K II
$\mathbf{B}_{19}$	=	$x_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2$	=	$-x_8 a\,\hat{\mathbf{x}} + x_8 a\,\hat{\mathbf{y}}$	(8*h*)	K II
$\mathbf{B}_{20}$	=	$-x_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2$	=	$x_8 a\,\hat{\mathbf{x}} - x_8 a\,\hat{\mathbf{y}}$	(8*h*)	K II
$\mathbf{B}_{21}$	=	$x_9\,\mathbf{a}_2 + x_9\,\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{x}}$	(8*i*)	Cl III
$\mathbf{B}_{22}$	=	$-x_9\,\mathbf{a}_2 - x_9\,\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{x}}$	(8*i*)	Cl III
$\mathbf{B}_{23}$	=	$x_9\,\mathbf{a}_1 + x_9\,\mathbf{a}_3$	=	$x_9 a\,\hat{\mathbf{y}}$	(8*i*)	Cl III
$\mathbf{B}_{24}$	=	$-x_9\,\mathbf{a}_1 - x_9\,\mathbf{a}_3$	=	$-x_9 a\,\hat{\mathbf{y}}$	(8*i*)	Cl III
$\mathbf{B}_{25}$	=	$\frac{1}{2}\,\mathbf{a}_1 + x_{10}\,\mathbf{a}_2 + \left(\frac{1}{2} + x_{10}\right)\,\mathbf{a}_3$	=	$x_{10} a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$	(8*j*)	H_2O II
$\mathbf{B}_{26}$	=	$\frac{1}{2}\,\mathbf{a}_1 - x_{10}\,\mathbf{a}_2 + \left(\frac{1}{2} - x_{10}\right)\,\mathbf{a}_3$	=	$-x_{10} a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$	(8*j*)	H_2O II
$\mathbf{B}_{27}$	=	$x_{10}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \left(\frac{1}{2} + x_{10}\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + x_{10} a\,\hat{\mathbf{y}}$	(8*j*)	H_2O II
$\mathbf{B}_{28}$	=	$-x_{10}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \left(\frac{1}{2} - x_{10}\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - x_{10} a\,\hat{\mathbf{y}}$	(8*j*)	H_2O II
$\mathbf{B}_{29}$	=	$\left(\frac{3}{4} + x_{11}\right)\,\mathbf{a}_1 + \left(\frac{1}{4} + x_{11}\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + 2x_{11}\right)\,\mathbf{a}_3$	=	$x_{11} a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_{11}\right) a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(16*k*)	Cl IV
$\mathbf{B}_{30}$	=	$\left(\frac{3}{4} - x_{11}\right)\,\mathbf{a}_1 + \left(\frac{1}{4} - x_{11}\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - 2x_{11}\right)\,\mathbf{a}_3$	=	$-x_{11} a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_{11}\right) a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(16*k*)	Cl IV
$\mathbf{B}_{31}$	=	$\left(\frac{1}{4} + x_{11}\right)\,\mathbf{a}_1 + \left(\frac{3}{4} - x_{11}\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{11}\right) a\,\hat{\mathbf{x}} + x_{11} a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(16*k*)	Cl IV
$\mathbf{B}_{32}$	=	$\left(\frac{1}{4} - x_{11}\right)\,\mathbf{a}_1 + \left(\frac{3}{4} + x_{11}\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{11}\right) a\,\hat{\mathbf{x}} - x_{11} a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(16*k*)	Cl IV
$\mathbf{B}_{33}$	=	$\left(\frac{1}{4} - x_{11}\right)\,\mathbf{a}_1 + \left(\frac{3}{4} - x_{11}\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - 2x_{11}\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{11}\right) a\,\hat{\mathbf{x}} - x_{11} a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(16*k*)	Cl IV

$\mathbf{B}_{34}$	=	$\left(\frac{1}{4}+x_{11}\right)\mathbf{a}_1+\left(\frac{3}{4}+x_{11}\right)\mathbf{a}_2+\left(\frac{1}{2}+2x_{11}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{11}\right)a\,\hat{\mathbf{x}}+x_{11}a\,\hat{\mathbf{y}}+\frac{1}{4}c\,\hat{\mathbf{z}}$	(16*k*)	Cl IV
$\mathbf{B}_{35}$	=	$\left(\frac{3}{4}-x_{11}\right)\mathbf{a}_1+\left(\frac{1}{4}+x_{11}\right)\mathbf{a}_2+\frac{1}{2}\mathbf{a}_3$	=	$x_{11}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_{11}\right)a\,\hat{\mathbf{y}}+\frac{1}{4}c\,\hat{\mathbf{z}}$	(16*k*)	Cl IV
$\mathbf{B}_{36}$	=	$\left(\frac{3}{4}+x_{11}\right)\mathbf{a}_1+\left(\frac{1}{4}-x_{11}\right)\mathbf{a}_2+\frac{1}{2}\mathbf{a}_3$	=	$-x_{11}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_{11}\right)a\,\hat{\mathbf{y}}+\frac{1}{4}c\,\hat{\mathbf{z}}$	(16*k*)	Cl IV
$\mathbf{B}_{37}$	=	$y_{12}\,\mathbf{a}_1+x_{12}\,\mathbf{a}_2+(x_{12}+y_{12})\,\mathbf{a}_3$	=	$x_{12}a\,\hat{\mathbf{x}}+y_{12}a\,\hat{\mathbf{y}}$	(16*l*)	Cl V
$\mathbf{B}_{38}$	=	$-y_{12}\,\mathbf{a}_1-x_{12}\,\mathbf{a}_2+(-x_{12}-y_{12})\,\mathbf{a}_3$	=	$-x_{12}a\,\hat{\mathbf{x}}-y_{12}a\,\hat{\mathbf{y}}$	(16*l*)	Cl V
$\mathbf{B}_{39}$	=	$x_{12}\,\mathbf{a}_1-y_{12}\,\mathbf{a}_2+(x_{12}-y_{12})\,\mathbf{a}_3$	=	$-y_{12}a\,\hat{\mathbf{x}}+x_{12}a\,\hat{\mathbf{y}}$	(16*l*)	Cl V
$\mathbf{B}_{40}$	=	$-x_{12}\,\mathbf{a}_1+y_{12}\,\mathbf{a}_2+(-x_{12}+y_{12})\,\mathbf{a}_3$	=	$y_{12}a\,\hat{\mathbf{x}}-x_{12}a\,\hat{\mathbf{y}}$	(16*l*)	Cl V
$\mathbf{B}_{41}$	=	$y_{12}\,\mathbf{a}_1-x_{12}\,\mathbf{a}_2+(-x_{12}+y_{12})\,\mathbf{a}_3$	=	$-x_{12}a\,\hat{\mathbf{x}}+y_{12}a\,\hat{\mathbf{y}}$	(16*l*)	Cl V
$\mathbf{B}_{42}$	=	$-y_{12}\,\mathbf{a}_1+x_{12}\,\mathbf{a}_2+(x_{12}-y_{12})\,\mathbf{a}_3$	=	$x_{12}a\,\hat{\mathbf{x}}-y_{12}a\,\hat{\mathbf{y}}$	(16*l*)	Cl V
$\mathbf{B}_{43}$	=	$x_{12}\,\mathbf{a}_1+y_{12}\,\mathbf{a}_2+(x_{12}+y_{12})\,\mathbf{a}_3$	=	$y_{12}a\,\hat{\mathbf{x}}+x_{12}a\,\hat{\mathbf{y}}$	(16*l*)	Cl V
$\mathbf{B}_{44}$	=	$-x_{12}\,\mathbf{a}_1-y_{12}\,\mathbf{a}_2+(-x_{12}-y_{12})\,\mathbf{a}_3$	=	$-y_{12}a\,\hat{\mathbf{x}}-x_{12}a\,\hat{\mathbf{y}}$	(16*l*)	Cl V
$\mathbf{B}_{45}$	=	$(x_{13}+z_{13})\,\mathbf{a}_1+(x_{13}+z_{13})\,\mathbf{a}_2+2x_{13}\,\mathbf{a}_3$	=	$x_{13}a\,\hat{\mathbf{x}}+x_{13}a\,\hat{\mathbf{y}}+z_{13}c\,\hat{\mathbf{z}}$	(16*m*)	Cl VI
$\mathbf{B}_{46}$	=	$(-x_{13}+z_{13})\,\mathbf{a}_1+(-x_{13}+z_{13})\,\mathbf{a}_2-2x_{13}\,\mathbf{a}_3$	=	$-x_{13}a\,\hat{\mathbf{x}}-x_{13}a\,\hat{\mathbf{y}}+z_{13}c\,\hat{\mathbf{z}}$	(16*m*)	Cl VI
$\mathbf{B}_{47}$	=	$(x_{13}+z_{13})\,\mathbf{a}_1+(-x_{13}+z_{13})\,\mathbf{a}_2$	=	$-x_{13}a\,\hat{\mathbf{x}}+x_{13}a\,\hat{\mathbf{y}}+z_{13}c\,\hat{\mathbf{z}}$	(16*m*)	Cl VI
$\mathbf{B}_{48}$	=	$(-x_{13}+z_{13})\,\mathbf{a}_1+(x_{13}+z_{13})\,\mathbf{a}_2$	=	$x_{13}a\,\hat{\mathbf{x}}-x_{13}a\,\hat{\mathbf{y}}+z_{13}c\,\hat{\mathbf{z}}$	(16*m*)	Cl VI
$\mathbf{B}_{49}$	=	$(x_{13}-z_{13})\,\mathbf{a}_1+(-x_{13}-z_{13})\,\mathbf{a}_2$	=	$-x_{13}a\,\hat{\mathbf{x}}+x_{13}a\,\hat{\mathbf{y}}-z_{13}c\,\hat{\mathbf{z}}$	(16*m*)	Cl VI
$\mathbf{B}_{50}$	=	$(-x_{13}-z_{13})\,\mathbf{a}_1+(x_{13}-z_{13})\,\mathbf{a}_2$	=	$x_{13}a\,\hat{\mathbf{x}}-x_{13}a\,\hat{\mathbf{y}}-z_{13}c\,\hat{\mathbf{z}}$	(16*m*)	Cl VI
$\mathbf{B}_{51}$	=	$(x_{13}-z_{13})\,\mathbf{a}_1+(x_{13}-z_{13})\,\mathbf{a}_2+2x_{13}\,\mathbf{a}_3$	=	$x_{13}a\,\hat{\mathbf{x}}+x_{13}a\,\hat{\mathbf{y}}-z_{13}c\,\hat{\mathbf{z}}$	(16*m*)	Cl VI
$\mathbf{B}_{52}$	=	$(-x_{13}-z_{13})\,\mathbf{a}_1+(-x_{13}-z_{13})\,\mathbf{a}_2-2x_{13}\,\mathbf{a}_3$	=	$-x_{13}a\,\hat{\mathbf{x}}-x_{13}a\,\hat{\mathbf{y}}-z_{13}c\,\hat{\mathbf{z}}$	(16*m*)	Cl VI
$\mathbf{B}_{53}$	=	$(x_{14}+z_{14})\,\mathbf{a}_1+(x_{14}+z_{14})\,\mathbf{a}_2+2x_{14}\,\mathbf{a}_3$	=	$x_{14}a\,\hat{\mathbf{x}}+x_{14}a\,\hat{\mathbf{y}}+z_{14}c\,\hat{\mathbf{z}}$	(16*m*)	Cl VII
$\mathbf{B}_{54}$	=	$(-x_{14}+z_{14})\,\mathbf{a}_1+(-x_{14}+z_{14})\,\mathbf{a}_2-2x_{14}\,\mathbf{a}_3$	=	$-x_{14}a\,\hat{\mathbf{x}}-x_{14}a\,\hat{\mathbf{y}}+z_{14}c\,\hat{\mathbf{z}}$	(16*m*)	Cl VII
$\mathbf{B}_{55}$	=	$(x_{14}+z_{14})\,\mathbf{a}_1+(-x_{14}+z_{14})\,\mathbf{a}_2$	=	$-x_{14}a\,\hat{\mathbf{x}}+x_{14}a\,\hat{\mathbf{y}}+z_{14}c\,\hat{\mathbf{z}}$	(16*m*)	Cl VII
$\mathbf{B}_{56}$	=	$(-x_{14}+z_{14})\,\mathbf{a}_1+(x_{14}+z_{14})\,\mathbf{a}_2$	=	$x_{14}a\,\hat{\mathbf{x}}-x_{14}a\,\hat{\mathbf{y}}+z_{14}c\,\hat{\mathbf{z}}$	(16*m*)	Cl VII
$\mathbf{B}_{57}$	=	$(x_{14}-z_{14})\,\mathbf{a}_1+(-x_{14}-z_{14})\,\mathbf{a}_2$	=	$-x_{14}a\,\hat{\mathbf{x}}+x_{14}a\,\hat{\mathbf{y}}-z_{14}c\,\hat{\mathbf{z}}$	(16*m*)	Cl VII
$\mathbf{B}_{58}$	=	$(-x_{14}-z_{14})\,\mathbf{a}_1+(x_{14}-z_{14})\,\mathbf{a}_2$	=	$x_{14}a\,\hat{\mathbf{x}}-x_{14}a\,\hat{\mathbf{y}}-z_{14}c\,\hat{\mathbf{z}}$	(16*m*)	Cl VII
$\mathbf{B}_{59}$	=	$(x_{14}-z_{14})\,\mathbf{a}_1+(x_{14}-z_{14})\,\mathbf{a}_2+2x_{14}\,\mathbf{a}_3$	=	$x_{14}a\,\hat{\mathbf{x}}+x_{14}a\,\hat{\mathbf{y}}-z_{14}c\,\hat{\mathbf{z}}$	(16*m*)	Cl VII
$\mathbf{B}_{60}$	=	$(-x_{14}-z_{14})\,\mathbf{a}_1+(-x_{14}-z_{14})\,\mathbf{a}_2-2x_{14}\,\mathbf{a}_3$	=	$-x_{14}a\,\hat{\mathbf{x}}-x_{14}a\,\hat{\mathbf{y}}-z_{14}c\,\hat{\mathbf{z}}$	(16*m*)	Cl VII
$\mathbf{B}_{61}$	=	$(y_{15}+z_{15})\,\mathbf{a}_1+z_{15}\,\mathbf{a}_2+y_{15}\,\mathbf{a}_3$	=	$y_{15}a\,\hat{\mathbf{y}}+z_{15}c\,\hat{\mathbf{z}}$	(16*n*)	H_2O III
$\mathbf{B}_{62}$	=	$(-y_{15}+z_{15})\,\mathbf{a}_1+z_{15}\,\mathbf{a}_2-y_{15}\,\mathbf{a}_3$	=	$-y_{15}a\,\hat{\mathbf{y}}+z_{15}c\,\hat{\mathbf{z}}$	(16*n*)	H_2O III
$\mathbf{B}_{63}$	=	$z_{15}\,\mathbf{a}_1+(-y_{15}+z_{15})\,\mathbf{a}_2-y_{15}\,\mathbf{a}_3$	=	$-y_{15}a\,\hat{\mathbf{x}}+z_{15}c\,\hat{\mathbf{z}}$	(16*n*)	H_2O III
$\mathbf{B}_{64}$	=	$z_{15}\,\mathbf{a}_1+(y_{15}+z_{15})\,\mathbf{a}_2+y_{15}\,\mathbf{a}_3$	=	$y_{15}a\,\hat{\mathbf{x}}+z_{15}c\,\hat{\mathbf{z}}$	(16*n*)	H_2O III
$\mathbf{B}_{65}$	=	$(y_{15}-z_{15})\,\mathbf{a}_1-z_{15}\,\mathbf{a}_2+y_{15}\,\mathbf{a}_3$	=	$y_{15}a\,\hat{\mathbf{y}}-z_{15}c\,\hat{\mathbf{z}}$	(16*n*)	H_2O III
$\mathbf{B}_{66}$	=	$(-y_{15}-z_{15})\,\mathbf{a}_1-z_{15}\,\mathbf{a}_2-y_{15}\,\mathbf{a}_3$	=	$-y_{15}a\,\hat{\mathbf{y}}-z_{15}c\,\hat{\mathbf{z}}$	(16*n*)	H_2O III
$\mathbf{B}_{67}$	=	$-z_{15}\,\mathbf{a}_1+(y_{15}-z_{15})\,\mathbf{a}_2+y_{15}\,\mathbf{a}_3$	=	$y_{15}a\,\hat{\mathbf{x}}-z_{15}c\,\hat{\mathbf{z}}$	(16*n*)	H_2O III

$\mathbf{B}_{68}$	=	$-z_{15}\,\mathbf{a}_1 + (-y_{15} - z_{15})\,\mathbf{a}_2 - y_{15}\,\mathbf{a}_3$	=	$-y_{15}a\,\hat{\mathbf{x}} - z_{15}c\,\hat{\mathbf{z}}$	$(16n)$	H_2O III
$\mathbf{B}_{69}$	=	$(y_{16} + z_{16})\,\mathbf{a}_1 + z_{16}\,\mathbf{a}_2 + y_{16}\,\mathbf{a}_3$	=	$y_{16}a\,\hat{\mathbf{y}} + z_{16}c\,\hat{\mathbf{z}}$	$(16n)$	K III
$\mathbf{B}_{70}$	=	$(-y_{16} + z_{16})\,\mathbf{a}_1 + z_{16}\,\mathbf{a}_2 - y_{16}\,\mathbf{a}_3$	=	$-y_{16}a\,\hat{\mathbf{y}} + z_{16}c\,\hat{\mathbf{z}}$	$(16n)$	K III
$\mathbf{B}_{71}$	=	$z_{16}\,\mathbf{a}_1 + (-y_{16} + z_{16})\,\mathbf{a}_2 - y_{16}\,\mathbf{a}_3$	=	$-y_{16}a\,\hat{\mathbf{x}} + z_{16}c\,\hat{\mathbf{z}}$	$(16n)$	K III
$\mathbf{B}_{72}$	=	$z_{16}\,\mathbf{a}_1 + (y_{16} + z_{16})\,\mathbf{a}_2 + y_{16}\,\mathbf{a}_3$	=	$y_{16}a\,\hat{\mathbf{x}} + z_{16}c\,\hat{\mathbf{z}}$	$(16n)$	K III
$\mathbf{B}_{73}$	=	$(y_{16} - z_{16})\,\mathbf{a}_1 - z_{16}\,\mathbf{a}_2 + y_{16}\,\mathbf{a}_3$	=	$y_{16}a\,\hat{\mathbf{y}} - z_{16}c\,\hat{\mathbf{z}}$	$(16n)$	K III
$\mathbf{B}_{74}$	=	$(-y_{16} - z_{16})\,\mathbf{a}_1 - z_{16}\,\mathbf{a}_2 - y_{16}\,\mathbf{a}_3$	=	$-y_{16}a\,\hat{\mathbf{y}} - z_{16}c\,\hat{\mathbf{z}}$	$(16n)$	K III
$\mathbf{B}_{75}$	=	$-z_{16}\,\mathbf{a}_1 + (y_{16} - z_{16})\,\mathbf{a}_2 + y_{16}\,\mathbf{a}_3$	=	$y_{16}a\,\hat{\mathbf{x}} - z_{16}c\,\hat{\mathbf{z}}$	$(16n)$	K III
$\mathbf{B}_{76}$	=	$-z_{16}\,\mathbf{a}_1 + (-y_{16} - z_{16})\,\mathbf{a}_2 - y_{16}\,\mathbf{a}_3$	=	$-y_{16}a\,\hat{\mathbf{x}} - z_{16}c\,\hat{\mathbf{z}}$	$(16n)$	K III
$\mathbf{B}_{77}$	=	$(y_{17} + z_{17})\,\mathbf{a}_1 + z_{17}\,\mathbf{a}_2 + y_{17}\,\mathbf{a}_3$	=	$y_{17}a\,\hat{\mathbf{y}} + z_{17}c\,\hat{\mathbf{z}}$	$(16n)$	K IV
$\mathbf{B}_{78}$	=	$(-y_{17} + z_{17})\,\mathbf{a}_1 + z_{17}\,\mathbf{a}_2 - y_{17}\,\mathbf{a}_3$	=	$-y_{17}a\,\hat{\mathbf{y}} + z_{17}c\,\hat{\mathbf{z}}$	$(16n)$	K IV
$\mathbf{B}_{79}$	=	$z_{17}\,\mathbf{a}_1 + (-y_{17} + z_{17})\,\mathbf{a}_2 - y_{17}\,\mathbf{a}_3$	=	$-y_{17}a\,\hat{\mathbf{x}} + z_{17}c\,\hat{\mathbf{z}}$	$(16n)$	K IV
$\mathbf{B}_{80}$	=	$z_{17}\,\mathbf{a}_1 + (y_{17} + z_{17})\,\mathbf{a}_2 + y_{17}\,\mathbf{a}_3$	=	$y_{17}a\,\hat{\mathbf{x}} + z_{17}c\,\hat{\mathbf{z}}$	$(16n)$	K IV
$\mathbf{B}_{81}$	=	$(y_{17} - z_{17})\,\mathbf{a}_1 - z_{17}\,\mathbf{a}_2 + y_{17}\,\mathbf{a}_3$	=	$y_{17}a\,\hat{\mathbf{y}} - z_{17}c\,\hat{\mathbf{z}}$	$(16n)$	K IV
$\mathbf{B}_{82}$	=	$(-y_{17} - z_{17})\,\mathbf{a}_1 - z_{17}\,\mathbf{a}_2 - y_{17}\,\mathbf{a}_3$	=	$-y_{17}a\,\hat{\mathbf{y}} - z_{17}c\,\hat{\mathbf{z}}$	$(16n)$	K IV
$\mathbf{B}_{83}$	=	$-z_{17}\,\mathbf{a}_1 + (y_{17} - z_{17})\,\mathbf{a}_2 + y_{17}\,\mathbf{a}_3$	=	$y_{17}a\,\hat{\mathbf{x}} - z_{17}c\,\hat{\mathbf{z}}$	$(16n)$	K IV
$\mathbf{B}_{84}$	=	$-z_{17}\,\mathbf{a}_1 + (-y_{17} - z_{17})\,\mathbf{a}_2 - y_{17}\,\mathbf{a}_3$	=	$-y_{17}a\,\hat{\mathbf{x}} - z_{17}c\,\hat{\mathbf{z}}$	$(16n)$	K IV

References:

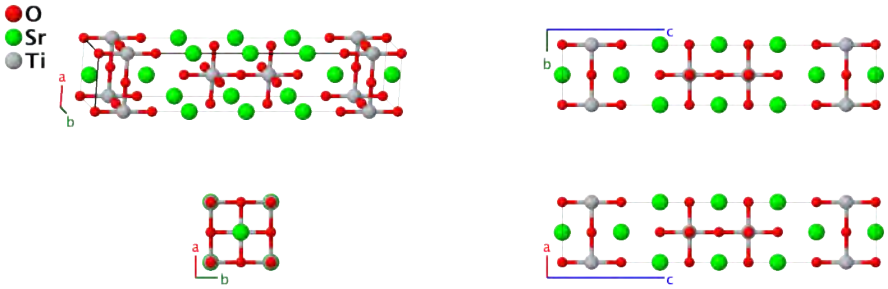
- J. L. Hoard and L. Goldstein, *The Structure of Potassium Hexachlorothalliate Dihydrate*, J. Chem. Phys. **3**, 645–649 (1935), doi:10.1063/1.1749568.

Geometry files:

- CIF: pp. 1708
- POSCAR: pp. 1708

$Sr_3Ti_2O_7$ Structure: A7B3C2_tI24_139_aeg_be_e

http://aflow.org/prototype-encyclopedia/A7B3C2_tI24_139_aeg_be_e



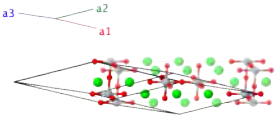
Prototype	:	$O_7Sr_3Ti_2$
AFLOW prototype label	:	A7B3C2_tI24_139_aeg_be_e
***Strukturbericht* designation**	:	None
Pearson symbol	:	tI24
Space group number	:	139
Space group symbol	:	$I4/mmm$
AFLOW prototype command	:	`aflow --proto=A7B3C2_tI24_139_aeg_be_e --params=`$a, c/a, z_3, z_4, z_5, z_6$

Other compounds with this structure

- $Ca_3Ti_2O_7$, $Sr_3Ru_2O_7$, $BaLa_2Fe_2O_7$, and $SrTb_2Fe_2O_7$

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}a\,\hat{\mathbf{x}} - \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} - \tfrac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	O I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	$(2b)$	Sr I
$\mathbf{B}_3$	=	$z_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_2$	=	$z_3 c\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_4$	=	$-z_3\,\mathbf{a}_1 - z_3\,\mathbf{a}_2$	=	$-z_3 c\,\hat{\mathbf{z}}$	$(4e)$	O II
$\mathbf{B}_5$	=	$z_4\,\mathbf{a}_1 + z_4\,\mathbf{a}_2$	=	$z_4 c\,\hat{\mathbf{z}}$	$(4e)$	Sr II
$\mathbf{B}_6$	=	$-z_4\,\mathbf{a}_1 - z_4\,\mathbf{a}_2$	=	$-z_4 c\,\hat{\mathbf{z}}$	$(4e)$	Sr II
$\mathbf{B}_7$	=	$z_5\,\mathbf{a}_1 + z_5\,\mathbf{a}_2$	=	$z_5 c\,\hat{\mathbf{z}}$	$(4e)$	Ti
$\mathbf{B}_8$	=	$-z_5\,\mathbf{a}_1 - z_5\,\mathbf{a}_2$	=	$-z_5 c\,\hat{\mathbf{z}}$	$(4e)$	Ti
$\mathbf{B}_9$	=	$\left(\frac{1}{2} + z_6\right)\mathbf{a}_1 + z_6\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(8g)$	O III
$\mathbf{B}_{10}$	=	$z_6\,\mathbf{a}_1 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + z_6 c\,\hat{\mathbf{z}}$	$(8g)$	O III

$\mathbf{B}_{11}$	=	$\left(\frac{1}{2} - z_6\right)\mathbf{a}_1 - z_6\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	($8g$)	O III
$\mathbf{B}_{12}$	=	$-z_6\,\mathbf{a}_1 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - z_6 c\,\hat{\mathbf{z}}$	($8g$)	O III

References:

- S. N. Ruddlesden and P. Popper, *The compound $Sr_3Ti_2O_7$ and its structure*, Acta Cryst. **11**, 54–55 (1958), doi:10.1107/S0365110X58000128.

Found in:

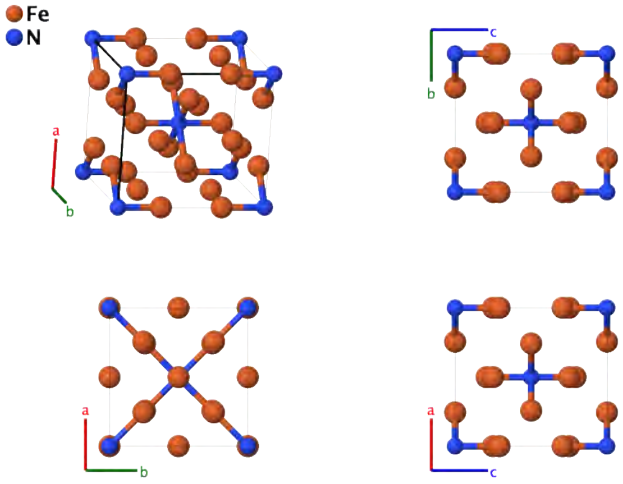
- Wikipedia, *Ruddlesden-Popper phase*, https://en.wikipedia.org/wiki/Ruddlesden-Popper_phase. $A_3B_2X_7$ series.

Geometry files:

- CIF: pp. 1709
- POSCAR: pp. 1709

Fe_8N ($D2_g$) Structure: A8B_tI18_139_deh_a

http://aflow.org/prototype-encyclopedia/A8B_tI18_139_deh_a

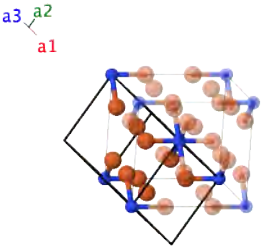


Prototype	:	Fe_8N
AFLOW prototype label	:	A8B_tI18_139_deh_a
***Strukturbericht* designation**	:	$D2_g$
Pearson symbol	:	tI18
Space group number	:	139
Space group symbol	:	$I4/mmm$
AFLOW prototype command	:	`aflow --proto=A8B_tI18_139_deh_a --params=`$a, c/a, z_3, x_4$

- (Jack, 1951) and others refer to this structure as α''-$Fe_{16}N_2$.
- (Yamashita, 2012) determined the structure by examining crystals containing various amounts of α''-$Fe_{16}N_2$ and α-Fe (bcc iron). We use their data from the sample that was 90% α''-$Fe_{16}N_2$. Their measurements were confirmed by first principles calculations, and are in agreement with the first principles calculations of (Sims, 2012).

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}a\,\hat{\mathbf{x}} - \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} - \tfrac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	N
$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(4d)$	Fe I

$\mathbf{B}_3$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*d*)	Fe I
$\mathbf{B}_4$	=	$z_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_2$	=	$z_3 c\,\hat{\mathbf{z}}$	(4*e*)	Fe II
$\mathbf{B}_5$	=	$-z_3\,\mathbf{a}_1 - z_3\,\mathbf{a}_2$	=	$-z_3 c\,\hat{\mathbf{z}}$	(4*e*)	Fe II
$\mathbf{B}_6$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + 2x_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}}$	(8*h*)	Fe III
$\mathbf{B}_7$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - 2x_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}}$	(8*h*)	Fe III
$\mathbf{B}_8$	=	$x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2$	=	$-x_4 a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}}$	(8*h*)	Fe III
$\mathbf{B}_9$	=	$-x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2$	=	$x_4 a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}}$	(8*h*)	Fe III

References:

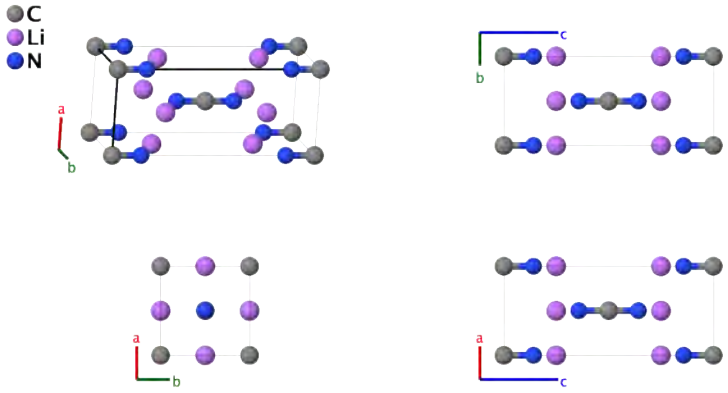
- S. Yamashita, Y. Masubuchi, Y. Nakazawa, T. Okayama, M. Tsuchiya, and S. Kikkawa, *Crystal structure and magnetic properties of "α″-$Fe_{16}N_2$" containing reisdual α-Fe prepared by low-temperature ammonia nitridation*, J. Solid State Chem. **194**, 76–79 (2012), doi:10.1016/j.jssc.2012.07.025.
- K. H. Jack, *The occurrence and the crystal structure of α″-iron nitride; a new type of interstitial alloy formed during the tempering of nitrogen-martensite*, Proc. Roy. Soc. Lond. A **208**, 216–224 (1951), doi:10.1098/rspa.1951.0155.
- H. Sims, W. H. Butler, M. Richter, K. Koepernik, E. Şaşıoğlu, C. Friedrich, and S. Blügel, *Theoretical investigation into the possibility of very large moments in $Fe_{16}N_2$*, Phys. Rev. B **86**, 174422 (2012), doi:10.1103/PhysRevB.86.174422.

Geometry files:

- CIF: pp. 1709
- POSCAR: pp. 1710

Li_2CN_2 Structure: AB2C2_tI10_139_a_d_e

http://aflow.org/prototype-encyclopedia/AB2C2_tI10_139_a_d_e.Li2CN2

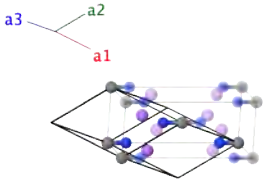


Prototype	:	CLi_2N_2
AFLOW prototype label	:	AB2C2_tI10_139_a_d_e
***Strukturbericht* designation**	:	None
Pearson symbol	:	tI10
Space group number	:	139
Space group symbol	:	$I4/mmm$
AFLOW prototype command	:	`aflow --proto=AB2C2_tI10_139_a_d_e --params=`$a, c/a, z_3$

- This structure has the same AFLOW prototype label, AB2C2_tI10_139_a_d_e, as autunite, $Ca(UO_2)_2(PO_4)_2{\cdot}10.5H_2O$ ($H5_9$). They are generated by the same symmetry operations with different sets of parameters (`--params`) specified in their corresponding CIF files.

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}a\,\hat{\mathbf{x}} - \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} - \tfrac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	C
$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(4d)$	Li
$\mathbf{B}_3$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(4d)$	Li
$\mathbf{B}_4$	=	$z_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_2$	=	$z_3 c\,\hat{\mathbf{z}}$	$(4e)$	N
$\mathbf{B}_5$	=	$-z_3\,\mathbf{a}_1 - z_3\,\mathbf{a}_2$	=	$-z_3 c\,\hat{\mathbf{z}}$	$(4e)$	N

References:

- M. G. Down, M. J. Haley, P. Hubberstey, R. J. Pulham, and A. E. Thunder, *Solutions of lithium salts in liquid lithium: preparation and X-ray crystal structure of the dilithium salt of carbodi-imide (cyanamide)*, J. Chem. Soc., Dalton Trans. pp. 1407–1411 (1978), doi:10.1039/DT9780001407.

Found in:

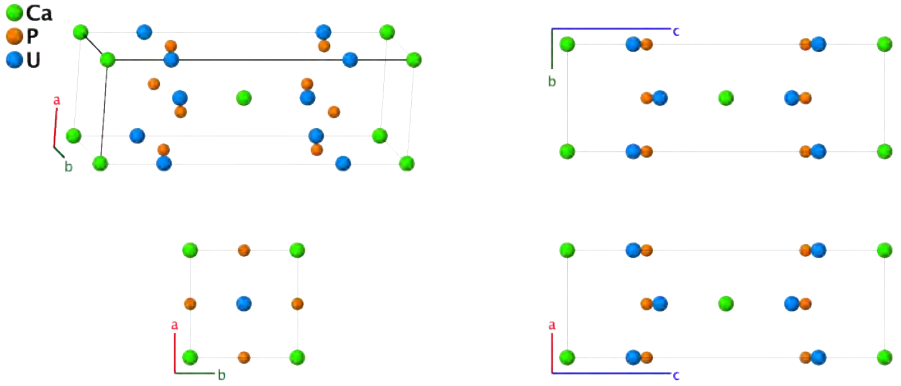
- A. Jain, S. P. Ong, G. Hautier, W. Chen, W. D. Richards, S. Dacek, S. Cholia, D. Gunter, D. Skinner, G. Ceder, and K. A. Persson, *Commentary: The Materials Project: A materials genome approach to accelerating materials innovation*, APL Mater. **1**, 011002 (2013), doi:10.1063/1.4812323.

Geometry files:

- CIF: pp. 1710
- POSCAR: pp. 1710

$H5_9$ [Autunite, $Ca(UO_2)_2(PO_4)_2{\cdot}10\frac{1}{2}H_2O$] (*obsolete*) Structure: AB2C2_tI10_139_a_d_e

http://aflow.org/prototype-encyclopedia/AB2C2_tI10_139_a_d_e

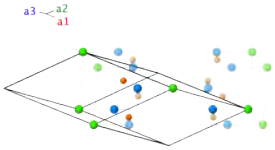


Prototype	:	$Ca(UO_2)_2(PO_4)_2{\cdot}10\frac{1}{2}H_2O$
AFLOW prototype label	:	AB2C2_tI10_139_a_d_e
***Strukturbericht* designation**	:	$H5_9$
Pearson symbol	:	tI10
Space group number	:	139
Space group symbol	:	$I4/mmm$
AFLOW prototype command	:	`aflow --proto=AB2C2_tI10_139_a_d_e --params=`$a, c/a, z_3$

- Autunite $Ca(UO_2)_2(PO_4)_2{\cdot}nH_2O$, is found in three varieties: naturally occurring Autunite, with $n \gtrsim 10$, and meta-autunite (I), which is partially dehydrated, $6 \gtrsim n \gtrsim 10$. Further dehydration in the laboratory produces meta-autunite (II).
- (Beintema, 1938) proposed this structure for autunite, which (Herrmann, 1941) designated $H5_9$. This original tetragonal structure did not locate the oxygen atoms or the water molecules. Later determinations of this structure (Locock, 2003) located all the atoms in the crystal for $n = 11$, and found the structure to be orthorhombic.

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}a\,\hat{\mathbf{x}} - \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} - \tfrac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	Ca
$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(4d)$	P
$\mathbf{B}_3$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(4d)$	P

$\mathbf{B}_4$	=	$z_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_2$	=	$z_3 c\,\hat{\mathbf{z}}$	($4e$)	U
$\mathbf{B}_5$	=	$-z_3\,\mathbf{a}_1 - z_3\,\mathbf{a}_2$	=	$-z_3 c\,\hat{\mathbf{z}}$	($4e$)	U

References:

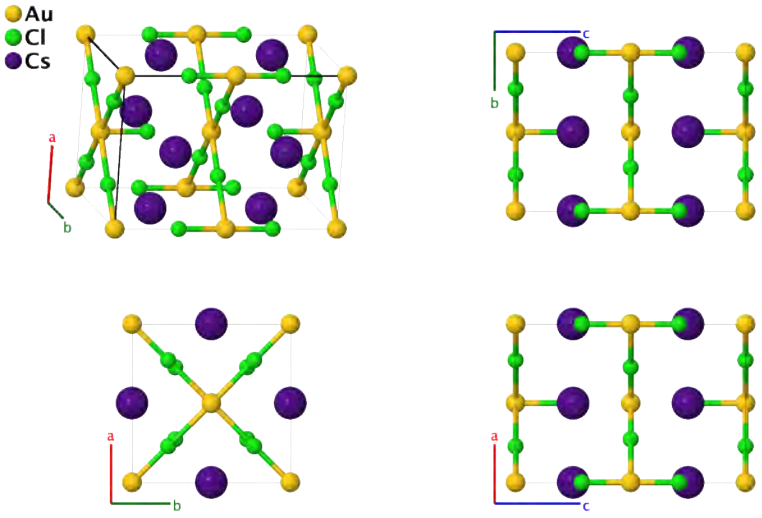
- J. Beintema, *On the composition and the crystallography of autunite and the meta-autunites*, Rec. Trav. Chim. Pays-Bas **57**, 155–175 (1938), doi:10.1002/recl.19380570206.
- A. J. Locock and P. C. Burns, *The crystal structure of synthetic autunite, $Ca[(UO_2)(PO_4)]_2(H_2O)_{11}$*, Am. Mineral. **88**, 240–244 (2003), doi:10.2138/am-2003-0128.

Geometry files:

- CIF: pp. 1710
- POSCAR: pp. 1711

$AuCsCl_3$ ($K7_6$) Structure: AB3C_tI20_139_ab_eh_d

http://aflow.org/prototype-encyclopedia/AB3C_tI20_139_ab_eh_d



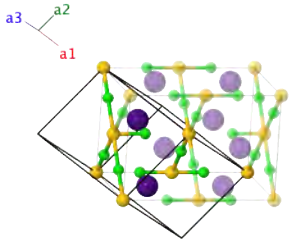
Prototype	:	$AuCl_3Cs$
AFLOW prototype label	:	AB3C_tI20_139_ab_eh_d
***Strukturbericht* designation**	:	$K7_6$
Pearson symbol	:	tI20
Space group number	:	139
Space group symbol	:	$I4/mmm$
AFLOW prototype command	:	`aflow --proto=AB3C_tI20_139_ab_eh_d --params=`$a, c/a, z_4, x_5$

Other compounds with this structure

- $AgAuCs_2Cl_6$

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}a\,\hat{\mathbf{x}} - \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} - \tfrac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($2a$)	Au I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	($2b$)	Au II
$\mathbf{B}_3$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($4d$)	Cs
$\mathbf{B}_4$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($4d$)	Cs

$\mathbf{B}_5$	=	$z_4\,\mathbf{a}_1 + z_4\,\mathbf{a}_2$	=	$z_4 c\,\hat{\mathbf{z}}$	($4e$)	Cl I
$\mathbf{B}_6$	=	$-z_4\,\mathbf{a}_1 - z_4\,\mathbf{a}_2$	=	$-z_4 c\,\hat{\mathbf{z}}$	($4e$)	Cl I
$\mathbf{B}_7$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + 2x_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}}$	($8h$)	Cl II
$\mathbf{B}_8$	=	$-x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - 2x_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}}$	($8h$)	Cl II
$\mathbf{B}_9$	=	$x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2$	=	$-x_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}}$	($8h$)	Cl II
$\mathbf{B}_{10}$	=	$-x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2$	=	$x_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}}$	($8h$)	Cl II

References:

- N. Elliott and L. Pauling, *The Crystal Structure of Cesium Aurous Auric Chloride, $Cs_2AuAuCl_6$, and Cesium Argentous Auric Chloride, $Cs_2AgAuCl_6$*, J. Am. Chem. Soc. **60**, 1846–1851 (1938), doi:10.1021/ja01275a037.

Found in:

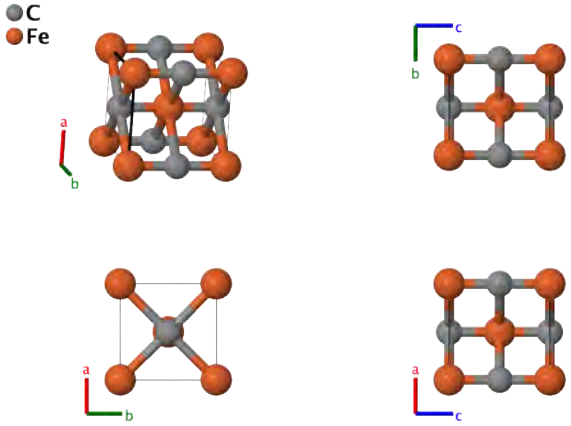
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1711
- POSCAR: pp. 1711

"Martensite Type" FeC_x ($x \leq 0.06$) ($L'2_0$) Structure: AB_tI4_139_b_a

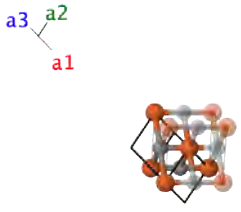
http://aflow.org/prototype-encyclopedia/AB_tI4_139_b_a



Prototype	:	C_xFe
AFLOW prototype label	:	AB_tI4_139_b_a
***Strukturbericht* designation**	:	$L'2_0$
Pearson symbol	:	tI4
Space group number	:	139
Space group symbol	:	$I4/mmm$
AFLOW prototype command	:	`aflow --proto=AB_tI4_139_b_a --params=`$a, c/a$

- This should *not* be considered a true model of martensite, a very complex material whose structure is sensitive to composition (Blackwell, 1996; Sherby, 2008). However, (Ewald, 1931) assign this the label $L'2_0$, saying that this structure is "α-Fe with 6 atomic percent C." Although Ewald and Hermann present a long discussion of martensite (pp. 489, 576-583), they do not give this exact crystal structure, which we take from (Brandes, 1992) and (Pearson, 1967).
- (Villars, 1995) and (Westbrook, 1995) ignore this structure and give the $L'2$ label to the ThH_2 structure. We previously followed this notation, but in the interest of historical accuracy will relabel that structure with the name given to it by (Brandes, 1992) and (Pearson, 1967), $L'2_b$.
- We should note that these last two references list the current structure simply as $L'2$.

Body-centered Tetragonal primitive vectors:



$$\mathbf{a}_1 = -\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} - \frac{1}{2}c\,\hat{\mathbf{z}}$$

Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($2a$)	Fe
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	($2b$)	C

References:

- E. A. Brandes and G. B. Brook, eds., *Smithells Metals Reference Book* (Butterworth Heinemann, 1992), chap. 6, pp. 6–63, 7e edn.
- R. Blackwell, *Internal Friction Effects in Tempered Martensite*, Nature **211**, 733–734 (1966), doi:10.1038/211733a0.
- O. D. Sherby, J. Wadsworth, D. R. Lesuer, and C. K. Syn, *Revisiting the Structure of Martensite in Iron-Carbon Steels*, Mater. Trans. **49**, 2016–2027 (2008), doi:10.2320/matertrans.MRA2007338.
- P. Villars and L. Calvert, *Pearson's Handbook of Crystallographic Data for Intermetallic Phases* (ASM International, Materials Park, OH, 1991), 2nd edn.
- J. H. Westbrook and R. L. Fleischer, eds., *Intermetallic Compounds – Principles and Practice* (John Wiley & Sons, Ltd., Chichester, England, 1995). Two Volumes.
- W. B. Pearson, *A Handbook of Lattice Spacings and Structures of Metals and Alloys, Volume 2*, *International Series of Monographs on Metal Physics and Physical Metallurgy*, vol. 8 (Pergamon Press, Oxford, London, Edinburgh, New York, Toronto, Sydney, Paris, Braunschweig, 1967). N. R. C. No. 8752.

Found in:

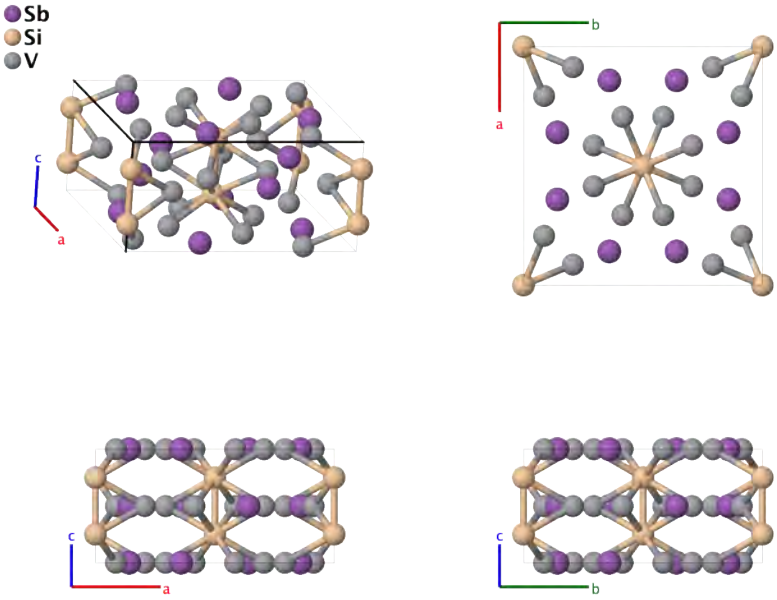
- P. P. Ewald and C. Hermann, eds., *Strukturbericht 1913-1928* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1931).

Geometry files:

- CIF: pp. 1711
- POSCAR: pp. 1712

V_4SiSb_2 Structure: A2BC4_tI28_140_h_a_k

http://aflow.org/prototype-encyclopedia/A2BC4_tI28_140_h_a_k



Prototype	:	Sb_2SiV_4
AFLOW prototype label	:	A2BC4_tI28_140_h_a_k
***Strukturbericht* designation**	:	None
Pearson symbol	:	tI28
Space group number	:	140
Space group symbol	:	$I4/mcm$
AFLOW prototype command	:	`aflow --proto=A2BC4_tI28_140_h_a_k --params=`$a, c/a, x_2, x_3, y_3$

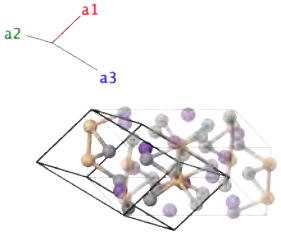
Other compounds with this structure

- Ti_4CoBi_2, Ti_4CrBi_2, Ti_4FeBi_2, Ti_4MnBi_2, and Ti_4NiBi_2

- This is the ternary version of the $D2_c$ U_6Mn structure. This can also be identified as a defected version of the $D8_m$ W_5Si_3 structure.

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{y}} + \tfrac{1}{2}\,c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2}\,a\,\hat{\mathbf{y}} + \tfrac{1}{2}\,c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{y}} - \tfrac{1}{2}\,c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2$	=	$\frac{1}{4}c\,\hat{\mathbf{z}}$	(4*a*)	Si
$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2$	=	$\frac{3}{4}c\,\hat{\mathbf{z}}$	(4*a*)	Si
$\mathbf{B}_3$	=	$\left(\frac{1}{2}+x_2\right)\mathbf{a}_1 + x_2\,\mathbf{a}_2 + \left(\frac{1}{2}+2x_2\right)\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_2\right)a\,\hat{\mathbf{y}}$	(8*h*)	Sb
$\mathbf{B}_4$	=	$\left(\frac{1}{2}-x_2\right)\mathbf{a}_1 - x_2\,\mathbf{a}_2 + \left(\frac{1}{2}-2x_2\right)\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_2\right)a\,\hat{\mathbf{y}}$	(8*h*)	Sb
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2}-x_2\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_2\right)a\,\hat{\mathbf{x}} + x_2 a\,\hat{\mathbf{y}}$	(8*h*)	Sb
$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2}+x_2\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_2\right)a\,\hat{\mathbf{x}} - x_2 a\,\hat{\mathbf{y}}$	(8*h*)	Sb
$\mathbf{B}_7$	=	$y_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + (x_3+y_3)\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 a\,\hat{\mathbf{y}}$	(16*k*)	V
$\mathbf{B}_8$	=	$-y_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + (-x_3-y_3)\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 a\,\hat{\mathbf{y}}$	(16*k*)	V
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + (x_3-y_3)\,\mathbf{a}_3$	=	$-y_3 a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}}$	(16*k*)	V
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + (-x_3+y_3)\,\mathbf{a}_3$	=	$y_3 a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}}$	(16*k*)	V
$\mathbf{B}_{11}$	=	$\left(\frac{1}{2}+y_3\right)\mathbf{a}_1 + \left(\frac{1}{2}-x_3\right)\mathbf{a}_2 + (-x_3+y_3)\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + y_3 a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(16*k*)	V
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2}-y_3\right)\mathbf{a}_1 + \left(\frac{1}{2}+x_3\right)\mathbf{a}_2 + (x_3-y_3)\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} - y_3 a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(16*k*)	V
$\mathbf{B}_{13}$	=	$\left(\frac{1}{2}+x_3\right)\mathbf{a}_1 + \left(\frac{1}{2}+y_3\right)\mathbf{a}_2 + (x_3+y_3)\,\mathbf{a}_3$	=	$y_3 a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(16*k*)	V
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2}-x_3\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_3\right)\mathbf{a}_2 + (-x_3-y_3)\,\mathbf{a}_3$	=	$-y_3 a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(16*k*)	V

References:

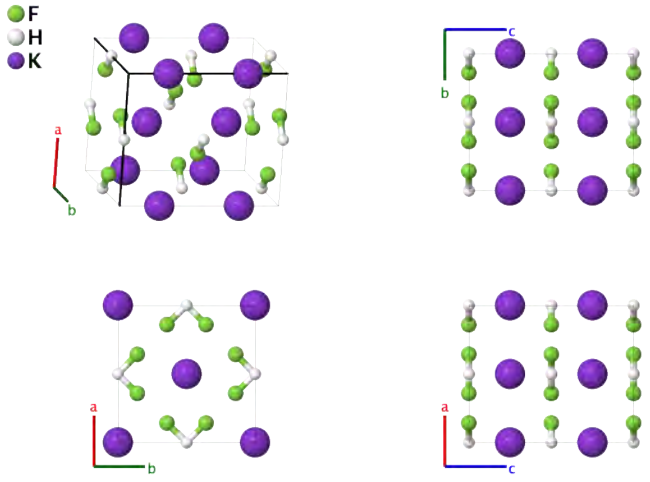
- P. Wollesen and W. Jeitschko, *V_4SiSb_2, a vanadium silicide antimonide crystallizing with a defect variant of the W_5Si_3-type structure*, J. Alloys Compd. **243**, 67–69 (1996), doi:10.1016/S0925-8388(96)02397-3.

Geometry files:

- CIF: pp. 1712
- POSCAR: pp. 1712

KHF_2 ($F5_2$) Structure: A2BC_tI16_140_h_d_a

http://aflow.org/prototype-encyclopedia/A2BC_tI16_140_h_d_a



Prototype	:	F_2HK
AFLOW prototype label	:	A2BC_tI16_140_h_d_a
***Strukturbericht* designation**	:	$F5_2$
Pearson symbol	:	tI16
Space group number	:	140
Space group symbol	:	$I4/mcm$
AFLOW prototype command	:	`aflow --proto=A2BC_tI16_140_h_d_a --params=`$a, c/a, x_3$

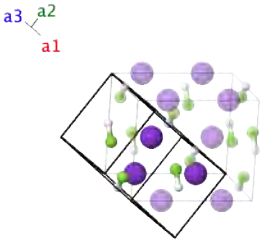
Other compounds with this structure

- KN_3 and CsN_3

- (Bozorth, 1923) originally determined the lattice constants of KHF_2 along with the positions of the potassium and fluorine atoms. He also assumed that the hydrogen atoms were on the (4*d*) Wyckoff sites. Both (Peterson, 1952) and (Ibers, 1964) confirmed his data. All three papers point out that it is possible that the hydrogen atoms are on half-filled (8*h*) sites, which would reduce to the (4*d*) site as the x coordinate approached zero. In KN_3 and CsN_3 the nitrogen atoms occupy the (4*d*) and (8*h*) Wyckoff sites, while the cation occupies the (4*a*) site.

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} - \frac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2$	=	$\frac{1}{4}c\,\hat{\mathbf{z}}$	(4*a*)	K
$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2$	=	$\frac{3}{4}c\,\hat{\mathbf{z}}$	(4*a*)	K
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}}$	(4*d*)	H
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	(4*d*)	H
$\mathbf{B}_5$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \left(\frac{1}{2} + 2x_3\right)\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}}$	(8*h*)	F
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\frac{1}{2} - 2x_3\right)\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}}$	(8*h*)	F
$\mathbf{B}_7$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}}$	(8*h*)	F
$\mathbf{B}_8$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}}$	(8*h*)	F

References:

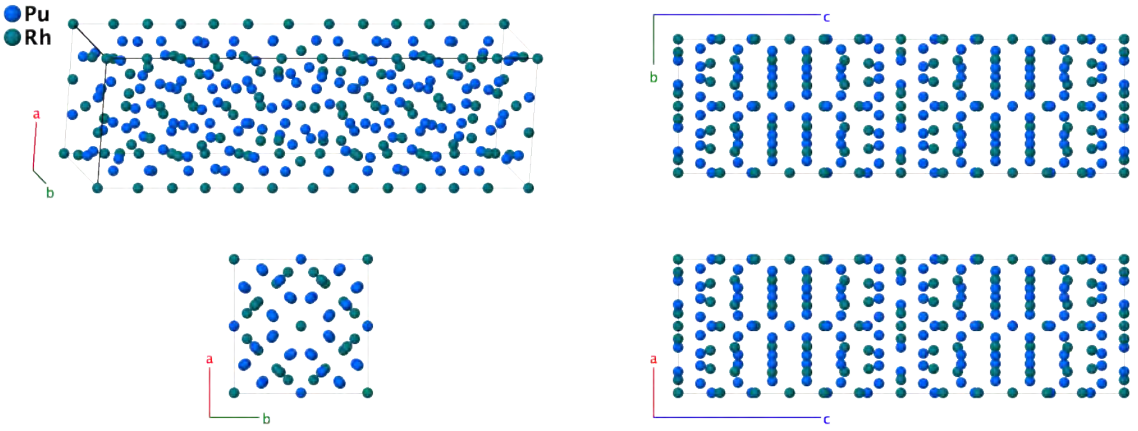
- J. A. Ibers, *Refinement of Peterson and Levy's Neutron Diffraction Data on KHF_2*, J. Chem. Phys. **40**, 402–404 (1964), doi:10.1063/1.1725126.
- S. W. Peterson and H. A. Levy, *A Single Crystal Neutron Diffraction Determination of the Hydrogen Position in Potassium Bifluoride*, J. Chem. Phys. **20**, 704–707 (1952), doi:10.1063/1.1700520.
- R. M. Bozorth, *The crystal structure of potassium hydrogen fluoride*, J. Am. Chem. Soc. **45**, 2128–2132 (1923), doi:10.1021/ja01662a013.

Geometry files:

- CIF: pp. 1712
- POSCAR: pp. 1713

$Pu_{31}Rh_{20}$ Structure: A31B20_tI204_140_b2gh3m_ac2fh3l

http://aflow.org/prototype-encyclopedia/A31B20_tI204_140_b2gh3m_ac2fh3l



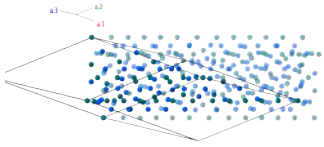
Prototype	:	$Pu_{31}Rh_{20}$
AFLOW prototype label	:	A31B20_tI204_140_b2gh3m_ac2fh3l
***Strukturbericht* designation**	:	None
Pearson symbol	:	tI204
Space group number	:	140
Space group symbol	:	$I4/mcm$
AFLOW prototype command	:	`aflow --proto=A31B20_tI204_140_b2gh3m_ac2fh3l --params=`$a, c/a, z_4, z_5, z_6, z_7, x_8, x_9, x_{10}, z_{10}, x_{11}, z_{11}, x_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}$

Other compounds with this structure

- $Pu_{31}Pt_{20}$ and $Ca_{31}Sn_{20}$

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}a\,\hat{\mathbf{x}} - \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} - \tfrac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\tfrac{1}{4}\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2$	=	$\tfrac{1}{4}c\,\hat{\mathbf{z}}$	($4a$)	Rh I
$\mathbf{B}_2$	=	$\tfrac{3}{4}\mathbf{a}_1 + \tfrac{3}{4}\mathbf{a}_2$	=	$\tfrac{3}{4}c\,\hat{\mathbf{z}}$	($4a$)	Rh I
$\mathbf{B}_3$	=	$\tfrac{3}{4}\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2 + \tfrac{1}{2}\mathbf{a}_3$	=	$\tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{4}c\,\hat{\mathbf{z}}$	($4b$)	Pu I
$\mathbf{B}_4$	=	$\tfrac{1}{4}\mathbf{a}_1 + \tfrac{3}{4}\mathbf{a}_2 + \tfrac{1}{2}\mathbf{a}_3$	=	$\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{4}c\,\hat{\mathbf{z}}$	($4b$)	Pu I
$\mathbf{B}_5$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($4c$)	Rh II
$\mathbf{B}_6$	=	$\tfrac{1}{2}\mathbf{a}_1 + \tfrac{1}{2}\mathbf{a}_2$	=	$\tfrac{1}{2}c\,\hat{\mathbf{z}}$	($4c$)	Rh II
$\mathbf{B}_7$	=	$z_4\,\mathbf{a}_1 + z_4\,\mathbf{a}_2$	=	$z_4 c\,\hat{\mathbf{z}}$	($8f$)	Rh III

$\mathbf{B}_8$	=	$\left(\frac{1}{2}-z_4\right)\mathbf{a}_1+\left(\frac{1}{2}-z_4\right)\mathbf{a}_2$	=	$\left(\frac{1}{2}-z_4\right)c\,\hat{\mathbf{z}}$	$(8f)$	Rh III
$\mathbf{B}_9$	=	$-z_4\,\mathbf{a}_1-z_4\,\mathbf{a}_2$	=	$-z_4c\,\hat{\mathbf{z}}$	$(8f)$	Rh III
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2}+z_4\right)\mathbf{a}_1+\left(\frac{1}{2}+z_4\right)\mathbf{a}_2$	=	$\left(\frac{1}{2}+z_4\right)c\,\hat{\mathbf{z}}$	$(8f)$	Rh III
$\mathbf{B}_{11}$	=	$z_5\,\mathbf{a}_1+z_5\,\mathbf{a}_2$	=	$z_5c\,\hat{\mathbf{z}}$	$(8f)$	Rh IV
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2}-z_5\right)\mathbf{a}_1+\left(\frac{1}{2}-z_5\right)\mathbf{a}_2$	=	$\left(\frac{1}{2}-z_5\right)c\,\hat{\mathbf{z}}$	$(8f)$	Rh IV
$\mathbf{B}_{13}$	=	$-z_5\,\mathbf{a}_1-z_5\,\mathbf{a}_2$	=	$-z_5c\,\hat{\mathbf{z}}$	$(8f)$	Rh IV
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2}+z_5\right)\mathbf{a}_1+\left(\frac{1}{2}+z_5\right)\mathbf{a}_2$	=	$\left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	$(8f)$	Rh IV
$\mathbf{B}_{15}$	=	$\left(\frac{1}{2}+z_6\right)\mathbf{a}_1+z_6\,\mathbf{a}_2+\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}}+z_6c\,\hat{\mathbf{z}}$	$(8g)$	Pu II
$\mathbf{B}_{16}$	=	$z_6\,\mathbf{a}_1+\left(\frac{1}{2}+z_6\right)\mathbf{a}_2+\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}+z_6c\,\hat{\mathbf{z}}$	$(8g)$	Pu II
$\mathbf{B}_{17}$	=	$-z_6\,\mathbf{a}_1+\left(\frac{1}{2}-z_6\right)\mathbf{a}_2+\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}-z_6c\,\hat{\mathbf{z}}$	$(8g)$	Pu II
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2}-z_6\right)\mathbf{a}_1-z_6\,\mathbf{a}_2+\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}}-z_6c\,\hat{\mathbf{z}}$	$(8g)$	Pu II
$\mathbf{B}_{19}$	=	$\left(\frac{1}{2}+z_7\right)\mathbf{a}_1+z_7\,\mathbf{a}_2+\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}}+z_7c\,\hat{\mathbf{z}}$	$(8g)$	Pu III
$\mathbf{B}_{20}$	=	$z_7\,\mathbf{a}_1+\left(\frac{1}{2}+z_7\right)\mathbf{a}_2+\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}+z_7c\,\hat{\mathbf{z}}$	$(8g)$	Pu III
$\mathbf{B}_{21}$	=	$-z_7\,\mathbf{a}_1+\left(\frac{1}{2}-z_7\right)\mathbf{a}_2+\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}-z_7c\,\hat{\mathbf{z}}$	$(8g)$	Pu III
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2}-z_7\right)\mathbf{a}_1-z_7\,\mathbf{a}_2+\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}}-z_7c\,\hat{\mathbf{z}}$	$(8g)$	Pu III
$\mathbf{B}_{23}$	=	$\left(\frac{1}{2}+x_8\right)\mathbf{a}_1+x_8\,\mathbf{a}_2+\left(\frac{1}{2}+2x_8\right)\mathbf{a}_3$	=	$x_8a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_8\right)a\,\hat{\mathbf{y}}$	$(8h)$	Pu IV
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2}-x_8\right)\mathbf{a}_1-x_8\,\mathbf{a}_2+\left(\frac{1}{2}-2x_8\right)\mathbf{a}_3$	=	$-x_8a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_8\right)a\,\hat{\mathbf{y}}$	$(8h)$	Pu IV
$\mathbf{B}_{25}$	=	$x_8\,\mathbf{a}_1+\left(\frac{1}{2}-x_8\right)\mathbf{a}_2+\frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_8\right)a\,\hat{\mathbf{x}}+x_8a\,\hat{\mathbf{y}}$	$(8h)$	Pu IV
$\mathbf{B}_{26}$	=	$-x_8\,\mathbf{a}_1+\left(\frac{1}{2}+x_8\right)\mathbf{a}_2+\frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_8\right)a\,\hat{\mathbf{x}}-x_8a\,\hat{\mathbf{y}}$	$(8h)$	Pu IV
$\mathbf{B}_{27}$	=	$\left(\frac{1}{2}+x_9\right)\mathbf{a}_1+x_9\,\mathbf{a}_2+\left(\frac{1}{2}+2x_9\right)\mathbf{a}_3$	=	$x_9a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_9\right)a\,\hat{\mathbf{y}}$	$(8h)$	Rh V
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2}-x_9\right)\mathbf{a}_1-x_9\,\mathbf{a}_2+\left(\frac{1}{2}-2x_9\right)\mathbf{a}_3$	=	$-x_9a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_9\right)a\,\hat{\mathbf{y}}$	$(8h)$	Rh V
$\mathbf{B}_{29}$	=	$x_9\,\mathbf{a}_1+\left(\frac{1}{2}-x_9\right)\mathbf{a}_2+\frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_9\right)a\,\hat{\mathbf{x}}+x_9a\,\hat{\mathbf{y}}$	$(8h)$	Rh V
$\mathbf{B}_{30}$	=	$-x_9\,\mathbf{a}_1+\left(\frac{1}{2}+x_9\right)\mathbf{a}_2+\frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_9\right)a\,\hat{\mathbf{x}}-x_9a\,\hat{\mathbf{y}}$	$(8h)$	Rh V
$\mathbf{B}_{31}$	=	$\left(\frac{1}{2}+x_{10}+z_{10}\right)\mathbf{a}_1+(x_{10}+z_{10})\,\mathbf{a}_2+\left(\frac{1}{2}+2x_{10}\right)\mathbf{a}_3$	=	$x_{10}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}}$	$(16l)$	Rh VI
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2}-x_{10}+z_{10}\right)\mathbf{a}_1+(-x_{10}+z_{10})\,\mathbf{a}_2+\left(\frac{1}{2}-2x_{10}\right)\mathbf{a}_3$	=	$-x_{10}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}}$	$(16l)$	Rh VI
$\mathbf{B}_{33}$	=	$(x_{10}+z_{10})\,\mathbf{a}_1+\left(\frac{1}{2}-x_{10}+z_{10}\right)\mathbf{a}_2+\frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{x}}+x_{10}a\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}}$	$(16l)$	Rh VI
$\mathbf{B}_{34}$	=	$(-x_{10}+z_{10})\,\mathbf{a}_1+\left(\frac{1}{2}+x_{10}+z_{10}\right)\mathbf{a}_2+\frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{x}}-x_{10}a\,\hat{\mathbf{y}}+z_{10}c\,\hat{\mathbf{z}}$	$(16l)$	Rh VI
$\mathbf{B}_{35}$	=	$(x_{10}-z_{10})\,\mathbf{a}_1+\left(\frac{1}{2}-x_{10}-z_{10}\right)\mathbf{a}_2+\frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{x}}+x_{10}a\,\hat{\mathbf{y}}-z_{10}c\,\hat{\mathbf{z}}$	$(16l)$	Rh VI
$\mathbf{B}_{36}$	=	$(-x_{10}-z_{10})\,\mathbf{a}_1+\left(\frac{1}{2}+x_{10}-z_{10}\right)\mathbf{a}_2+\frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{x}}-x_{10}a\,\hat{\mathbf{y}}-z_{10}c\,\hat{\mathbf{z}}$	$(16l)$	Rh VI
$\mathbf{B}_{37}$	=	$\left(\frac{1}{2}+x_{10}-z_{10}\right)\mathbf{a}_1+(x_{10}-z_{10})\,\mathbf{a}_2+\left(\frac{1}{2}+2x_{10}\right)\mathbf{a}_3$	=	$x_{10}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{y}}-z_{10}c\,\hat{\mathbf{z}}$	$(16l)$	Rh VI
$\mathbf{B}_{38}$	=	$\left(\frac{1}{2}-x_{10}-z_{10}\right)\mathbf{a}_1+(-x_{10}-z_{10})\,\mathbf{a}_2+\left(\frac{1}{2}-2x_{10}\right)\mathbf{a}_3$	=	$-x_{10}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{y}}-z_{10}c\,\hat{\mathbf{z}}$	$(16l)$	Rh VI

$$\mathbf{B}_{39} = \left(\tfrac{1}{2} + x_{11} + z_{11}\right)\mathbf{a}_1 + (x_{11} + z_{11})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + 2x_{11}\right)\mathbf{a}_3 = x_{11}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_{11}\right)a\,\hat{\mathbf{y}} + z_{11}c\,\hat{\mathbf{z}} \quad (16l) \quad \text{Rh VII}$$

$$\mathbf{B}_{40} = \left(\tfrac{1}{2} - x_{11} + z_{11}\right)\mathbf{a}_1 + (-x_{11} + z_{11})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - 2x_{11}\right)\mathbf{a}_3 = -x_{11}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_{11}\right)a\,\hat{\mathbf{y}} + z_{11}c\,\hat{\mathbf{z}} \quad (16l) \quad \text{Rh VII}$$

$$\mathbf{B}_{41} = (x_{11} + z_{11})\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_{11} + z_{11}\right)\mathbf{a}_2 + \tfrac{1}{2}\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{11}\right)a\,\hat{\mathbf{x}} + x_{11}a\,\hat{\mathbf{y}} + z_{11}c\,\hat{\mathbf{z}} \quad (16l) \quad \text{Rh VII}$$

$$\mathbf{B}_{42} = (-x_{11} + z_{11})\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_{11} + z_{11}\right)\mathbf{a}_2 + \tfrac{1}{2}\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{11}\right)a\,\hat{\mathbf{x}} - x_{11}a\,\hat{\mathbf{y}} + z_{11}c\,\hat{\mathbf{z}} \quad (16l) \quad \text{Rh VII}$$

$$\mathbf{B}_{43} = (x_{11} - z_{11})\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_{11} - z_{11}\right)\mathbf{a}_2 + \tfrac{1}{2}\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{11}\right)a\,\hat{\mathbf{x}} + x_{11}a\,\hat{\mathbf{y}} - z_{11}c\,\hat{\mathbf{z}} \quad (16l) \quad \text{Rh VII}$$

$$\mathbf{B}_{44} = (-x_{11} - z_{11})\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_{11} - z_{11}\right)\mathbf{a}_2 + \tfrac{1}{2}\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{11}\right)a\,\hat{\mathbf{x}} - x_{11}a\,\hat{\mathbf{y}} - z_{11}c\,\hat{\mathbf{z}} \quad (16l) \quad \text{Rh VII}$$

$$\mathbf{B}_{45} = \left(\tfrac{1}{2} + x_{11} - z_{11}\right)\mathbf{a}_1 + (x_{11} - z_{11})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + 2x_{11}\right)\mathbf{a}_3 = x_{11}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_{11}\right)a\,\hat{\mathbf{y}} - z_{11}c\,\hat{\mathbf{z}} \quad (16l) \quad \text{Rh VII}$$

$$\mathbf{B}_{46} = \left(\tfrac{1}{2} - x_{11} - z_{11}\right)\mathbf{a}_1 + (-x_{11} - z_{11})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - 2x_{11}\right)\mathbf{a}_3 = -x_{11}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_{11}\right)a\,\hat{\mathbf{y}} - z_{11}c\,\hat{\mathbf{z}} \quad (16l) \quad \text{Rh VII}$$

$$\mathbf{B}_{47} = \left(\tfrac{1}{2} + x_{12} + z_{12}\right)\mathbf{a}_1 + (x_{12} + z_{12})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + 2x_{12}\right)\mathbf{a}_3 = x_{12}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_{12}\right)a\,\hat{\mathbf{y}} + z_{12}c\,\hat{\mathbf{z}} \quad (16l) \quad \text{Rh VIII}$$

$$\mathbf{B}_{48} = \left(\tfrac{1}{2} - x_{12} + z_{12}\right)\mathbf{a}_1 + (-x_{12} + z_{12})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - 2x_{12}\right)\mathbf{a}_3 = -x_{12}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_{12}\right)a\,\hat{\mathbf{y}} + z_{12}c\,\hat{\mathbf{z}} \quad (16l) \quad \text{Rh VIII}$$

$$\mathbf{B}_{49} = (x_{12} + z_{12})\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_{12} + z_{12}\right)\mathbf{a}_2 + \tfrac{1}{2}\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{12}\right)a\,\hat{\mathbf{x}} + x_{12}a\,\hat{\mathbf{y}} + z_{12}c\,\hat{\mathbf{z}} \quad (16l) \quad \text{Rh VIII}$$

$$\mathbf{B}_{50} = (-x_{12} + z_{12})\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_{12} + z_{12}\right)\mathbf{a}_2 + \tfrac{1}{2}\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{12}\right)a\,\hat{\mathbf{x}} - x_{12}a\,\hat{\mathbf{y}} + z_{12}c\,\hat{\mathbf{z}} \quad (16l) \quad \text{Rh VIII}$$

$$\mathbf{B}_{51} = (x_{12} - z_{12})\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_{12} - z_{12}\right)\mathbf{a}_2 + \tfrac{1}{2}\mathbf{a}_3 = \left(\tfrac{1}{2} - x_{12}\right)a\,\hat{\mathbf{x}} + x_{12}a\,\hat{\mathbf{y}} - z_{12}c\,\hat{\mathbf{z}} \quad (16l) \quad \text{Rh VIII}$$

$$\mathbf{B}_{52} = (-x_{12} - z_{12})\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_{12} - z_{12}\right)\mathbf{a}_2 + \tfrac{1}{2}\mathbf{a}_3 = \left(\tfrac{1}{2} + x_{12}\right)a\,\hat{\mathbf{x}} - x_{12}a\,\hat{\mathbf{y}} - z_{12}c\,\hat{\mathbf{z}} \quad (16l) \quad \text{Rh VIII}$$

$$\mathbf{B}_{53} = \left(\tfrac{1}{2} + x_{12} - z_{12}\right)\mathbf{a}_1 + (x_{12} - z_{12})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + 2x_{12}\right)\mathbf{a}_3 = x_{12}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_{12}\right)a\,\hat{\mathbf{y}} - z_{12}c\,\hat{\mathbf{z}} \quad (16l) \quad \text{Rh VIII}$$

$$\mathbf{B}_{54} = \left(\tfrac{1}{2} - x_{12} - z_{12}\right)\mathbf{a}_1 + (-x_{12} - z_{12})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - 2x_{12}\right)\mathbf{a}_3 = -x_{12}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_{12}\right)a\,\hat{\mathbf{y}} - z_{12}c\,\hat{\mathbf{z}} \quad (16l) \quad \text{Rh VIII}$$

$$\mathbf{B}_{55} = (y_{13} + z_{13})\,\mathbf{a}_1 + (x_{13} + z_{13})\,\mathbf{a}_2 + (x_{13} + y_{13})\,\mathbf{a}_3 = x_{13}a\,\hat{\mathbf{x}} + y_{13}a\,\hat{\mathbf{y}} + z_{13}c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu V}$$

$$\mathbf{B}_{56} = (-y_{13} + z_{13})\,\mathbf{a}_1 + (-x_{13} + z_{13})\,\mathbf{a}_2 + (-x_{13} - y_{13})\,\mathbf{a}_3 = -x_{13}a\,\hat{\mathbf{x}} - y_{13}a\,\hat{\mathbf{y}} + z_{13}c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu V}$$

$$\mathbf{B}_{57} = (x_{13} + z_{13})\,\mathbf{a}_1 + (-y_{13} + z_{13})\,\mathbf{a}_2 + (x_{13} - y_{13})\,\mathbf{a}_3 = -y_{13}a\,\hat{\mathbf{x}} + x_{13}a\,\hat{\mathbf{y}} + z_{13}c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu V}$$

$$\mathbf{B}_{58} = (-x_{13} + z_{13})\,\mathbf{a}_1 + (y_{13} + z_{13})\,\mathbf{a}_2 + (-x_{13} + y_{13})\,\mathbf{a}_3 = y_{13}a\,\hat{\mathbf{x}} - x_{13}a\,\hat{\mathbf{y}} + z_{13}c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu V}$$

$$\mathbf{B}_{59} = \left(\tfrac{1}{2} + y_{13} - z_{13}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_{13} - z_{13}\right)\mathbf{a}_2 + (-x_{13} + y_{13})\,\mathbf{a}_3 = -x_{13}a\,\hat{\mathbf{x}} + y_{13}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{13}\right)c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu V}$$

$$\mathbf{B}_{60} = \left(\tfrac{1}{2} - y_{13} - z_{13}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_{13} - z_{13}\right)\mathbf{a}_2 + (x_{13} - y_{13})\,\mathbf{a}_3 = x_{13}a\,\hat{\mathbf{x}} - y_{13}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{13}\right)c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu V}$$

$$\mathbf{B}_{61} = \left(\tfrac{1}{2} + x_{13} - z_{13}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{13} - z_{13}\right)\mathbf{a}_2 + (x_{13} + y_{13})\,\mathbf{a}_3 = y_{13}a\,\hat{\mathbf{x}} + x_{13}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{13}\right)c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu V}$$

$$\mathbf{B}_{62} = \left(\tfrac{1}{2} - x_{13} - z_{13}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{13} - z_{13}\right)\mathbf{a}_2 + (-x_{13} - y_{13})\,\mathbf{a}_3 = -y_{13}a\,\hat{\mathbf{x}} - x_{13}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{13}\right)c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu V}$$

$$\mathbf{B}_{63} = (-y_{13} - z_{13})\,\mathbf{a}_1 + (-x_{13} - z_{13})\,\mathbf{a}_2 + (-x_{13} - y_{13})\,\mathbf{a}_3 = -x_{13}a\,\hat{\mathbf{x}} - y_{13}a\,\hat{\mathbf{y}} - z_{13}c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu V}$$

$$\mathbf{B}_{64} = (y_{13} - z_{13})\,\mathbf{a}_1 + (x_{13} - z_{13})\,\mathbf{a}_2 + (x_{13} + y_{13})\,\mathbf{a}_3 = x_{13}a\,\hat{\mathbf{x}} + y_{13}a\,\hat{\mathbf{y}} - z_{13}c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu V}$$

$$\mathbf{B}_{65} = (-x_{13} - z_{13})\,\mathbf{a}_1 + (y_{13} - z_{13})\,\mathbf{a}_2 + (-x_{13} + y_{13})\,\mathbf{a}_3 = y_{13}a\,\hat{\mathbf{x}} - x_{13}a\,\hat{\mathbf{y}} - z_{13}c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu V}$$

$$\mathbf{B}_{66} = (x_{13} - z_{13})\,\mathbf{a}_1 + (-y_{13} - z_{13})\,\mathbf{a}_2 + (x_{13} - y_{13})\,\mathbf{a}_3 = -y_{13}a\,\hat{\mathbf{x}} + x_{13}a\,\hat{\mathbf{y}} - z_{13}c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu V}$$

$$\mathbf{B}_{67} = \left(\tfrac{1}{2} - y_{13} + z_{13}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_{13} + z_{13}\right)\mathbf{a}_2 + (x_{13} - y_{13})\,\mathbf{a}_3 = x_{13}a\,\hat{\mathbf{x}} - y_{13}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{13}\right)c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu V}$$

$$\mathbf{B}_{68} = \left(\tfrac{1}{2} + y_{13} + z_{13}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_{13} + z_{13}\right)\mathbf{a}_2 + (-x_{13} + y_{13})\,\mathbf{a}_3 = -x_{13}a\,\hat{\mathbf{x}} + y_{13}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{13}\right)c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu V}$$

$$\mathbf{B}_{69} = \left(\tfrac{1}{2} - x_{13} + z_{13}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{13} + z_{13}\right)\mathbf{a}_2 + (-x_{13} - y_{13})\,\mathbf{a}_3 = -y_{13}a\,\hat{\mathbf{x}} - x_{13}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{13}\right)c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu V}$$

$$\mathbf{B}_{70} = \left(\tfrac{1}{2} + x_{13} + z_{13}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{13} + z_{13}\right)\mathbf{a}_2 + (x_{13} + y_{13})\,\mathbf{a}_3 = y_{13}a\,\hat{\mathbf{x}} + x_{13}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{13}\right)c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu V}$$

$$\mathbf{B}_{71} = (y_{14} + z_{14})\,\mathbf{a}_1 + (x_{14} + z_{14})\,\mathbf{a}_2 + (x_{14} + y_{14})\,\mathbf{a}_3 = x_{14}a\,\hat{\mathbf{x}} + y_{14}a\,\hat{\mathbf{y}} + z_{14}c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu VI}$$

$$\mathbf{B}_{72} = (-y_{14} + z_{14})\,\mathbf{a}_1 + (-x_{14} + z_{14})\,\mathbf{a}_2 + (-x_{14} - y_{14})\,\mathbf{a}_3 = -x_{14}a\,\hat{\mathbf{x}} - y_{14}a\,\hat{\mathbf{y}} + z_{14}c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu VI}$$

$$\mathbf{B}_{73} = (x_{14} + z_{14})\,\mathbf{a}_1 + (-y_{14} + z_{14})\,\mathbf{a}_2 + (x_{14} - y_{14})\,\mathbf{a}_3 = -y_{14}a\,\hat{\mathbf{x}} + x_{14}a\,\hat{\mathbf{y}} + z_{14}c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu VI}$$

$$\mathbf{B}_{74} = (-x_{14} + z_{14})\,\mathbf{a}_1 + (y_{14} + z_{14})\,\mathbf{a}_2 + (-x_{14} + y_{14})\,\mathbf{a}_3 = y_{14}a\,\hat{\mathbf{x}} - x_{14}a\,\hat{\mathbf{y}} + z_{14}c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu VI}$$

$$\mathbf{B}_{75} = \left(\tfrac{1}{2} + y_{14} - z_{14}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_{14} - z_{14}\right)\mathbf{a}_2 + (-x_{14} + y_{14})\,\mathbf{a}_3 = -x_{14}a\,\hat{\mathbf{x}} + y_{14}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{14}\right)c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu VI}$$

$$\mathbf{B}_{76} = \left(\tfrac{1}{2} - y_{14} - z_{14}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_{14} - z_{14}\right)\mathbf{a}_2 + (x_{14} - y_{14})\,\mathbf{a}_3 = x_{14}a\,\hat{\mathbf{x}} - y_{14}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{14}\right)c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu VI}$$

$$\mathbf{B}_{77} = \left(\tfrac{1}{2} + x_{14} - z_{14}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{14} - z_{14}\right)\mathbf{a}_2 + (x_{14} + y_{14})\,\mathbf{a}_3 = y_{14}a\,\hat{\mathbf{x}} + x_{14}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{14}\right)c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu VI}$$

$$\mathbf{B}_{78} = \left(\tfrac{1}{2} - x_{14} - z_{14}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{14} - z_{14}\right)\mathbf{a}_2 + (-x_{14} - y_{14})\,\mathbf{a}_3 = -y_{14}a\,\hat{\mathbf{x}} - x_{14}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{14}\right)c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu VI}$$

$$\mathbf{B}_{79} = (-y_{14} - z_{14})\,\mathbf{a}_1 + (-x_{14} - z_{14})\,\mathbf{a}_2 + (-x_{14} - y_{14})\,\mathbf{a}_3 = -x_{14}a\,\hat{\mathbf{x}} - y_{14}a\,\hat{\mathbf{y}} - z_{14}c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu VI}$$

$$\mathbf{B}_{80} = (y_{14} - z_{14})\,\mathbf{a}_1 + (x_{14} - z_{14})\,\mathbf{a}_2 + (x_{14} + y_{14})\,\mathbf{a}_3 = x_{14}a\,\hat{\mathbf{x}} + y_{14}a\,\hat{\mathbf{y}} - z_{14}c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu VI}$$

$$\mathbf{B}_{81} = (-x_{14} - z_{14})\,\mathbf{a}_1 + (y_{14} - z_{14})\,\mathbf{a}_2 + (-x_{14} + y_{14})\,\mathbf{a}_3 = y_{14}a\,\hat{\mathbf{x}} - x_{14}a\,\hat{\mathbf{y}} - z_{14}c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu VI}$$

$$\mathbf{B}_{82} = (x_{14} - z_{14})\,\mathbf{a}_1 + (-y_{14} - z_{14})\,\mathbf{a}_2 + (x_{14} - y_{14})\,\mathbf{a}_3 = -y_{14}a\,\hat{\mathbf{x}} + x_{14}a\,\hat{\mathbf{y}} - z_{14}c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu VI}$$

$$\mathbf{B}_{83} = \left(\tfrac{1}{2} - y_{14} + z_{14}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_{14} + z_{14}\right)\mathbf{a}_2 + (x_{14} - y_{14})\,\mathbf{a}_3 = x_{14}a\,\hat{\mathbf{x}} - y_{14}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{14}\right)c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu VI}$$

$$\mathbf{B}_{84} = \left(\tfrac{1}{2} + y_{14} + z_{14}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_{14} + z_{14}\right)\mathbf{a}_2 + (-x_{14} + y_{14})\,\mathbf{a}_3 = -x_{14}a\,\hat{\mathbf{x}} + y_{14}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{14}\right)c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu VI}$$

$$\mathbf{B}_{85} = \left(\tfrac{1}{2} - x_{14} + z_{14}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{14} + z_{14}\right)\mathbf{a}_2 + (-x_{14} - y_{14})\,\mathbf{a}_3 = -y_{14}a\,\hat{\mathbf{x}} - x_{14}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{14}\right)c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu VI}$$

$$\mathbf{B}_{86} = \left(\tfrac{1}{2} + x_{14} + z_{14}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{14} + z_{14}\right)\mathbf{a}_2 + (x_{14} + y_{14})\,\mathbf{a}_3 = y_{14}a\,\hat{\mathbf{x}} + x_{14}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{14}\right)c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu VI}$$

$$\mathbf{B}_{87} = (y_{15} + z_{15})\,\mathbf{a}_1 + (x_{15} + z_{15})\,\mathbf{a}_2 + (x_{15} + y_{15})\,\mathbf{a}_3 = x_{15}a\,\hat{\mathbf{x}} + y_{15}a\,\hat{\mathbf{y}} + z_{15}c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu VII}$$

$$\mathbf{B}_{88} = (-y_{15} + z_{15})\,\mathbf{a}_1 + (-x_{15} + z_{15})\,\mathbf{a}_2 + (-x_{15} - y_{15})\,\mathbf{a}_3 = -x_{15}a\,\hat{\mathbf{x}} - y_{15}a\,\hat{\mathbf{y}} + z_{15}c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu VII}$$

$$\mathbf{B}_{89} = (x_{15} + z_{15})\,\mathbf{a}_1 + (-y_{15} + z_{15})\,\mathbf{a}_2 + (x_{15} - y_{15})\,\mathbf{a}_3 = -y_{15}a\,\hat{\mathbf{x}} + x_{15}a\,\hat{\mathbf{y}} + z_{15}c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu VII}$$

$$\mathbf{B}_{90} = (-x_{15} + z_{15})\,\mathbf{a}_1 + (y_{15} + z_{15})\,\mathbf{a}_2 + (-x_{15} + y_{15})\,\mathbf{a}_3 = y_{15}a\,\hat{\mathbf{x}} - x_{15}a\,\hat{\mathbf{y}} + z_{15}c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu VII}$$

$$\mathbf{B}_{91} = \left(\tfrac{1}{2} + y_{15} - z_{15}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_{15} - z_{15}\right)\mathbf{a}_2 + (-x_{15} + y_{15})\,\mathbf{a}_3 = -x_{15}a\,\hat{\mathbf{x}} + y_{15}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{15}\right)c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu VII}$$

$$\mathbf{B}_{92} = \left(\tfrac{1}{2} - y_{15} - z_{15}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_{15} - z_{15}\right)\mathbf{a}_2 + (x_{15} - y_{15})\,\mathbf{a}_3 = x_{15}a\,\hat{\mathbf{x}} - y_{15}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{15}\right)c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu VII}$$

$$\mathbf{B}_{93} = \left(\tfrac{1}{2} + x_{15} - z_{15}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{15} - z_{15}\right)\mathbf{a}_2 + (x_{15} + y_{15})\,\mathbf{a}_3 = y_{15}a\,\hat{\mathbf{x}} + x_{15}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{15}\right)c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu VII}$$

$$\mathbf{B}_{94} = \left(\tfrac{1}{2} - x_{15} - z_{15}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{15} - z_{15}\right)\mathbf{a}_2 + (-x_{15} - y_{15})\,\mathbf{a}_3 = -y_{15}a\,\hat{\mathbf{x}} - x_{15}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{15}\right)c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu VII}$$

$$\mathbf{B}_{95} = (-y_{15} - z_{15})\,\mathbf{a}_1 + (-x_{15} - z_{15})\,\mathbf{a}_2 + (-x_{15} - y_{15})\,\mathbf{a}_3 = -x_{15}a\,\hat{\mathbf{x}} - y_{15}a\,\hat{\mathbf{y}} - z_{15}c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu VII}$$

$$\mathbf{B}_{96} = (y_{15} - z_{15})\,\mathbf{a}_1 + (x_{15} - z_{15})\,\mathbf{a}_2 + (x_{15} + y_{15})\,\mathbf{a}_3 = x_{15}a\,\hat{\mathbf{x}} + y_{15}a\,\hat{\mathbf{y}} - z_{15}c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu VII}$$

$$\mathbf{B}_{97} = (-x_{15} - z_{15})\,\mathbf{a}_1 + (y_{15} - z_{15})\,\mathbf{a}_2 + (-x_{15} + y_{15})\,\mathbf{a}_3 = y_{15}a\,\hat{\mathbf{x}} - x_{15}a\,\hat{\mathbf{y}} - z_{15}c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu VII}$$

$$\mathbf{B}_{98} = (x_{15} - z_{15})\,\mathbf{a}_1 + (-y_{15} - z_{15})\,\mathbf{a}_2 + (x_{15} - y_{15})\,\mathbf{a}_3 = -y_{15}a\,\hat{\mathbf{x}} + x_{15}a\,\hat{\mathbf{y}} - z_{15}c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu VII}$$

$$\mathbf{B}_{99} = \left(\tfrac{1}{2} - y_{15} + z_{15}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_{15} + z_{15}\right)\mathbf{a}_2 + (x_{15} - y_{15})\,\mathbf{a}_3 = x_{15}a\,\hat{\mathbf{x}} - y_{15}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{15}\right)c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu VII}$$

$$\mathbf{B}_{100} = \left(\tfrac{1}{2} + y_{15} + z_{15}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_{15} + z_{15}\right)\mathbf{a}_2 + (-x_{15} + y_{15})\,\mathbf{a}_3 = -x_{15}a\,\hat{\mathbf{x}} + y_{15}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{15}\right)c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu VII}$$

$$\mathbf{B}_{101} = \left(\tfrac{1}{2} - x_{15} + z_{15}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{15} + z_{15}\right)\mathbf{a}_2 + (-x_{15} - y_{15})\,\mathbf{a}_3 = -y_{15}a\,\hat{\mathbf{x}} - x_{15}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{15}\right)c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu VII}$$

$$\mathbf{B}_{102} = \left(\tfrac{1}{2} + x_{15} + z_{15}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{15} + z_{15}\right)\mathbf{a}_2 + (x_{15} + y_{15})\,\mathbf{a}_3 = y_{15}a\,\hat{\mathbf{x}} + x_{15}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{15}\right)c\,\hat{\mathbf{z}} \quad (32m) \quad \text{Pu VII}$$

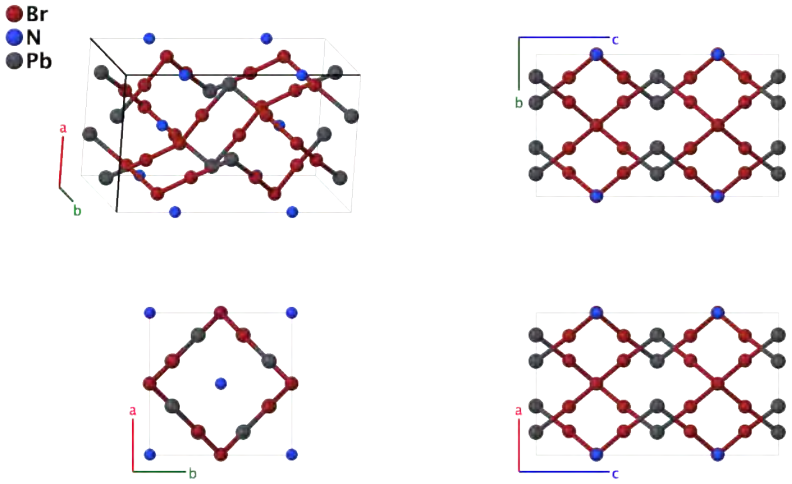
References:

- D. T. Cromer and A. C. Larson, *The Crystal Structure of $Pu_{31}Pt_{20}$ and $Pu_{31}Rh_{20}$*, Acta Crystallogr. Sect. B Struct. Sci. **33**, 2620–2627 (1977), doi:10.1107/S0567740877009030.

Geometry files:
- CIF: pp. 1713
- POSCAR: pp. 1713

$NH_4Pb_2Br_5$ ($K3_4$) Structure: A5BC2_tI32_140_bl_a_h

http://aflow.org/prototype-encyclopedia/A5BC2_tI32_140_bl_a_h



Prototype	:	$Br_5(NH_4)Pb_2$
AFLOW prototype label	:	A5BC2_tI32_140_bl_a_h
***Strukturbericht* designation**	:	$K3_4$
Pearson symbol	:	tI32
Space group number	:	140
Space group symbol	:	$I4/mcm$
AFLOW prototype command	:	`aflow --proto=A5BC2_tI32_140_bl_a_h --params=`$a, c/a, x_3, x_4, z_4$

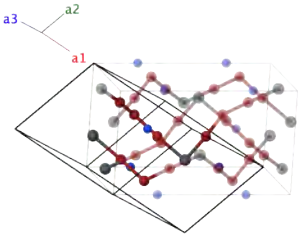
Other compounds with this structure

- KPb_2Br_5, $RbPb_2Br_5$, KSn_2Br_5, KSn_2I_5, $InSn_2Br_5$, $InSn_2I_5$, $CsSn_2Br_5$, $TlSn_2Br_5$, $RbSn_2Br_5$, $InPb_2I_5$, and ISe_2Tl_5

- The positions of the hydrogen atoms in the NH_4 ions were not determined, so we only provide the positions of the nitrogen atoms (labeled as NH_4).

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}a\,\hat{\mathbf{x}} - \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} - \tfrac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\tfrac{1}{4}\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2$	=	$\tfrac{1}{4}c\,\hat{\mathbf{z}}$	$(4a)$	NH_4
$\mathbf{B}_2$	=	$\tfrac{3}{4}\mathbf{a}_1 + \tfrac{3}{4}\mathbf{a}_2$	=	$\tfrac{3}{4}c\,\hat{\mathbf{z}}$	$(4a)$	NH_4

$\mathbf{B}_3$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*b*)	Br I
$\mathbf{B}_4$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*b*)	Br I
$\mathbf{B}_5$	=	$\left(\frac{1}{2}+x_3\right)\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \left(\frac{1}{2}+2x_3\right)\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{y}}$	(8*h*)	Pb
$\mathbf{B}_6$	=	$\left(\frac{1}{2}-x_3\right)\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\frac{1}{2}-2x_3\right)\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{y}}$	(8*h*)	Pb
$\mathbf{B}_7$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2}-x_3\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}}$	(8*h*)	Pb
$\mathbf{B}_8$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2}+x_3\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}}$	(8*h*)	Pb
$\mathbf{B}_9$	=	$\left(\frac{1}{2}+x_4+z_4\right)\mathbf{a}_1 + (x_4+z_4)\,\mathbf{a}_2 + \left(\frac{1}{2}+2x_4\right)\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	(16*l*)	Br II
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2}-x_4+z_4\right)\mathbf{a}_1 + (-x_4+z_4)\,\mathbf{a}_2 + \left(\frac{1}{2}-2x_4\right)\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	(16*l*)	Br II
$\mathbf{B}_{11}$	=	$(x_4+z_4)\,\mathbf{a}_1 + \left(\frac{1}{2}-x_4+z_4\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	(16*l*)	Br II
$\mathbf{B}_{12}$	=	$(-x_4+z_4)\,\mathbf{a}_1 + \left(\frac{1}{2}+x_4+z_4\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	(16*l*)	Br II
$\mathbf{B}_{13}$	=	$(x_4-z_4)\,\mathbf{a}_1 + \left(\frac{1}{2}-x_4-z_4\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	(16*l*)	Br II
$\mathbf{B}_{14}$	=	$(-x_4-z_4)\,\mathbf{a}_1 + \left(\frac{1}{2}+x_4-z_4\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	(16*l*)	Br II
$\mathbf{B}_{15}$	=	$\left(\frac{1}{2}+x_4-z_4\right)\mathbf{a}_1 + (x_4-z_4)\,\mathbf{a}_2 + \left(\frac{1}{2}+2x_4\right)\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	(16*l*)	Br II
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2}-x_4-z_4\right)\mathbf{a}_1 + (-x_4-z_4)\,\mathbf{a}_2 + \left(\frac{1}{2}-2x_4\right)\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	(16*l*)	Br II

References:

- H. M. Powell and H. S. Tasker, *The valency angle of bivalent lead: the crystal structure of ammonium, rubidium, and potassium pentabromodiplumbites*, J. Chem. Soc. pp. 119–123 (1937), doi:10.1039/JR9370000119.

Found in:

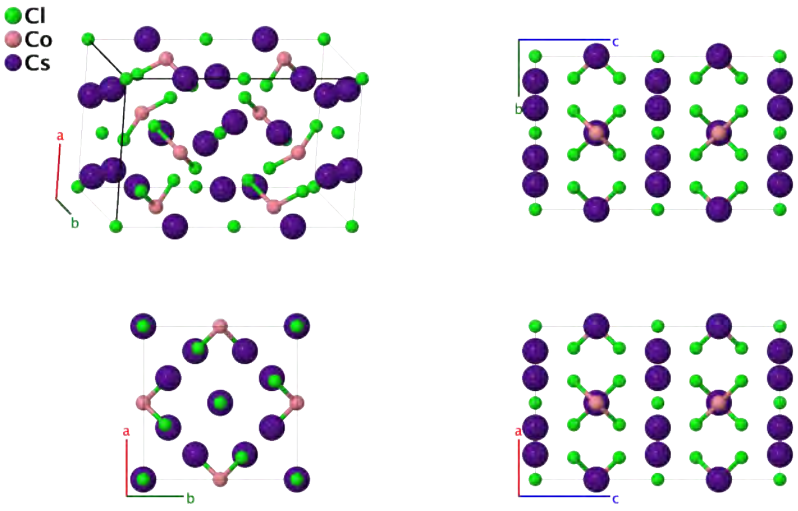
- C. Gottfried, ed., *Strukturbericht Band V 1937* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1940).

Geometry files:

- CIF: pp. 1714
- POSCAR: pp. 1714

Cs_3CoCl_5 ($K3_1$) Structure: A5BC3_tI36_140_cl_b_ah

http://aflow.org/prototype-encyclopedia/A5BC3_tI36_140_cl_b_ah



Prototype	:	Cl_5CoCs_3
AFLOW prototype label	:	A5BC3_tI36_140_cl_b_ah
***Strukturbericht* designation**	:	$K3_1$
Pearson symbol	:	tI36
Space group number	:	140
Space group symbol	:	$I4/mcm$
AFLOW prototype command	:	`aflow --proto=A5BC3_tI36_140_cl_b_ah --params=`$a, c/a, x_4, x_5, z_5$

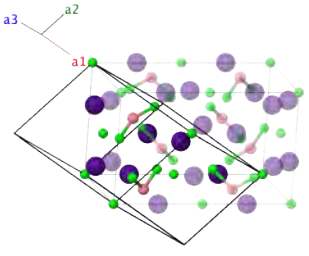
Other compounds with this structure

- Cs_3CoBr_5, $CsBa_5Ti_2Se_9Cl$, Rb_3ZnH_5, Cs_3ZnH_5, $(Sr_{3-x}A_x)AlO_4F$, (A = Ba, Ca), Ba_3MO_5, (M = tetravalent metal, and $M_3Mg(BH_4)_5$, (M = Rb, Cs)

- We show the structure using the data taken at 4.2 K.

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}a\,\hat{\mathbf{x}} - \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} - \tfrac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\tfrac{1}{4}\mathbf{a}_1 + \tfrac{1}{4}\mathbf{a}_2$	=	$\tfrac{1}{4}c\,\hat{\mathbf{z}}$	($4a$)	Cs I

$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2$	=	$\frac{3}{4}c\,\hat{\mathbf{z}}$	(4*a*)	Cs I
$\mathbf{B}_3$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*b*)	Co
$\mathbf{B}_4$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*b*)	Co
$\mathbf{B}_5$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(4*c*)	Cl I
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	(4*c*)	Cl I
$\mathbf{B}_7$	=	$\left(\frac{1}{2}+x_4\right)\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \left(\frac{1}{2}+2x_4\right)\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{y}}$	(8*h*)	Cs II
$\mathbf{B}_8$	=	$\left(\frac{1}{2}-x_4\right)\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \left(\frac{1}{2}-2x_4\right)\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{y}}$	(8*h*)	Cs II
$\mathbf{B}_9$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2}-x_4\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}}$	(8*h*)	Cs II
$\mathbf{B}_{10}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2}+x_4\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}}$	(8*h*)	Cs II
$\mathbf{B}_{11}$	=	$\left(\frac{1}{2}+x_5+z_5\right)\mathbf{a}_1 + (x_5+z_5)\,\mathbf{a}_2 + \left(\frac{1}{2}+2x_5\right)\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(16*l*)	Cl II
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2}-x_5+z_5\right)\mathbf{a}_1 + (-x_5+z_5)\,\mathbf{a}_2 + \left(\frac{1}{2}-2x_5\right)\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(16*l*)	Cl II
$\mathbf{B}_{13}$	=	$(x_5+z_5)\,\mathbf{a}_1 + \left(\frac{1}{2}-x_5+z_5\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(16*l*)	Cl II
$\mathbf{B}_{14}$	=	$(-x_5+z_5)\,\mathbf{a}_1 + \left(\frac{1}{2}+x_5+z_5\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(16*l*)	Cl II
$\mathbf{B}_{15}$	=	$(x_5-z_5)\,\mathbf{a}_1 + \left(\frac{1}{2}-x_5-z_5\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(16*l*)	Cl II
$\mathbf{B}_{16}$	=	$(-x_5-z_5)\,\mathbf{a}_1 + \left(\frac{1}{2}+x_5-z_5\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(16*l*)	Cl II
$\mathbf{B}_{17}$	=	$\left(\frac{1}{2}+x_5-z_5\right)\mathbf{a}_1 + (x_5-z_5)\,\mathbf{a}_2 + \left(\frac{1}{2}+2x_5\right)\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(16*l*)	Cl II
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2}-x_5-z_5\right)\mathbf{a}_1 + (-x_5-z_5)\,\mathbf{a}_2 + \left(\frac{1}{2}-2x_5\right)\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(16*l*)	Cl II

References:

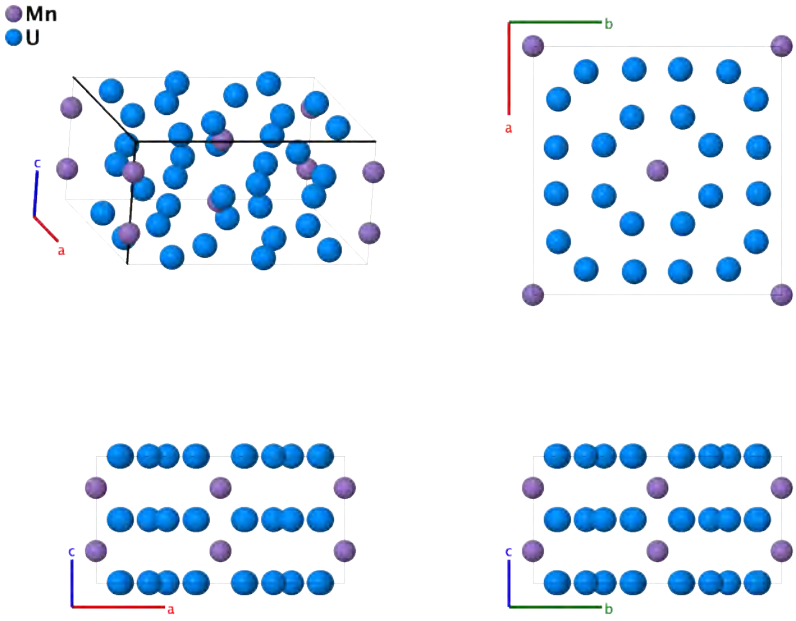
- B. N. Figgis, R. Mason, A. R. P. Smith, and G. A. Williams, *Neutron Diffraction Structure of Cs_3CoCl_5 at 4.2 K*, Acta Crystallogr. Sect. B Struct. Sci. **36**, 509–512 (1980), doi:10.1107/S0567740880003731.

Geometry files:

- CIF: pp. 1715
- POSCAR: pp. 1715

U_6Mn ($D2_c$) Structure: AB6_tI28_140_a_hk

http://aflow.org/prototype-encyclopedia/AB6_tI28_140_a_hk



Prototype	:	MnU_6
AFLOW prototype label	:	AB6_tI28_140_a_hk
***Strukturbericht* designation**	:	$D2_c$
Pearson symbol	:	tI28
Space group number	:	140
Space group symbol	:	$I4/mcm$
AFLOW prototype command	:	`aflow --proto=AB6_tI28_140_a_hk --params=`$a, c/a, x_2, x_3, y_3$

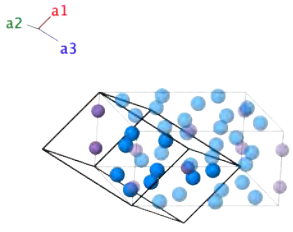
Other compounds with this structure

- U_6Co, U_6Fe, and U_6Ni

- This structure is closely related to the V_4SiSb_2 structure. This can also be identified as a defected version of the $D8_m$ W_5Si_3 structure.

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} - \frac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2$	=	$\frac{1}{4}c\,\hat{\mathbf{z}}$	(4*a*)	Mn
$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2$	=	$\frac{3}{4}c\,\hat{\mathbf{z}}$	(4*a*)	Mn
$\mathbf{B}_3$	=	$\left(\frac{1}{2}+x_2\right)\mathbf{a}_1 + x_2\,\mathbf{a}_2 + \left(\frac{1}{2}+2x_2\right)\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_2\right)a\,\hat{\mathbf{y}}$	(8*h*)	U I
$\mathbf{B}_4$	=	$\left(\frac{1}{2}-x_2\right)\mathbf{a}_1 - x_2\,\mathbf{a}_2 + \left(\frac{1}{2}-2x_2\right)\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_2\right)a\,\hat{\mathbf{y}}$	(8*h*)	U I
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2}-x_2\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_2\right)a\,\hat{\mathbf{x}} + x_2 a\,\hat{\mathbf{y}}$	(8*h*)	U I
$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2}+x_2\right)\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_2\right)a\,\hat{\mathbf{x}} - x_2 a\,\hat{\mathbf{y}}$	(8*h*)	U I
$\mathbf{B}_7$	=	$y_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + (x_3+y_3)\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 a\,\hat{\mathbf{y}}$	(16*k*)	U II
$\mathbf{B}_8$	=	$-y_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + (-x_3-y_3)\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 a\,\hat{\mathbf{y}}$	(16*k*)	U II
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + (x_3-y_3)\,\mathbf{a}_3$	=	$-y_3 a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}}$	(16*k*)	U II
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + (-x_3+y_3)\,\mathbf{a}_3$	=	$y_3 a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}}$	(16*k*)	U II
$\mathbf{B}_{11}$	=	$\left(\frac{1}{2}+y_3\right)\mathbf{a}_1 + \left(\frac{1}{2}-x_3\right)\mathbf{a}_2 + (-x_3+y_3)\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + y_3 a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(16*k*)	U II
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2}-y_3\right)\mathbf{a}_1 + \left(\frac{1}{2}+x_3\right)\mathbf{a}_2 + (x_3-y_3)\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} - y_3 a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(16*k*)	U II
$\mathbf{B}_{13}$	=	$\left(\frac{1}{2}+x_3\right)\mathbf{a}_1 + \left(\frac{1}{2}+y_3\right)\mathbf{a}_2 + (x_3+y_3)\,\mathbf{a}_3$	=	$y_3 a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(16*k*)	U II
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2}-x_3\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_3\right)\mathbf{a}_2 + (-x_3-y_3)\,\mathbf{a}_3$	=	$-y_3 a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(16*k*)	U II

References:

- N. C. Baenziger, R. E. Rundle, A. I. Snow, and A. S. Wilson, *Compounds of uranium with the transition metals of the first long period*, Acta Cryst. **3**, 34–40 (1950), doi:10.1107/S0365110X50000082.

Found in:

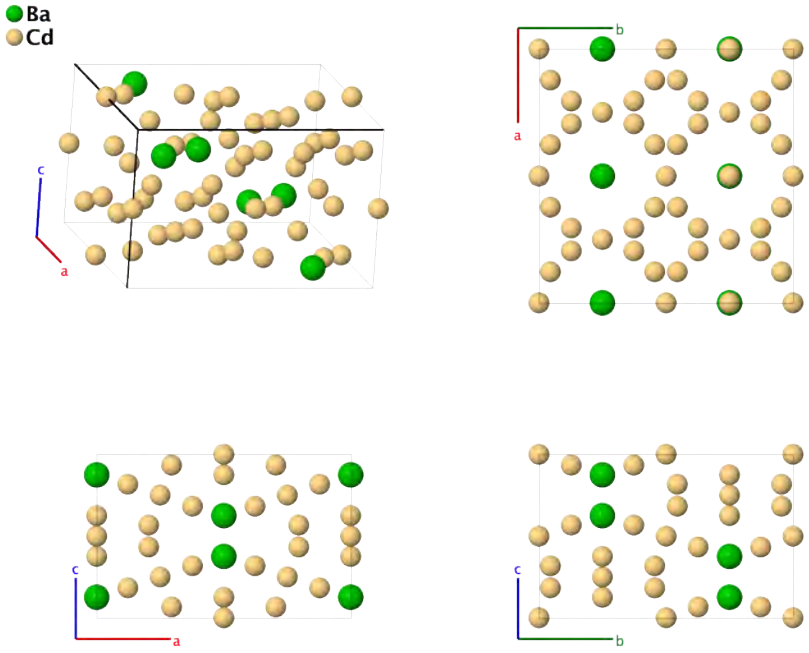
- P. Villars (Chief Editor), *PAULING FILE*, http://materials.springer.com/isp/crystallographic/docs/sd_1250751 (2016). In: Inorganic Solid Phases, SpringerMaterials (online database), Springer, Heidelberg.

Geometry files:

- CIF: pp. 1715
- POSCAR: pp. 1716

$BaCd_{11}$ Structure: AB11_tI48_141_a_bdi

http://aflow.org/prototype-encyclopedia/AB11_tI48_141_a_bdi



Prototype	:	$BaCd_{11}$
AFLOW prototype label	:	AB11_tI48_141_a_bdi
***Strukturbericht* designation**	:	None
Pearson symbol	:	tI48
Space group number	:	141
Space group symbol	:	$I4_1/amd$
AFLOW prototype command	:	`aflow --proto=AB11_tI48_141_a_bdi --params=`$a, c/a, x_4, y_4, z_4$

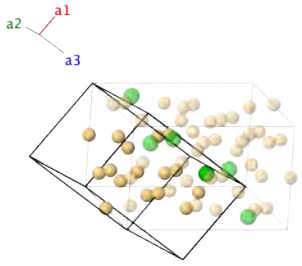
Other compounds with this structure

- $SrCd_{11}$, $LaZn_{11}$, $CeZn_{11}$, and $PrZn_{11}$

- (Sanderson, 1953) gave the atomic positions in setting 1 of space group #141. We used FINDSYM to transform this to the standard setting 2.

Body-centered Tetragonal primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= -\tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{y}} + \tfrac{1}{2}\,c\,\hat{\mathbf{z}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2}\,a\,\hat{\mathbf{y}} + \tfrac{1}{2}\,c\,\hat{\mathbf{z}}\\ \mathbf{a}_3 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{y}} - \tfrac{1}{2}\,c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{7}{8}\,\mathbf{a}_1 + \frac{1}{8}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{y}} + \frac{1}{8}c\,\hat{\mathbf{z}}$	(4*a*)	Ba
$\mathbf{B}_2$	=	$\frac{1}{8}\,\mathbf{a}_1 + \frac{7}{8}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{3}{8}c\,\hat{\mathbf{z}}$	(4*a*)	Ba
$\mathbf{B}_3$	=	$\frac{5}{8}\,\mathbf{a}_1 + \frac{3}{8}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{y}} + \frac{3}{8}c\,\hat{\mathbf{z}}$	(4*b*)	Cd I
$\mathbf{B}_4$	=	$\frac{3}{8}\,\mathbf{a}_1 + \frac{5}{8}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{8}c\,\hat{\mathbf{z}}$	(4*b*)	Cd I
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	(8*d*)	Cd II
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	(8*d*)	Cd II
$\mathbf{B}_7$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$-\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(8*d*)	Cd II
$\mathbf{B}_8$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(8*d*)	Cd II
$\mathbf{B}_9$	=	$(y_4 + z_4)\,\mathbf{a}_1 + (x_4 + z_4)\,\mathbf{a}_2 + (x_4 + y_4)\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + y_4a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	(32*i*)	Cd III
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - y_4 + z_4\right)\mathbf{a}_1 + (-x_4 + z_4)\,\mathbf{a}_2 + \left(\frac{1}{2} - x_4 - y_4\right)\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	(32*i*)	Cd III
$\mathbf{B}_{11}$	=	$(x_4 + z_4)\,\mathbf{a}_1 + \left(\frac{1}{2} - y_4 + z_4\right)\mathbf{a}_2 + (x_4 - y_4)\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - y_4\right)a\,\hat{\mathbf{x}} + \left(\frac{3}{4} + x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_4\right)c\,\hat{\mathbf{z}}$	(32*i*)	Cd III
$\mathbf{B}_{12}$	=	$(-x_4 + z_4)\,\mathbf{a}_1 + (y_4 + z_4)\,\mathbf{a}_2 + \left(\frac{1}{2} - x_4 + y_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + y_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{3}{4} + z_4\right)c\,\hat{\mathbf{z}}$	(32*i*)	Cd III
$\mathbf{B}_{13}$	=	$\left(\frac{1}{2} + y_4 - z_4\right)\mathbf{a}_1 + (-x_4 - z_4)\,\mathbf{a}_2 + \left(\frac{1}{2} - x_4 + y_4\right)\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	(32*i*)	Cd III
$\mathbf{B}_{14}$	=	$(-y_4 - z_4)\,\mathbf{a}_1 + (x_4 - z_4)\,\mathbf{a}_2 + (x_4 - y_4)\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} - y_4a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	(32*i*)	Cd III
$\mathbf{B}_{15}$	=	$(x_4 - z_4)\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4 - z_4\right)\mathbf{a}_2 + (x_4 + y_4)\,\mathbf{a}_3$	=	$\left(\frac{1}{4} + y_4\right)a\,\hat{\mathbf{x}} + \left(\frac{3}{4} + x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_4\right)c\,\hat{\mathbf{z}}$	(32*i*)	Cd III
$\mathbf{B}_{16}$	=	$(-x_4 - z_4)\,\mathbf{a}_1 + (-y_4 - z_4)\,\mathbf{a}_2 + \left(\frac{1}{2} - x_4 - y_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - y_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{3}{4} - z_4\right)c\,\hat{\mathbf{z}}$	(32*i*)	Cd III
$\mathbf{B}_{17}$	=	$(-y_4 - z_4)\,\mathbf{a}_1 + (-x_4 - z_4)\,\mathbf{a}_2 + (-x_4 - y_4)\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} - y_4a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	(32*i*)	Cd III
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} + y_4 - z_4\right)\mathbf{a}_1 + (x_4 - z_4)\,\mathbf{a}_2 + \left(\frac{1}{2} + x_4 + y_4\right)\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	(32*i*)	Cd III
$\mathbf{B}_{19}$	=	$(-x_4 - z_4)\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4 - z_4\right)\mathbf{a}_2 + (-x_4 + y_4)\,\mathbf{a}_3$	=	$\left(\frac{1}{4} + y_4\right)a\,\hat{\mathbf{x}} - \left(x_4 + \frac{1}{4}\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_4\right)c\,\hat{\mathbf{z}}$	(32*i*)	Cd III

$\mathbf{B}_{20}$	=	$(x_4 - z_4)\,\mathbf{a}_1 + (-y_4 - z_4)\,\mathbf{a}_2 + \left(\frac{1}{2} + x_4 - y_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - y_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_4\right)a\,\hat{\mathbf{y}} - \left(z_4 + \frac{1}{4}\right)c\,\hat{\mathbf{z}}$	(32*i*)	Cd III
$\mathbf{B}_{21}$	=	$\left(\frac{1}{2} - y_4 + z_4\right)\mathbf{a}_1 + (x_4 + z_4)\,\mathbf{a}_2 + \left(\frac{1}{2} + x_4 - y_4\right)\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(32*i*)	Cd III
$\mathbf{B}_{22}$	=	$(y_4 + z_4)\,\mathbf{a}_1 + (-x_4 + z_4)\,\mathbf{a}_2 + (-x_4 + y_4)\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + y_4 a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(32*i*)	Cd III
$\mathbf{B}_{23}$	=	$(-x_4 + z_4)\,\mathbf{a}_1 + \left(\frac{1}{2} - y_4 + z_4\right)\mathbf{a}_2 + (-x_4 - y_4)\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - y_4\right)a\,\hat{\mathbf{x}} - \left(x_4 + \frac{1}{4}\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_4\right)c\,\hat{\mathbf{z}}$	(32*i*)	Cd III
$\mathbf{B}_{24}$	=	$(x_4 + z_4)\,\mathbf{a}_1 + (y_4 + z_4)\,\mathbf{a}_2 + \left(\frac{1}{2} + x_4 + y_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + y_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_4\right)a\,\hat{\mathbf{y}} + \left(-\frac{1}{4} + z_4\right)c\,\hat{\mathbf{z}}$	(32*i*)	Cd III

References:

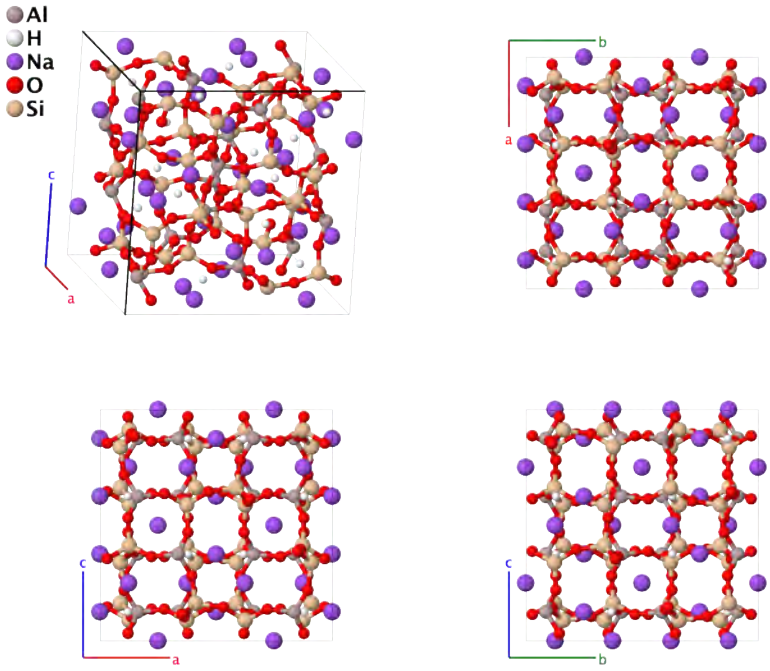
- M. J. Sanderson and N. C. Baenziger, *The Crystal Structure of* $BaCd_{11}$, Acta Cryst. **6**, 627–631 (1953), doi:10.1107/S0365110X53001745.

Geometry files:

- CIF: pp. 1716
- POSCAR: pp. 1716

Analcime ($NaAlSi_2O_6 \cdot H_2O$, $S6_1$) Structure: A2B2C3D12E4_tI184_142_f_f_be_3g_g

http://aflow.org/prototype-encyclopedia/A2B2C3D12E4_tI184_142_f_f_be_3g_g

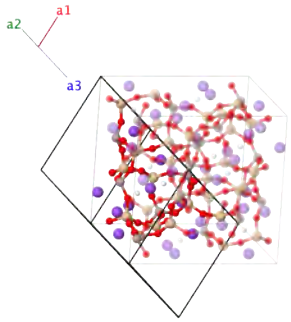


Prototype	:	$Al(H_2O)NaO_6Si_2$
AFLOW prototype label	:	A2B2C3D12E4_tI184_142_f_f_be_3g_g
***Strukturbericht* designation**	:	$S6_1$
Pearson symbol	:	tI184
Space group number	:	142
Space group symbol	:	$I4_1/acd$
AFLOW prototype command	:	`aflow --proto=A2B2C3D12E4_tI184_142_f_f_be_3g_g --params=`$a, c/a, x_2, x_3, x_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8$

- Analcime is pseudo-cubic, space group $Ia\bar{3}d$ #230, but (Hartwig, 1931) showed that the structure actually is tetragonal, space group $I4_1/acd$ #142. (Hermann, 1937) listed both structures in defining *Strukturbericht* type $S6_1$.
- Analcime can have many crystal structures, as found in *e.g.,* (Mazzi, 1978) and (Pechar, 1988). We take our prototype from (Mazzi, 1978), using the data from their sample ANA 1, which is in space group $I4_1/acd$ #142. The major difference between this structure and that found by (Hartwig, 1931) is that sodium atoms were found at the (8*b*) Wyckoff position (Na-I). (Mazzi, 1978) recover the stoichiometry by noting that site Na-I (8*b*) is only occupied 23% of the time, while site Na-II (16*e*) has 82% occupancy. In addition, our Si site (32*g*) is actually 47% aluminum and 53% silicon, and the site which we, following (Hartwig, 1931), label Al (16*f*) is actually 98% silicon and only 2% aluminum. These fractions are highly dependent upon the choice of sample.
- Removing all sodium from the (8*b*) site results in the original structure found by (Hartwig, 1931) and defined as $S6_1$ by (Hermann, 1937).
- The positions of the hydrogen atoms in the water molecules were not determined, so we only provide the positions of the oxygen atoms (labeled as H_2O).

Body-centered Tetragonal primitive vectors:

$$
\begin{aligned}
\mathbf{a}_1 &= -\tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{y}} + \tfrac{1}{2}\,c\,\hat{\mathbf{z}} \\
\mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2}\,a\,\hat{\mathbf{y}} + \tfrac{1}{2}\,c\,\hat{\mathbf{z}} \\
\mathbf{a}_3 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{y}} - \tfrac{1}{2}\,c\,\hat{\mathbf{z}}
\end{aligned}
$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\tfrac{3}{8}\,\mathbf{a}_1 + \tfrac{1}{8}\,\mathbf{a}_2 + \tfrac{1}{4}\,\mathbf{a}_3$	=	$\tfrac{1}{4}a\,\hat{\mathbf{y}} + \tfrac{1}{8}c\,\hat{\mathbf{z}}$	(8*b*)	Na I
$\mathbf{B}_2$	=	$\tfrac{1}{8}\,\mathbf{a}_1 + \tfrac{3}{8}\,\mathbf{a}_2 + \tfrac{3}{4}\,\mathbf{a}_3$	=	$\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{4}a\,\hat{\mathbf{y}} - \tfrac{1}{8}c\,\hat{\mathbf{z}}$	(8*b*)	Na I
$\mathbf{B}_3$	=	$\tfrac{5}{8}\,\mathbf{a}_1 + \tfrac{7}{8}\,\mathbf{a}_2 + \tfrac{3}{4}\,\mathbf{a}_3$	=	$\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{4}a\,\hat{\mathbf{y}} + \tfrac{3}{8}c\,\hat{\mathbf{z}}$	(8*b*)	Na I
$\mathbf{B}_4$	=	$\tfrac{7}{8}\,\mathbf{a}_1 + \tfrac{5}{8}\,\mathbf{a}_2 + \tfrac{1}{4}\,\mathbf{a}_3$	=	$\tfrac{1}{4}a\,\hat{\mathbf{y}} + \tfrac{5}{8}c\,\hat{\mathbf{z}}$	(8*b*)	Na I
$\mathbf{B}_5$	=	$\tfrac{1}{4}\,\mathbf{a}_1 + \left(\tfrac{1}{4} + x_2\right)\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$x_2a\,\hat{\mathbf{x}} + \tfrac{1}{4}c\,\hat{\mathbf{z}}$	(16*e*)	Na II
$\mathbf{B}_6$	=	$\tfrac{3}{4}\,\mathbf{a}_1 + \left(\tfrac{1}{4} - x_2\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_2\right)\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{4}c\,\hat{\mathbf{z}}$	(16*e*)	Na II
$\mathbf{B}_7$	=	$\left(\tfrac{1}{4} + x_2\right)\mathbf{a}_1 + \tfrac{3}{4}\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$\tfrac{1}{4}a\,\hat{\mathbf{x}} + \left(\tfrac{3}{4} + x_2\right)a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$	(16*e*)	Na II
$\mathbf{B}_8$	=	$\left(\tfrac{1}{4} - x_2\right)\mathbf{a}_1 + \tfrac{1}{4}\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_2\right)\mathbf{a}_3$	=	$\tfrac{1}{4}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_2\right)a\,\hat{\mathbf{y}}$	(16*e*)	Na II
$\mathbf{B}_9$	=	$\tfrac{3}{4}\,\mathbf{a}_1 + \left(\tfrac{3}{4} - x_2\right)\mathbf{a}_2 - x_2\,\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}} + \tfrac{3}{4}c\,\hat{\mathbf{z}}$	(16*e*)	Na II
$\mathbf{B}_{10}$	=	$\tfrac{1}{4}\,\mathbf{a}_1 + \left(\tfrac{3}{4} + x_2\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_2\right)\mathbf{a}_3$	=	$\left(\tfrac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \tfrac{1}{4}c\,\hat{\mathbf{z}}$	(16*e*)	Na II
$\mathbf{B}_{11}$	=	$\left(\tfrac{3}{4} - x_2\right)\mathbf{a}_1 + \tfrac{1}{4}\,\mathbf{a}_2 - x_2\,\mathbf{a}_3$	=	$\tfrac{3}{4}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_2\right)a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$	(16*e*)	Na II
$\mathbf{B}_{12}$	=	$\left(\tfrac{3}{4} + x_2\right)\mathbf{a}_1 + \tfrac{3}{4}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_2\right)\mathbf{a}_3$	=	$\tfrac{1}{4}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_2\right)a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$	(16*e*)	Na II
$\mathbf{B}_{13}$	=	$\left(\tfrac{3}{8} + x_3\right)\mathbf{a}_1 + \left(\tfrac{1}{8} + x_3\right)\mathbf{a}_2 + \left(\tfrac{1}{4} + 2x_3\right)\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_3\right)a\,\hat{\mathbf{y}} + \tfrac{1}{8}c\,\hat{\mathbf{z}}$	(16*f*)	Al
$\mathbf{B}_{14}$	=	$\left(\tfrac{3}{8} - x_3\right)\mathbf{a}_1 + \left(\tfrac{1}{8} - x_3\right)\mathbf{a}_2 + \left(\tfrac{1}{4} - 2x_3\right)\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_3\right)a\,\hat{\mathbf{y}} + \tfrac{1}{8}c\,\hat{\mathbf{z}}$	(16*f*)	Al
$\mathbf{B}_{15}$	=	$\left(\tfrac{1}{8} + x_3\right)\mathbf{a}_1 + \left(\tfrac{3}{8} - x_3\right)\mathbf{a}_2 + \tfrac{3}{4}\,\mathbf{a}_3$	=	$\left(\tfrac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_3\right)a\,\hat{\mathbf{y}} - \tfrac{1}{8}c\,\hat{\mathbf{z}}$	(16*f*)	Al
$\mathbf{B}_{16}$	=	$\left(\tfrac{1}{8} - x_3\right)\mathbf{a}_1 + \left(\tfrac{3}{8} + x_3\right)\mathbf{a}_2 + \tfrac{3}{4}\,\mathbf{a}_3$	=	$\left(\tfrac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_3\right)a\,\hat{\mathbf{y}} + \tfrac{7}{8}c\,\hat{\mathbf{z}}$	(16*f*)	Al
$\mathbf{B}_{17}$	=	$\left(\tfrac{5}{8} - x_3\right)\mathbf{a}_1 + \left(\tfrac{7}{8} - x_3\right)\mathbf{a}_2 + \left(\tfrac{3}{4} - 2x_3\right)\mathbf{a}_3$	=	$\left(\tfrac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_3\right)a\,\hat{\mathbf{y}} + \tfrac{3}{8}c\,\hat{\mathbf{z}}$	(16*f*)	Al
$\mathbf{B}_{18}$	=	$\left(\tfrac{5}{8} + x_3\right)\mathbf{a}_1 + \left(\tfrac{7}{8} + x_3\right)\mathbf{a}_2 + \left(\tfrac{3}{4} + 2x_3\right)\mathbf{a}_3$	=	$\left(\tfrac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_3\right)a\,\hat{\mathbf{y}} + \tfrac{3}{8}c\,\hat{\mathbf{z}}$	(16*f*)	Al
$\mathbf{B}_{19}$	=	$\left(\tfrac{7}{8} - x_3\right)\mathbf{a}_1 + \left(\tfrac{5}{8} + x_3\right)\mathbf{a}_2 + \tfrac{1}{4}\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_3\right)a\,\hat{\mathbf{y}} + \tfrac{5}{8}c\,\hat{\mathbf{z}}$	(16*f*)	Al
$\mathbf{B}_{20}$	=	$\left(\tfrac{7}{8} + x_3\right)\mathbf{a}_1 + \left(\tfrac{5}{8} - x_3\right)\mathbf{a}_2 + \tfrac{1}{4}\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_3\right)a\,\hat{\mathbf{y}} + \tfrac{5}{8}c\,\hat{\mathbf{z}}$	(16*f*)	Al
$\mathbf{B}_{21}$	=	$\left(\tfrac{3}{8} + x_4\right)\mathbf{a}_1 + \left(\tfrac{1}{8} + x_4\right)\mathbf{a}_2 + \left(\tfrac{1}{4} + 2x_4\right)\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_4\right)a\,\hat{\mathbf{y}} + \tfrac{1}{8}c\,\hat{\mathbf{z}}$	(16*f*)	H_2O
$\mathbf{B}_{22}$	=	$\left(\tfrac{3}{8} - x_4\right)\mathbf{a}_1 + \left(\tfrac{1}{8} - x_4\right)\mathbf{a}_2 + \left(\tfrac{1}{4} - 2x_4\right)\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_4\right)a\,\hat{\mathbf{y}} + \tfrac{1}{8}c\,\hat{\mathbf{z}}$	(16*f*)	H_2O

$\mathbf{B}_{23}$	=	$\left(\frac{1}{8}+x_4\right)\mathbf{a}_1+\left(\frac{3}{8}-x_4\right)\mathbf{a}_2+\frac{3}{4}\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+x_4\right)a\,\hat{\mathbf{y}}-\frac{1}{8}c\,\hat{\mathbf{z}}$	(16*f*)	H_2O
$\mathbf{B}_{24}$	=	$\left(\frac{1}{8}-x_4\right)\mathbf{a}_1+\left(\frac{3}{8}+x_4\right)\mathbf{a}_2+\frac{3}{4}\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-x_4\right)a\,\hat{\mathbf{y}}+\frac{7}{8}c\,\hat{\mathbf{z}}$	(16*f*)	H_2O
$\mathbf{B}_{25}$	=	$\left(\frac{5}{8}-x_4\right)\mathbf{a}_1+\left(\frac{7}{8}-x_4\right)\mathbf{a}_2+\left(\frac{3}{4}-2x_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-x_4\right)a\,\hat{\mathbf{y}}+\frac{3}{8}c\,\hat{\mathbf{z}}$	(16*f*)	H_2O
$\mathbf{B}_{26}$	=	$\left(\frac{5}{8}+x_4\right)\mathbf{a}_1+\left(\frac{7}{8}+x_4\right)\mathbf{a}_2+\left(\frac{3}{4}+2x_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+x_4\right)a\,\hat{\mathbf{y}}+\frac{3}{8}c\,\hat{\mathbf{z}}$	(16*f*)	H_2O
$\mathbf{B}_{27}$	=	$\left(\frac{7}{8}-x_4\right)\mathbf{a}_1+\left(\frac{5}{8}+x_4\right)\mathbf{a}_2+\frac{1}{4}\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-x_4\right)a\,\hat{\mathbf{y}}+\frac{5}{8}c\,\hat{\mathbf{z}}$	(16*f*)	H_2O
$\mathbf{B}_{28}$	=	$\left(\frac{7}{8}+x_4\right)\mathbf{a}_1+\left(\frac{5}{8}-x_4\right)\mathbf{a}_2+\frac{1}{4}\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+x_4\right)a\,\hat{\mathbf{y}}+\frac{5}{8}c\,\hat{\mathbf{z}}$	(16*f*)	H_2O
$\mathbf{B}_{29}$	=	$(y_5+z_5)\,\mathbf{a}_1+(x_5+z_5)\,\mathbf{a}_2+(x_5+y_5)\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}}+y_5a\,\hat{\mathbf{y}}+z_5c\,\hat{\mathbf{z}}$	(32*g*)	O I
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2}-y_5+z_5\right)\mathbf{a}_1+(-x_5+z_5)\,\mathbf{a}_2+\left(\frac{1}{2}-x_5-y_5\right)\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_5\right)a\,\hat{\mathbf{y}}+z_5c\,\hat{\mathbf{z}}$	(32*g*)	O I
$\mathbf{B}_{31}$	=	$(x_5+z_5)\,\mathbf{a}_1+\left(\frac{1}{2}-y_5+z_5\right)\mathbf{a}_2+(x_5-y_5)\,\mathbf{a}_3$	=	$\left(\frac{1}{4}-y_5\right)a\,\hat{\mathbf{x}}+\left(\frac{3}{4}+x_5\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+z_5\right)c\,\hat{\mathbf{z}}$	(32*g*)	O I
$\mathbf{B}_{32}$	=	$(-x_5+z_5)\,\mathbf{a}_1+(y_5+z_5)\,\mathbf{a}_2+\left(\frac{1}{2}-x_5+y_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}+y_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-x_5\right)a\,\hat{\mathbf{y}}+\left(\frac{3}{4}+z_5\right)c\,\hat{\mathbf{z}}$	(32*g*)	O I
$\mathbf{B}_{33}$	=	$(y_5-z_5)\,\mathbf{a}_1+\left(\frac{1}{2}-x_5-z_5\right)\mathbf{a}_2+\left(\frac{1}{2}-x_5+y_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{x}}+y_5a\,\hat{\mathbf{y}}-z_5c\,\hat{\mathbf{z}}$	(32*g*)	O I
$\mathbf{B}_{34}$	=	$\left(\frac{1}{2}-y_5-z_5\right)\mathbf{a}_1+\left(\frac{1}{2}+x_5-z_5\right)\mathbf{a}_2+(x_5-y_5)\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}}-y_5a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_5\right)c\,\hat{\mathbf{z}}$	(32*g*)	O I
$\mathbf{B}_{35}$	=	$\left(\frac{1}{2}+x_5-z_5\right)\mathbf{a}_1+(y_5-z_5)\,\mathbf{a}_2+(x_5+y_5)\,\mathbf{a}_3$	=	$\left(-\frac{1}{4}+y_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+x_5\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-z_5\right)c\,\hat{\mathbf{z}}$	(32*g*)	O I
$\mathbf{B}_{36}$	=	$\left(\frac{1}{2}-x_5-z_5\right)\mathbf{a}_1+\left(\frac{1}{2}-y_5-z_5\right)\mathbf{a}_2+\left(\frac{1}{2}-x_5-y_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}-y_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-x_5\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-z_5\right)c\,\hat{\mathbf{z}}$	(32*g*)	O I
$\mathbf{B}_{37}$	=	$(-y_5-z_5)\,\mathbf{a}_1+(-x_5-z_5)\,\mathbf{a}_2+(-x_5-y_5)\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}}-y_5a\,\hat{\mathbf{y}}-z_5c\,\hat{\mathbf{z}}$	(32*g*)	O I
$\mathbf{B}_{38}$	=	$\left(\frac{1}{2}+y_5-z_5\right)\mathbf{a}_1+(x_5-z_5)\,\mathbf{a}_2+\left(\frac{1}{2}+x_5+y_5\right)\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_5\right)a\,\hat{\mathbf{y}}-z_5c\,\hat{\mathbf{z}}$	(32*g*)	O I
$\mathbf{B}_{39}$	=	$(-x_5-z_5)\,\mathbf{a}_1+\left(\frac{1}{2}+y_5-z_5\right)\mathbf{a}_2+(-x_5+y_5)\,\mathbf{a}_3$	=	$\left(\frac{1}{4}+y_5\right)a\,\hat{\mathbf{x}}-\left(x_5+\frac{1}{4}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-z_5\right)c\,\hat{\mathbf{z}}$	(32*g*)	O I
$\mathbf{B}_{40}$	=	$(x_5-z_5)\,\mathbf{a}_1+(-y_5-z_5)\,\mathbf{a}_2+\left(\frac{1}{2}+x_5-y_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}-y_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+x_5\right)a\,\hat{\mathbf{y}}-\left(z_5+\frac{1}{4}\right)c\,\hat{\mathbf{z}}$	(32*g*)	O I
$\mathbf{B}_{41}$	=	$(-y_5+z_5)\,\mathbf{a}_1+\left(\frac{1}{2}+x_5+z_5\right)\mathbf{a}_2+\left(\frac{1}{2}+x_5-y_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{x}}-y_5a\,\hat{\mathbf{y}}+z_5c\,\hat{\mathbf{z}}$	(32*g*)	O I
$\mathbf{B}_{42}$	=	$\left(\frac{1}{2}+y_5+z_5\right)\mathbf{a}_1+\left(\frac{1}{2}-x_5+z_5\right)\mathbf{a}_2+(-x_5+y_5)\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}}+y_5a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	(32*g*)	O I
$\mathbf{B}_{43}$	=	$\left(\frac{1}{2}-x_5+z_5\right)\mathbf{a}_1+(-y_5+z_5)\,\mathbf{a}_2+(-x_5-y_5)\,\mathbf{a}_3$	=	$\left(\frac{3}{4}-y_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-x_5\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+z_5\right)c\,\hat{\mathbf{z}}$	(32*g*)	O I
$\mathbf{B}_{44}$	=	$\left(\frac{1}{2}+x_5+z_5\right)\mathbf{a}_1+\left(\frac{1}{2}+y_5+z_5\right)\mathbf{a}_2+\left(\frac{1}{2}+x_5+y_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}+y_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+x_5\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+z_5\right)c\,\hat{\mathbf{z}}$	(32*g*)	O I

$$\mathbf{B}_{45} = (y_6 + z_6)\,\mathbf{a}_1 + (x_6 + z_6)\,\mathbf{a}_2 + (x_6 + y_6)\,\mathbf{a}_3 = x_6 a\,\hat{\mathbf{x}} + y_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}} \quad (32g) \quad \text{O II}$$

$$\mathbf{B}_{46} = \left(\tfrac{1}{2} - y_6 + z_6\right)\mathbf{a}_1 + (-x_6 + z_6)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_6 - y_6\right)\mathbf{a}_3 = -x_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_6\right) a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}} \quad (32g) \quad \text{O II}$$

$$\mathbf{B}_{47} = (x_6 + z_6)\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_6 + z_6\right)\mathbf{a}_2 + (x_6 - y_6)\,\mathbf{a}_3 = \left(\tfrac{1}{4} - y_6\right) a\,\hat{\mathbf{x}} + \left(\tfrac{3}{4} + x_6\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_6\right) c\,\hat{\mathbf{z}} \quad (32g) \quad \text{O II}$$

$$\mathbf{B}_{48} = (-x_6 + z_6)\,\mathbf{a}_1 + (y_6 + z_6)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_6 + y_6\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + y_6\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_6\right) a\,\hat{\mathbf{y}} + \left(\tfrac{3}{4} + z_6\right) c\,\hat{\mathbf{z}} \quad (32g) \quad \text{O II}$$

$$\mathbf{B}_{49} = (y_6 - z_6)\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_6 - z_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_6 + y_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} + y_6 a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}} \quad (32g) \quad \text{O II}$$

$$\mathbf{B}_{50} = \left(\tfrac{1}{2} - y_6 - z_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_6 - z_6\right)\mathbf{a}_2 + (x_6 - y_6)\,\mathbf{a}_3 = x_6 a\,\hat{\mathbf{x}} - y_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_6\right) c\,\hat{\mathbf{z}} \quad (32g) \quad \text{O II}$$

$$\mathbf{B}_{51} = \left(\tfrac{1}{2} + x_6 - z_6\right)\mathbf{a}_1 + (y_6 - z_6)\,\mathbf{a}_2 + (x_6 + y_6)\,\mathbf{a}_3 = \left(-\tfrac{1}{4} + y_6\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_6\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - z_6\right) c\,\hat{\mathbf{z}} \quad (32g) \quad \text{O II}$$

$$\mathbf{B}_{52} = \left(\tfrac{1}{2} - x_6 - z_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_6 - z_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_6 - y_6\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - y_6\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_6\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - z_6\right) c\,\hat{\mathbf{z}} \quad (32g) \quad \text{O II}$$

$$\mathbf{B}_{53} = (-y_6 - z_6)\,\mathbf{a}_1 + (-x_6 - z_6)\,\mathbf{a}_2 + (-x_6 - y_6)\,\mathbf{a}_3 = -x_6 a\,\hat{\mathbf{x}} - y_6 a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}} \quad (32g) \quad \text{O II}$$

$$\mathbf{B}_{54} = \left(\tfrac{1}{2} + y_6 - z_6\right)\mathbf{a}_1 + (x_6 - z_6)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_6 + y_6\right)\mathbf{a}_3 = x_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_6\right) a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}} \quad (32g) \quad \text{O II}$$

$$\mathbf{B}_{55} = (-x_6 - z_6)\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_6 - z_6\right)\mathbf{a}_2 + (-x_6 + y_6)\,\mathbf{a}_3 = \left(\tfrac{1}{4} + y_6\right) a\,\hat{\mathbf{x}} - \left(x_6 + \tfrac{1}{4}\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - z_6\right) c\,\hat{\mathbf{z}} \quad (32g) \quad \text{O II}$$

$$\mathbf{B}_{56} = (x_6 - z_6)\,\mathbf{a}_1 + (-y_6 - z_6)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_6 - y_6\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - y_6\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_6\right) a\,\hat{\mathbf{y}} - \left(z_6 + \tfrac{1}{4}\right) c\,\hat{\mathbf{z}} \quad (32g) \quad \text{O II}$$

$$\mathbf{B}_{57} = (-y_6 + z_6)\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_6 + z_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_6 - y_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} - y_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}} \quad (32g) \quad \text{O II}$$

$$\mathbf{B}_{58} = \left(\tfrac{1}{2} + y_6 + z_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_6 + z_6\right)\mathbf{a}_2 + (-x_6 + y_6)\,\mathbf{a}_3 = -x_6 a\,\hat{\mathbf{x}} + y_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right) c\,\hat{\mathbf{z}} \quad (32g) \quad \text{O II}$$

$$\mathbf{B}_{59} = \left(\tfrac{1}{2} - x_6 + z_6\right)\mathbf{a}_1 + (-y_6 + z_6)\,\mathbf{a}_2 + (-x_6 - y_6)\,\mathbf{a}_3 = \left(\tfrac{3}{4} - y_6\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_6\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_6\right) c\,\hat{\mathbf{z}} \quad (32g) \quad \text{O II}$$

$$\mathbf{B}_{60} = \left(\tfrac{1}{2} + x_6 + z_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_6 + z_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_6 + y_6\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + y_6\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_6\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_6\right) c\,\hat{\mathbf{z}} \quad (32g) \quad \text{O II}$$

$$\mathbf{B}_{61} = (y_7 + z_7)\,\mathbf{a}_1 + (x_7 + z_7)\,\mathbf{a}_2 + (x_7 + y_7)\,\mathbf{a}_3 = x_7 a\,\hat{\mathbf{x}} + y_7 a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}} \quad (32g) \quad \text{O III}$$

$$\mathbf{B}_{62} = \left(\tfrac{1}{2} - y_7 + z_7\right)\mathbf{a}_1 + (-x_7 + z_7)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_7 - y_7\right)\mathbf{a}_3 = -x_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_7\right) a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}} \quad (32g) \quad \text{O III}$$

$$\mathbf{B}_{63} = (x_7 + z_7)\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_7 + z_7\right)\mathbf{a}_2 + (x_7 - y_7)\,\mathbf{a}_3 = \left(\tfrac{1}{4} - y_7\right) a\,\hat{\mathbf{x}} + \left(\tfrac{3}{4} + x_7\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_7\right) c\,\hat{\mathbf{z}} \quad (32g) \quad \text{O III}$$

$$\mathbf{B}_{64} = (-x_7 + z_7)\,\mathbf{a}_1 + (y_7 + z_7)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_7 + y_7\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + y_7\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_7\right) a\,\hat{\mathbf{y}} + \left(\tfrac{3}{4} + z_7\right) c\,\hat{\mathbf{z}} \quad (32g) \quad \text{O III}$$

$$\mathbf{B}_{65} = (y_7 - z_7)\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_7 - z_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_7 + y_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_7\right) a\,\hat{\mathbf{x}} + y_7 a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}} \quad (32g) \quad \text{O III}$$

$$\mathbf{B}_{66} = \left(\tfrac{1}{2} - y_7 - z_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_7 - z_7\right)\mathbf{a}_2 + (x_7 - y_7)\,\mathbf{a}_3 = x_7 a\,\hat{\mathbf{x}} - y_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_7\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{O III}$$

$$\mathbf{B}_{67} = \left(\tfrac{1}{2} + x_7 - z_7\right)\mathbf{a}_1 + (y_7 - z_7)\,\mathbf{a}_2 + (x_7 + y_7)\,\mathbf{a}_3 = \left(-\tfrac{1}{4} + y_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - z_7\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{O III}$$

$$\mathbf{B}_{68} = \left(\tfrac{1}{2} - x_7 - z_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_7 - z_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_7 - y_7\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - y_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - z_7\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{O III}$$

$$\mathbf{B}_{69} = (-y_7 - z_7)\,\mathbf{a}_1 + (-x_7 - z_7)\,\mathbf{a}_2 + (-x_7 - y_7)\,\mathbf{a}_3 = -x_7 a\,\hat{\mathbf{x}} - y_7 a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}} \quad (32g) \quad \text{O III}$$

$$\mathbf{B}_{70} = \left(\tfrac{1}{2} + y_7 - z_7\right)\mathbf{a}_1 + (x_7 - z_7)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_7 + y_7\right)\mathbf{a}_3 = x_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_7\right)a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}} \quad (32g) \quad \text{O III}$$

$$\mathbf{B}_{71} = (-x_7 - z_7)\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_7 - z_7\right)\mathbf{a}_2 + (-x_7 + y_7)\,\mathbf{a}_3 = \left(\tfrac{1}{4} + y_7\right)a\,\hat{\mathbf{x}} - \left(x_7 + \tfrac{1}{4}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - z_7\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{O III}$$

$$\mathbf{B}_{72} = (x_7 - z_7)\,\mathbf{a}_1 + (-y_7 - z_7)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_7 - y_7\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - y_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_7\right)a\,\hat{\mathbf{y}} - \left(z_7 + \tfrac{1}{4}\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{O III}$$

$$\mathbf{B}_{73} = (-y_7 + z_7)\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_7 + z_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_7 - y_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_7\right)a\,\hat{\mathbf{x}} - y_7 a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}} \quad (32g) \quad \text{O III}$$

$$\mathbf{B}_{74} = \left(\tfrac{1}{2} + y_7 + z_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_7 + z_7\right)\mathbf{a}_2 + (-x_7 + y_7)\,\mathbf{a}_3 = -x_7 a\,\hat{\mathbf{x}} + y_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_7\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{O III}$$

$$\mathbf{B}_{75} = \left(\tfrac{1}{2} - x_7 + z_7\right)\mathbf{a}_1 + (-y_7 + z_7)\,\mathbf{a}_2 + (-x_7 - y_7)\,\mathbf{a}_3 = \left(\tfrac{3}{4} - y_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_7\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{O III}$$

$$\mathbf{B}_{76} = \left(\tfrac{1}{2} + x_7 + z_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_7 + z_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_7 + y_7\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + y_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_7\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{O III}$$

$$\mathbf{B}_{77} = (y_8 + z_8)\,\mathbf{a}_1 + (x_8 + z_8)\,\mathbf{a}_2 + (x_8 + y_8)\,\mathbf{a}_3 = x_8 a\,\hat{\mathbf{x}} + y_8 a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Si}$$

$$\mathbf{B}_{78} = \left(\tfrac{1}{2} - y_8 + z_8\right)\mathbf{a}_1 + (-x_8 + z_8)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_8 - y_8\right)\mathbf{a}_3 = -x_8 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_8\right)a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Si}$$

$$\mathbf{B}_{79} = (x_8 + z_8)\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_8 + z_8\right)\mathbf{a}_2 + (x_8 - y_8)\,\mathbf{a}_3 = \left(\tfrac{1}{4} - y_8\right)a\,\hat{\mathbf{x}} + \left(\tfrac{3}{4} + x_8\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_8\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Si}$$

$$\mathbf{B}_{80} = (-x_8 + z_8)\,\mathbf{a}_1 + (y_8 + z_8)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_8 + y_8\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + y_8\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_8\right)a\,\hat{\mathbf{y}} + \left(\tfrac{3}{4} + z_8\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Si}$$

$$\mathbf{B}_{81} = (y_8 - z_8)\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_8 - z_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_8 + y_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_8\right)a\,\hat{\mathbf{x}} + y_8 a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Si}$$

$$\mathbf{B}_{82} = \left(\tfrac{1}{2} - y_8 - z_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_8 - z_8\right)\mathbf{a}_2 + (x_8 - y_8)\,\mathbf{a}_3 = x_8 a\,\hat{\mathbf{x}} - y_8 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_8\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Si}$$

$$\mathbf{B}_{83} = \left(\tfrac{1}{2} + x_8 - z_8\right)\mathbf{a}_1 + (y_8 - z_8)\,\mathbf{a}_2 + (x_8 + y_8)\,\mathbf{a}_3 = \left(-\tfrac{1}{4} + y_8\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_8\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - z_8\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Si}$$

$$\mathbf{B}_{84} = \left(\tfrac{1}{2} - x_8 - z_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_8 - z_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_8 - y_8\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - y_8\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_8\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - z_8\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Si}$$

$$\mathbf{B}_{85} = (-y_8 - z_8)\,\mathbf{a}_1 + (-x_8 - z_8)\,\mathbf{a}_2 + (-x_8 - y_8)\,\mathbf{a}_3 = -x_8 a\,\hat{\mathbf{x}} - y_8 a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Si}$$

$$\mathbf{B}_{86} = \left(\tfrac{1}{2} + y_8 - z_8\right)\mathbf{a}_1 + (x_8 - z_8)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_8 + y_8\right)\mathbf{a}_3 = x_8 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_8\right)a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Si}$$

$\mathbf{B}_{87}$	=	$(-x_8 - z_8)\,\mathbf{a}_1 + \left(\frac{1}{2} + y_8 - z_8\right)\mathbf{a}_2 + (-x_8 + y_8)\,\mathbf{a}_3$	=	$\left(\frac{1}{4} + y_8\right)a\,\hat{\mathbf{x}} - \left(x_8 + \frac{1}{4}\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_8\right)c\,\hat{\mathbf{z}}$	(32*g*)	Si
$\mathbf{B}_{88}$	=	$(x_8 - z_8)\,\mathbf{a}_1 + (-y_8 - z_8)\,\mathbf{a}_2 + \left(\frac{1}{2} + x_8 - y_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - y_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_8\right)a\,\hat{\mathbf{y}} - \left(z_8 + \frac{1}{4}\right)c\,\hat{\mathbf{z}}$	(32*g*)	Si
$\mathbf{B}_{89}$	=	$(-y_8 + z_8)\,\mathbf{a}_1 + \left(\frac{1}{2} + x_8 + z_8\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_8 - y_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_8\right)a\,\hat{\mathbf{x}} - y_8 a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	(32*g*)	Si
$\mathbf{B}_{90}$	=	$\left(\frac{1}{2} + y_8 + z_8\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_8 + z_8\right)\mathbf{a}_2 + (-x_8 + y_8)\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} + y_8 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right)c\,\hat{\mathbf{z}}$	(32*g*)	Si
$\mathbf{B}_{91}$	=	$\left(\frac{1}{2} - x_8 + z_8\right)\mathbf{a}_1 + (-y_8 + z_8)\,\mathbf{a}_2 + (-x_8 - y_8)\,\mathbf{a}_3$	=	$\left(\frac{3}{4} - y_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_8\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_8\right)c\,\hat{\mathbf{z}}$	(32*g*)	Si
$\mathbf{B}_{92}$	=	$\left(\frac{1}{2} + x_8 + z_8\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_8 + z_8\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_8 + y_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + y_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_8\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_8\right)c\,\hat{\mathbf{z}}$	(32*g*)	Si

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- W. Hartwig, *Zur Strukturbestimmung des Analcims*, Zeitschrift für Kristallographie - Crystalline Materials **78**, 173–207 (1931), doi:10.1524/zkri.1931.78.1.173.
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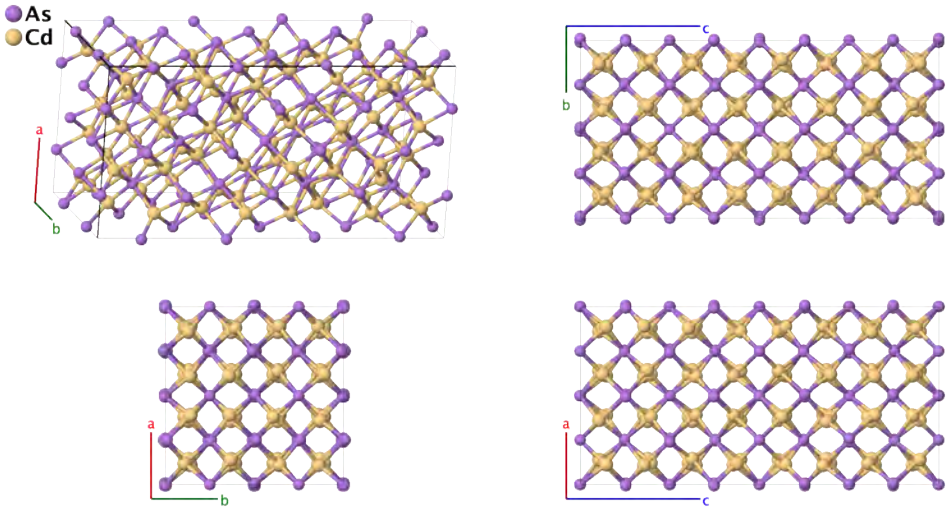
- F. Pechar, *The crystal structure of natural monoclinic analcime ($NaAlSi_2O_6{\cdot}H_2O$)*, Zeitschrift für Kristallographie - Crystalline Materials **184**, 63–70 (1988), doi:10.1524/zkri.1988.184.14.63.

Geometry files:

- CIF: pp. 1716
- POSCAR: pp. 1717

Cd_3As_2 Structure: A2B3_tI160_142_deg_3g

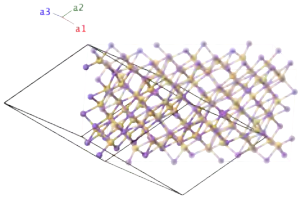
http://aflow.org/prototype-encyclopedia/A2B3_tI160_142_deg_3g



Prototype	:	As_2Cd_3
AFLOW prototype label	:	A2B3_tI160_142_deg_3g
***Strukturbericht* designation**	:	None
Pearson symbol	:	tI160
Space group number	:	142
Space group symbol	:	$I4_1/acd$
AFLOW prototype command	:	`aflow --proto=A2B3_tI160_142_deg_3g --params=`$a, c/a, z_1, x_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6$

Body-centered Tetragonal primitive vectors:

$$\mathbf{a}_1 = -\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}a\,\hat{\mathbf{x}} - \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} - \tfrac{1}{2}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\left(\frac{1}{4}+z_1\right)\mathbf{a}_1 + z_1\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	$(16d)$	As I
$\mathbf{B}_2$	=	$z_1\,\mathbf{a}_1 + \left(\frac{1}{4}+z_1\right)\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \left(-\frac{1}{4}+z_1\right)c\,\hat{\mathbf{z}}$	$(16d)$	As I
$\mathbf{B}_3$	=	$\left(\frac{1}{4}-z_1\right)\mathbf{a}_1 + \left(\frac{1}{2}-z_1\right)\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	$(16d)$	As I
$\mathbf{B}_4$	=	$\left(\frac{1}{2}-z_1\right)\mathbf{a}_1 + \left(\frac{1}{4}-z_1\right)\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{y}} + \left(\frac{1}{4}-z_1\right)c\,\hat{\mathbf{z}}$	$(16d)$	As I
$\mathbf{B}_5$	=	$\left(\frac{3}{4}-z_1\right)\mathbf{a}_1 - z_1\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	$(16d)$	As I
$\mathbf{B}_6$	=	$-z_1\,\mathbf{a}_1 + \left(\frac{3}{4}-z_1\right)\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{4}a\,\hat{\mathbf{y}} + \left(\frac{1}{4}-z_1\right)c\,\hat{\mathbf{z}}$	$(16d)$	As I
$\mathbf{B}_7$	=	$\left(\frac{3}{4}+z_1\right)\mathbf{a}_1 + \left(\frac{1}{2}+z_1\right)\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_1\right)c\,\hat{\mathbf{z}}$	$(16d)$	As I

$\mathbf{B}_8$	=	$\left(\frac{1}{2}+z_1\right)\mathbf{a}_1+\left(\frac{3}{4}+z_1\right)\mathbf{a}_2+\frac{3}{4}\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}+\frac{1}{4}a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+z_1\right)c\,\hat{\mathbf{z}}$	(16*d*)	As I
$\mathbf{B}_9$	=	$\frac{1}{4}\mathbf{a}_1+\left(\frac{1}{4}+x_2\right)\mathbf{a}_2+x_2\,\mathbf{a}_3$	=	$x_2a\,\hat{\mathbf{x}}+\frac{1}{4}c\,\hat{\mathbf{z}}$	(16*e*)	As II
$\mathbf{B}_{10}$	=	$\frac{3}{4}\mathbf{a}_1+\left(\frac{1}{4}-x_2\right)\mathbf{a}_2+\left(\frac{1}{2}-x_2\right)\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}}+\frac{1}{2}a\,\hat{\mathbf{y}}+\frac{1}{4}c\,\hat{\mathbf{z}}$	(16*e*)	As II
$\mathbf{B}_{11}$	=	$\left(\frac{1}{4}+x_2\right)\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2+x_2\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}}+\left(\frac{3}{4}+x_2\right)a\,\hat{\mathbf{y}}+\frac{1}{2}c\,\hat{\mathbf{z}}$	(16*e*)	As II
$\mathbf{B}_{12}$	=	$\left(\frac{1}{4}-x_2\right)\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+\left(\frac{1}{2}-x_2\right)\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-x_2\right)a\,\hat{\mathbf{y}}$	(16*e*)	As II
$\mathbf{B}_{13}$	=	$\frac{3}{4}\mathbf{a}_1+\left(\frac{3}{4}-x_2\right)\mathbf{a}_2-x_2\,\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}}+\frac{3}{4}c\,\hat{\mathbf{z}}$	(16*e*)	As II
$\mathbf{B}_{14}$	=	$\frac{1}{4}\mathbf{a}_1+\left(\frac{3}{4}+x_2\right)\mathbf{a}_2+\left(\frac{1}{2}+x_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_2\right)a\,\hat{\mathbf{x}}+\frac{1}{4}c\,\hat{\mathbf{z}}$	(16*e*)	As II
$\mathbf{B}_{15}$	=	$\left(\frac{3}{4}-x_2\right)\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2-x_2\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-x_2\right)a\,\hat{\mathbf{y}}+\frac{1}{2}c\,\hat{\mathbf{z}}$	(16*e*)	As II
$\mathbf{B}_{16}$	=	$\left(\frac{3}{4}+x_2\right)\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2+\left(\frac{1}{2}+x_2\right)\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+x_2\right)a\,\hat{\mathbf{y}}+\frac{1}{2}c\,\hat{\mathbf{z}}$	(16*e*)	As II
$\mathbf{B}_{17}$	=	$(y_3+z_3)\,\mathbf{a}_1+(x_3+z_3)\,\mathbf{a}_2+(x_3+y_3)\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}}+y_3a\,\hat{\mathbf{y}}+z_3c\,\hat{\mathbf{z}}$	(32*g*)	As III
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2}-y_3+z_3\right)\mathbf{a}_1+(-x_3+z_3)\,\mathbf{a}_2+\left(\frac{1}{2}-x_3-y_3\right)\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_3\right)a\,\hat{\mathbf{y}}+z_3c\,\hat{\mathbf{z}}$	(32*g*)	As III
$\mathbf{B}_{19}$	=	$(x_3+z_3)\,\mathbf{a}_1+\left(\frac{1}{2}-y_3+z_3\right)\mathbf{a}_2+(x_3-y_3)\,\mathbf{a}_3$	=	$\left(\frac{1}{4}-y_3\right)a\,\hat{\mathbf{x}}+\left(\frac{3}{4}+x_3\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+z_3\right)c\,\hat{\mathbf{z}}$	(32*g*)	As III
$\mathbf{B}_{20}$	=	$(-x_3+z_3)\,\mathbf{a}_1+(y_3+z_3)\,\mathbf{a}_2+\left(\frac{1}{2}-x_3+y_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}+y_3\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-x_3\right)a\,\hat{\mathbf{y}}+\left(\frac{3}{4}+z_3\right)c\,\hat{\mathbf{z}}$	(32*g*)	As III
$\mathbf{B}_{21}$	=	$(y_3-z_3)\,\mathbf{a}_1+\left(\frac{1}{2}-x_3-z_3\right)\mathbf{a}_2+\left(\frac{1}{2}-x_3+y_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{x}}+y_3a\,\hat{\mathbf{y}}-z_3c\,\hat{\mathbf{z}}$	(32*g*)	As III
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2}-y_3-z_3\right)\mathbf{a}_1+\left(\frac{1}{2}+x_3-z_3\right)\mathbf{a}_2+(x_3-y_3)\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}}-y_3a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_3\right)c\,\hat{\mathbf{z}}$	(32*g*)	As III
$\mathbf{B}_{23}$	=	$\left(\frac{1}{2}+x_3-z_3\right)\mathbf{a}_1+(y_3-z_3)\,\mathbf{a}_2+(x_3+y_3)\,\mathbf{a}_3$	=	$\left(-\frac{1}{4}+y_3\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+x_3\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-z_3\right)c\,\hat{\mathbf{z}}$	(32*g*)	As III
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2}-x_3-z_3\right)\mathbf{a}_1+\left(\frac{1}{2}-y_3-z_3\right)\mathbf{a}_2+\left(\frac{1}{2}-x_3-y_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}-y_3\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-x_3\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-z_3\right)c\,\hat{\mathbf{z}}$	(32*g*)	As III
$\mathbf{B}_{25}$	=	$(-y_3-z_3)\,\mathbf{a}_1+(-x_3-z_3)\,\mathbf{a}_2+(-x_3-y_3)\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}}-y_3a\,\hat{\mathbf{y}}-z_3c\,\hat{\mathbf{z}}$	(32*g*)	As III
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2}+y_3-z_3\right)\mathbf{a}_1+(x_3-z_3)\,\mathbf{a}_2+\left(\frac{1}{2}+x_3+y_3\right)\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_3\right)a\,\hat{\mathbf{y}}-z_3c\,\hat{\mathbf{z}}$	(32*g*)	As III
$\mathbf{B}_{27}$	=	$(-x_3-z_3)\,\mathbf{a}_1+\left(\frac{1}{2}+y_3-z_3\right)\mathbf{a}_2+(-x_3+y_3)\,\mathbf{a}_3$	=	$\left(\frac{1}{4}+y_3\right)a\,\hat{\mathbf{x}}-\left(x_3+\frac{1}{4}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-z_3\right)c\,\hat{\mathbf{z}}$	(32*g*)	As III
$\mathbf{B}_{28}$	=	$(x_3-z_3)\,\mathbf{a}_1+(-y_3-z_3)\,\mathbf{a}_2+\left(\frac{1}{2}+x_3-y_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}-y_3\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+x_3\right)a\,\hat{\mathbf{y}}-\left(z_3+\frac{1}{4}\right)c\,\hat{\mathbf{z}}$	(32*g*)	As III
$\mathbf{B}_{29}$	=	$(-y_3+z_3)\,\mathbf{a}_1+\left(\frac{1}{2}+x_3+z_3\right)\mathbf{a}_2+\left(\frac{1}{2}+x_3-y_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{x}}-y_3a\,\hat{\mathbf{y}}+z_3c\,\hat{\mathbf{z}}$	(32*g*)	As III
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2}+y_3+z_3\right)\mathbf{a}_1+\left(\frac{1}{2}-x_3+z_3\right)\mathbf{a}_2+(-x_3+y_3)\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}}+y_3a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_3\right)c\,\hat{\mathbf{z}}$	(32*g*)	As III
$\mathbf{B}_{31}$	=	$\left(\frac{1}{2}-x_3+z_3\right)\mathbf{a}_1+(-y_3+z_3)\,\mathbf{a}_2+(-x_3-y_3)\,\mathbf{a}_3$	=	$\left(\frac{3}{4}-y_3\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-x_3\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+z_3\right)c\,\hat{\mathbf{z}}$	(32*g*)	As III

$$\mathbf{B}_{32} = \left(\tfrac{1}{2}+x_3+z_3\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_3+z_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+x_3+y_3\right)\mathbf{a}_3 = \left(\tfrac{1}{4}+y_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4}+x_3\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4}+z_3\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{As III}$$

$$\mathbf{B}_{33} = (y_4+z_4)\,\mathbf{a}_1 + (x_4+z_4)\,\mathbf{a}_2 + (x_4+y_4)\,\mathbf{a}_3 = x_4a\,\hat{\mathbf{x}} + y_4a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd I}$$

$$\mathbf{B}_{34} = \left(\tfrac{1}{2}-y_4+z_4\right)\mathbf{a}_1 + (-x_4+z_4)\,\mathbf{a}_2 + \left(\tfrac{1}{2}-x_4-y_4\right)\mathbf{a}_3 = -x_4a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_4\right)a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd I}$$

$$\mathbf{B}_{35} = (x_4+z_4)\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_4+z_4\right)\mathbf{a}_2 + (x_4-y_4)\,\mathbf{a}_3 = \left(\tfrac{1}{4}-y_4\right)a\,\hat{\mathbf{x}} + \left(\tfrac{3}{4}+x_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4}+z_4\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd I}$$

$$\mathbf{B}_{36} = (-x_4+z_4)\,\mathbf{a}_1 + (y_4+z_4)\,\mathbf{a}_2 + \left(\tfrac{1}{2}-x_4+y_4\right)\mathbf{a}_3 = \left(\tfrac{1}{4}+y_4\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4}-x_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{3}{4}+z_4\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd I}$$

$$\mathbf{B}_{37} = (y_4-z_4)\,\mathbf{a}_1 + \left(\tfrac{1}{2}-x_4-z_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-x_4+y_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_4\right)a\,\hat{\mathbf{x}} + y_4a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd I}$$

$$\mathbf{B}_{38} = \left(\tfrac{1}{2}-y_4-z_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+x_4-z_4\right)\mathbf{a}_2 + (x_4-y_4)\,\mathbf{a}_3 = x_4a\,\hat{\mathbf{x}} - y_4a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_4\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd I}$$

$$\mathbf{B}_{39} = \left(\tfrac{1}{2}+x_4-z_4\right)\mathbf{a}_1 + (y_4-z_4)\,\mathbf{a}_2 + (x_4+y_4)\,\mathbf{a}_3 = \left(-\tfrac{1}{4}+y_4\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4}+x_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4}-z_4\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd I}$$

$$\mathbf{B}_{40} = \left(\tfrac{1}{2}-x_4-z_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_4-z_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-x_4-y_4\right)\mathbf{a}_3 = \left(\tfrac{1}{4}-y_4\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4}-x_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4}-z_4\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd I}$$

$$\mathbf{B}_{41} = (-y_4-z_4)\,\mathbf{a}_1 + (-x_4-z_4)\,\mathbf{a}_2 + (-x_4-y_4)\,\mathbf{a}_3 = -x_4a\,\hat{\mathbf{x}} - y_4a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd I}$$

$$\mathbf{B}_{42} = \left(\tfrac{1}{2}+y_4-z_4\right)\mathbf{a}_1 + (x_4-z_4)\,\mathbf{a}_2 + \left(\tfrac{1}{2}+x_4+y_4\right)\mathbf{a}_3 = x_4a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_4\right)a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd I}$$

$$\mathbf{B}_{43} = (-x_4-z_4)\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_4-z_4\right)\mathbf{a}_2 + (-x_4+y_4)\,\mathbf{a}_3 = \left(\tfrac{1}{4}+y_4\right)a\,\hat{\mathbf{x}} - \left(x_4+\tfrac{1}{4}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4}-z_4\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd I}$$

$$\mathbf{B}_{44} = (x_4-z_4)\,\mathbf{a}_1 + (-y_4-z_4)\,\mathbf{a}_2 + \left(\tfrac{1}{2}+x_4-y_4\right)\mathbf{a}_3 = \left(\tfrac{1}{4}-y_4\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4}+x_4\right)a\,\hat{\mathbf{y}} - \left(z_4+\tfrac{1}{4}\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd I}$$

$$\mathbf{B}_{45} = (-y_4+z_4)\,\mathbf{a}_1 + \left(\tfrac{1}{2}+x_4+z_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+x_4-y_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_4\right)a\,\hat{\mathbf{x}} - y_4a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd I}$$

$$\mathbf{B}_{46} = \left(\tfrac{1}{2}+y_4+z_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-x_4+z_4\right)\mathbf{a}_2 + (-x_4+y_4)\,\mathbf{a}_3 = -x_4a\,\hat{\mathbf{x}} + y_4a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_4\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd I}$$

$$\mathbf{B}_{47} = \left(\tfrac{1}{2}-x_4+z_4\right)\mathbf{a}_1 + (-y_4+z_4)\,\mathbf{a}_2 + (-x_4-y_4)\,\mathbf{a}_3 = \left(\tfrac{3}{4}-y_4\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4}-x_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4}+z_4\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd I}$$

$$\mathbf{B}_{48} = \left(\tfrac{1}{2}+x_4+z_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_4+z_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+x_4+y_4\right)\mathbf{a}_3 = \left(\tfrac{1}{4}+y_4\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4}+x_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4}+z_4\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd I}$$

$$\mathbf{B}_{49} = (y_5+z_5)\,\mathbf{a}_1 + (x_5+z_5)\,\mathbf{a}_2 + (x_5+y_5)\,\mathbf{a}_3 = x_5a\,\hat{\mathbf{x}} + y_5a\,\hat{\mathbf{y}} + z_5c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd II}$$

$$\mathbf{B}_{50} = \left(\tfrac{1}{2}-y_5+z_5\right)\mathbf{a}_1 + (-x_5+z_5)\,\mathbf{a}_2 + \left(\tfrac{1}{2}-x_5-y_5\right)\mathbf{a}_3 = -x_5a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_5\right)a\,\hat{\mathbf{y}} + z_5c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd II}$$

$$\mathbf{B}_{51} = (x_5+z_5)\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_5+z_5\right)\mathbf{a}_2 + (x_5-y_5)\,\mathbf{a}_3 = \left(\tfrac{1}{4}-y_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{3}{4}+x_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4}+z_5\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd II}$$

$$\mathbf{B}_{52} = (-x_5+z_5)\,\mathbf{a}_1 + (y_5+z_5)\,\mathbf{a}_2 + \left(\tfrac{1}{2}-x_5+y_5\right)\mathbf{a}_3 = \left(\tfrac{1}{4}+y_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4}-x_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{3}{4}+z_5\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd II}$$

$$\mathbf{B}_{53} = (y_5 - z_5)\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_5 - z_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_5 + y_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + y_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd II}$$

$$\mathbf{B}_{54} = \left(\tfrac{1}{2} - y_5 - z_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_5 - z_5\right)\mathbf{a}_2 + (x_5 - y_5)\,\mathbf{a}_3 = x_5 a\,\hat{\mathbf{x}} - y_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_5\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd II}$$

$$\mathbf{B}_{55} = \left(\tfrac{1}{2} + x_5 - z_5\right)\mathbf{a}_1 + (y_5 - z_5)\,\mathbf{a}_2 + (x_5 + y_5)\,\mathbf{a}_3 = \left(-\tfrac{1}{4} + y_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - z_5\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd II}$$

$$\mathbf{B}_{56} = \left(\tfrac{1}{2} - x_5 - z_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_5 - z_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_5 - y_5\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - y_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - z_5\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd II}$$

$$\mathbf{B}_{57} = (-y_5 - z_5)\,\mathbf{a}_1 + (-x_5 - z_5)\,\mathbf{a}_2 + (-x_5 - y_5)\,\mathbf{a}_3 = -x_5 a\,\hat{\mathbf{x}} - y_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd II}$$

$$\mathbf{B}_{58} = \left(\tfrac{1}{2} + y_5 - z_5\right)\mathbf{a}_1 + (x_5 - z_5)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_5 + y_5\right)\mathbf{a}_3 = x_5 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_5\right)a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd II}$$

$$\mathbf{B}_{59} = (-x_5 - z_5)\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_5 - z_5\right)\mathbf{a}_2 + (-x_5 + y_5)\,\mathbf{a}_3 = \left(\tfrac{1}{4} + y_5\right)a\,\hat{\mathbf{x}} - \left(x_5 + \tfrac{1}{4}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - z_5\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd II}$$

$$\mathbf{B}_{60} = (x_5 - z_5)\,\mathbf{a}_1 + (-y_5 - z_5)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_5 - y_5\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - y_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_5\right)a\,\hat{\mathbf{y}} - \left(z_5 + \tfrac{1}{4}\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd II}$$

$$\mathbf{B}_{61} = (-y_5 + z_5)\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_5 + z_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_5 - y_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} - y_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd II}$$

$$\mathbf{B}_{62} = \left(\tfrac{1}{2} + y_5 + z_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_5 + z_5\right)\mathbf{a}_2 + (-x_5 + y_5)\,\mathbf{a}_3 = -x_5 a\,\hat{\mathbf{x}} + y_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_5\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd II}$$

$$\mathbf{B}_{63} = \left(\tfrac{1}{2} - x_5 + z_5\right)\mathbf{a}_1 + (-y_5 + z_5)\,\mathbf{a}_2 + (-x_5 - y_5)\,\mathbf{a}_3 = \left(\tfrac{3}{4} - y_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_5\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd II}$$

$$\mathbf{B}_{64} = \left(\tfrac{1}{2} + x_5 + z_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_5 + z_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_5 + y_5\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + y_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_5\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd II}$$

$$\mathbf{B}_{65} = (y_6 + z_6)\,\mathbf{a}_1 + (x_6 + z_6)\,\mathbf{a}_2 + (x_6 + y_6)\,\mathbf{a}_3 = x_6 a\,\hat{\mathbf{x}} + y_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd III}$$

$$\mathbf{B}_{66} = \left(\tfrac{1}{2} - y_6 + z_6\right)\mathbf{a}_1 + (-x_6 + z_6)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_6 - y_6\right)\mathbf{a}_3 = -x_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_6\right)a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd III}$$

$$\mathbf{B}_{67} = (x_6 + z_6)\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_6 + z_6\right)\mathbf{a}_2 + (x_6 - y_6)\,\mathbf{a}_3 = \left(\tfrac{1}{4} - y_6\right)a\,\hat{\mathbf{x}} + \left(\tfrac{3}{4} + x_6\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_6\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd III}$$

$$\mathbf{B}_{68} = (-x_6 + z_6)\,\mathbf{a}_1 + (y_6 + z_6)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_6 + y_6\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + y_6\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_6\right)a\,\hat{\mathbf{y}} + \left(\tfrac{3}{4} + z_6\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd III}$$

$$\mathbf{B}_{69} = (y_6 - z_6)\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_6 - z_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_6 + y_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_6\right)a\,\hat{\mathbf{x}} + y_6 a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd III}$$

$$\mathbf{B}_{70} = \left(\tfrac{1}{2} - y_6 - z_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_6 - z_6\right)\mathbf{a}_2 + (x_6 - y_6)\,\mathbf{a}_3 = x_6 a\,\hat{\mathbf{x}} - y_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_6\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd III}$$

$$\mathbf{B}_{71} = \left(\tfrac{1}{2} + x_6 - z_6\right)\mathbf{a}_1 + (y_6 - z_6)\,\mathbf{a}_2 + (x_6 + y_6)\,\mathbf{a}_3 = \left(-\tfrac{1}{4} + y_6\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_6\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - z_6\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd III}$$

$$\mathbf{B}_{72} = \left(\tfrac{1}{2} - x_6 - z_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_6 - z_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_6 - y_6\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - y_6\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_6\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - z_6\right)c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd III}$$

$$\mathbf{B}_{73} = (-y_6 - z_6)\,\mathbf{a}_1 + (-x_6 - z_6)\,\mathbf{a}_2 + (-x_6 - y_6)\,\mathbf{a}_3 = -x_6 a\,\hat{\mathbf{x}} - y_6 a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}} \quad (32g) \quad \text{Cd III}$$

$\mathbf{B}_{74}$	=	$\left(\frac{1}{2}+y_6-z_6\right)\mathbf{a}_1+(x_6-z_6)\,\mathbf{a}_2+\left(\frac{1}{2}+x_6+y_6\right)\mathbf{a}_3$	=	$x_6a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_6\right)a\,\hat{\mathbf{y}}-z_6c\,\hat{\mathbf{z}}$	(32*g*)	Cd III
$\mathbf{B}_{75}$	=	$(-x_6-z_6)\,\mathbf{a}_1+\left(\frac{1}{2}+y_6-z_6\right)\mathbf{a}_2+(-x_6+y_6)\,\mathbf{a}_3$	=	$\left(\frac{1}{4}+y_6\right)a\,\hat{\mathbf{x}}-\left(x_6+\frac{1}{4}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-z_6\right)c\,\hat{\mathbf{z}}$	(32*g*)	Cd III
$\mathbf{B}_{76}$	=	$(x_6-z_6)\,\mathbf{a}_1+(-y_6-z_6)\,\mathbf{a}_2+\left(\frac{1}{2}+x_6-y_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}-y_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+x_6\right)a\,\hat{\mathbf{y}}-\left(z_6+\frac{1}{4}\right)c\,\hat{\mathbf{z}}$	(32*g*)	Cd III
$\mathbf{B}_{77}$	=	$(-y_6+z_6)\,\mathbf{a}_1+\left(\frac{1}{2}+x_6+z_6\right)\mathbf{a}_2+\left(\frac{1}{2}+x_6-y_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{x}}-y_6a\,\hat{\mathbf{y}}+z_6c\,\hat{\mathbf{z}}$	(32*g*)	Cd III
$\mathbf{B}_{78}$	=	$\left(\frac{1}{2}+y_6+z_6\right)\mathbf{a}_1+\left(\frac{1}{2}-x_6+z_6\right)\mathbf{a}_2+(-x_6+y_6)\,\mathbf{a}_3$	=	$-x_6a\,\hat{\mathbf{x}}+y_6a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	(32*g*)	Cd III
$\mathbf{B}_{79}$	=	$\left(\frac{1}{2}-x_6+z_6\right)\mathbf{a}_1+(-y_6+z_6)\,\mathbf{a}_2+(-x_6-y_6)\,\mathbf{a}_3$	=	$\left(\frac{3}{4}-y_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-x_6\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+z_6\right)c\,\hat{\mathbf{z}}$	(32*g*)	Cd III
$\mathbf{B}_{80}$	=	$\left(\frac{1}{2}+x_6+z_6\right)\mathbf{a}_1+\left(\frac{1}{2}+y_6+z_6\right)\mathbf{a}_2+\left(\frac{1}{2}+x_6+y_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}+y_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+x_6\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+z_6\right)c\,\hat{\mathbf{z}}$	(32*g*)	Cd III

References:

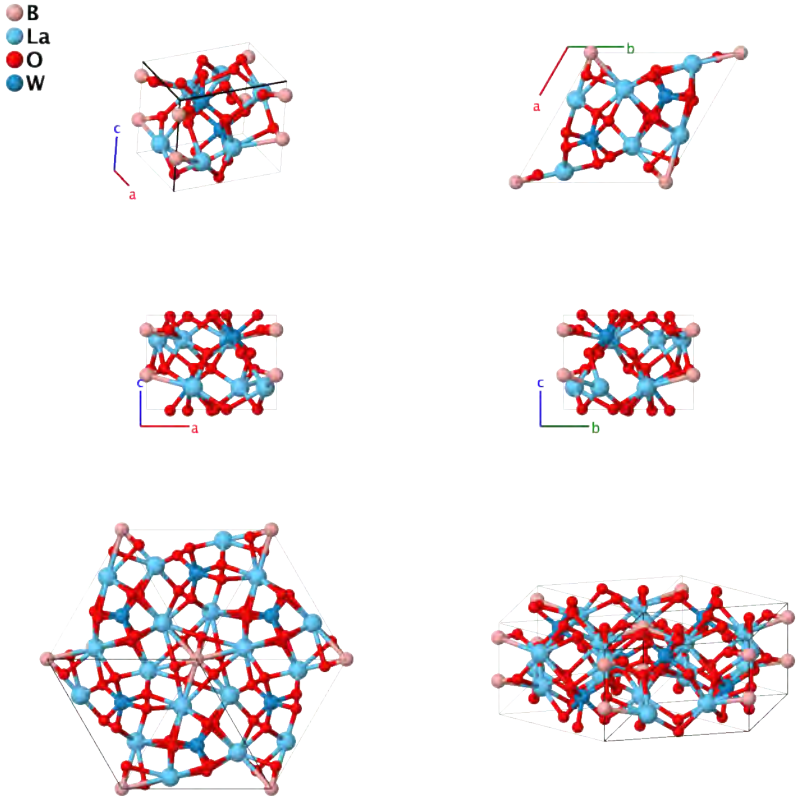
- M. N. Ali, Q. Gibson, S. Jeon, B. B. Zhou, A. Yazdani, and R. J. Cava, *The Crystal and Electronic Structures of Cd_3As_2, the Three-Dimensional Electronic Analogue of Graphene*, Inorg. Chem. **53**, 4062–4067 (2014), doi:10.1021/ic403163d.

Geometry files:

- CIF: pp. 1717
- POSCAR: pp. 1718

La_3BWO_9 ($P3$) Structure: AB3C9D_hP28_143_2a_2d_6d_bc

http://aflow.org/prototype-encyclopedia/AB3C9D_hP28_143_2a_2d_6d_bc



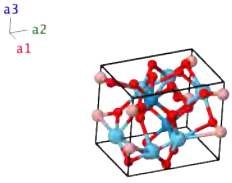
Prototype	:	BLa_3O_9W
AFLOW prototype label	:	AB3C9D_hP28_143_2a_2d_6d_bc
***Strukturbericht* designation**	:	None
Pearson symbol	:	hP28
Space group number	:	143
Space group symbol	:	$P3$
AFLOW prototype command	:	`aflow --proto=AB3C9D_hP28_143_2a_2d_6d_bc --params=`$a, c/a, z_1, z_2, z_3, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}$

- Most refinements of the BLa_3O_9W structure, including (Ashtar, 2020) place it in hexagonal space group $P6_3$ #173. (Han, 2018) find a better fit to the data by refining it in the trigonal $P3$ #143 space group, which places the lanthanum atoms on two independent crystallographic sites. As this may be due to the presence of bismuth impurities, which (Han, 2018) place on the lanthanum site, while (Ashtar, 2020) claim to have very pure samples, we withhold judgment on which structure is correct and present both.
- Space group $P3$ does not specify the origin of the z-axis. Here it is set so that the coordinate of the ($1b$) tungsten atom is $z_3 = 1/4$.
- (Han, 2018) mislabel the Wyckoff positions of the boron and tungsten atoms. We show the correct positions based on

the coordinates in their table, and the correspondence with the $P6_3$ structure.

Trigonal Hexagonal primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$z_1\,\mathbf{a}_3$	=	$z_1 c\,\hat{\mathbf{z}}$	$(1a)$	B I
$\mathbf{B}_2$	=	$z_2\,\mathbf{a}_3$	=	$z_2 c\,\hat{\mathbf{z}}$	$(1a)$	B II
$\mathbf{B}_3$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(1b)$	W I
$\mathbf{B}_4$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(1c)$	W II
$\mathbf{B}_5$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_5 + y_5)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_5 + y_5)\,a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(3d)$	La I
$\mathbf{B}_6$	=	$-y_5\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_5 - y_5\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(3d)$	La I
$\mathbf{B}_7$	=	$(-x_5 + y_5)\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\left(-x_5 + \frac{1}{2}y_5\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(3d)$	La I
$\mathbf{B}_8$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_6 + y_6)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_6 + y_6)\,a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(3d)$	La II
$\mathbf{B}_9$	=	$-y_6\,\mathbf{a}_1 + (x_6 - y_6)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_6 - y_6\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(3d)$	La II
$\mathbf{B}_{10}$	=	$(-x_6 + y_6)\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\left(-x_6 + \frac{1}{2}y_6\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(3d)$	La II
$\mathbf{B}_{11}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_7 + y_7)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_7 + y_7)\,a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	$(3d)$	O I
$\mathbf{B}_{12}$	=	$-y_7\,\mathbf{a}_1 + (x_7 - y_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_7 - y_7\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_7 a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	$(3d)$	O I
$\mathbf{B}_{13}$	=	$(-x_7 + y_7)\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$\left(-x_7 + \frac{1}{2}y_7\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_7 a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	$(3d)$	O I
$\mathbf{B}_{14}$	=	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_8 + y_8)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_8 + y_8)\,a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	$(3d)$	O II
$\mathbf{B}_{15}$	=	$-y_8\,\mathbf{a}_1 + (x_8 - y_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_8 - y_8\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_8 a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	$(3d)$	O II
$\mathbf{B}_{16}$	=	$(-x_8 + y_8)\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$\left(-x_8 + \frac{1}{2}y_8\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_8 a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	$(3d)$	O II
$\mathbf{B}_{17}$	=	$x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_9 + y_9)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_9 + y_9)\,a\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	$(3d)$	O III
$\mathbf{B}_{18}$	=	$-y_9\,\mathbf{a}_1 + (x_9 - y_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_9 - y_9\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_9 a\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	$(3d)$	O III
$\mathbf{B}_{19}$	=	$(-x_9 + y_9)\,\mathbf{a}_1 - x_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$\left(-x_9 + \frac{1}{2}y_9\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_9 a\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	$(3d)$	O III
$\mathbf{B}_{20}$	=	$x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_{10} + y_{10})\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_{10} + y_{10})\,a\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}}$	$(3d)$	O IV
$\mathbf{B}_{21}$	=	$-y_{10}\,\mathbf{a}_1 + (x_{10} - y_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_{10} - y_{10}\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_{10} a\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}}$	$(3d)$	O IV
$\mathbf{B}_{22}$	=	$(-x_{10} + y_{10})\,\mathbf{a}_1 - x_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$\left(-x_{10} + \frac{1}{2}y_{10}\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_{10} a\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}}$	$(3d)$	O IV
$\mathbf{B}_{23}$	=	$x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_{11} + y_{11})\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_{11} + y_{11})\,a\,\hat{\mathbf{y}} + z_{11} c\,\hat{\mathbf{z}}$	$(3d)$	O V
$\mathbf{B}_{24}$	=	$-y_{11}\,\mathbf{a}_1 + (x_{11} - y_{11})\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_{11} - y_{11}\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_{11} a\,\hat{\mathbf{y}} + z_{11} c\,\hat{\mathbf{z}}$	$(3d)$	O V

$$\mathbf{B}_{25} = (-x_{11}+y_{11})\,\mathbf{a}_1 - x_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 = \left(-x_{11}+\tfrac{1}{2}y_{11}\right)a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}y_{11}a\,\hat{\mathbf{y}} + z_{11}c\,\hat{\mathbf{z}} \quad (3d) \quad \text{O V}$$

$$\mathbf{B}_{26} = x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 = \tfrac{1}{2}(x_{12}+y_{12})\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}(-x_{12}+y_{12})\,a\,\hat{\mathbf{y}} + z_{12}c\,\hat{\mathbf{z}} \quad (3d) \quad \text{O VI}$$

$$\mathbf{B}_{27} = -y_{12}\,\mathbf{a}_1 + (x_{12}-y_{12})\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 = \left(\tfrac{1}{2}x_{12}-y_{12}\right)a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}x_{12}a\,\hat{\mathbf{y}} + z_{12}c\,\hat{\mathbf{z}} \quad (3d) \quad \text{O VI}$$

$$\mathbf{B}_{28} = (-x_{12}+y_{12})\,\mathbf{a}_1 - x_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3 = \left(-x_{12}+\tfrac{1}{2}y_{12}\right)a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}y_{12}a\,\hat{\mathbf{y}} + z_{12}c\,\hat{\mathbf{z}} \quad (3d) \quad \text{O VI}$$

References:

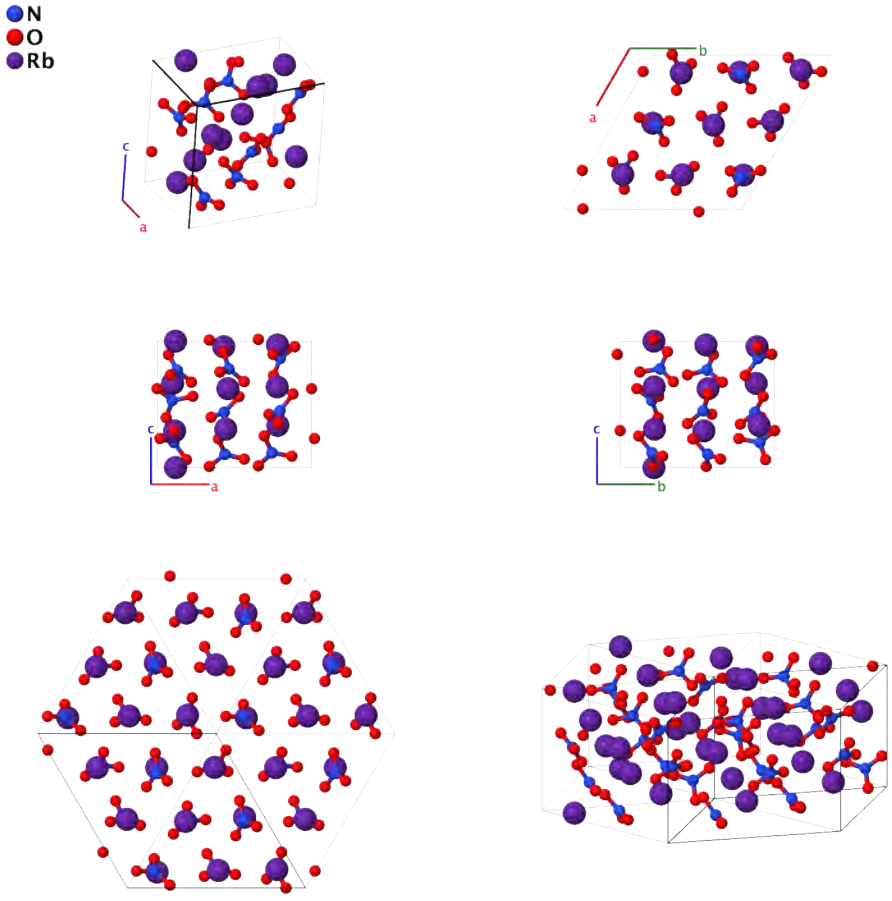
- J. Han, F. Pan, M. S. Molokeev, J. Dai, M. Peng, W. Zhou, and J. Wang, *Redefinition of Crystal Structure and Bi^{3+} Yellow Luminescence with Strong Near-Ultraviolet Excitation in La_3BWO_9:Bi^{3+} Phosphor for White Light-Emitting Diodes*, ACS Appl. Mater. Interfaces **10**, 13660–13668 (2018), doi:10.1021/acsami.8b00808.
- M. Ashtar, J. Guo, Z. Wan, Y. Wang, G. Gong, Y. Liu, Y. Su, and Z. Tian, *A new family of disorder-free Rare-Earth-based kagomé lattice magnets: structure and magnetic characterizations of RE_3BWO_9 (RE = Pr, Nd, Gd-Ho) Boratotungstates*, http://arxiv.org/abs/2002.05420. ArXiv:2002.05420 [cond-mat.mtrl-sci].

Geometry files:

- CIF: pp. 1718
- POSCAR: pp. 1719

$RbNO_3$ (IV) Structure: AB3C_hP45_144_3a_9a_3a

http://aflow.org/prototype-encyclopedia/AB3C_hP45_144_3a_9a_3a



Prototype	:	NO_3Rb
AFLOW prototype label	:	AB3C_hP45_144_3a_9a_3a
***Strukturbericht* designation**	:	None
Pearson symbol	:	hP45
Space group number	:	144
Space group symbol	:	$P3_1$
AFLOW prototype command	:	`aflow --proto=AB3C_hP45_144_3a_9a_3a --params=`$a, c/a, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}, x_{15}, y_{15}, z_{15}$

Other compounds with this structure

- $CsNO_3$ (II) and $TlNO_3$ (III)

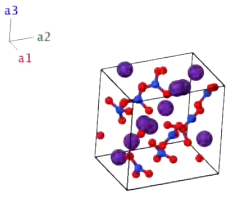
- $RbNO_3$ takes on four distinct phases with increasing temperature (Shamsuzzoha, 1988). Phase IV, presented here, is the ground state, stable up to 437 K. Phase III, stable in the range 437-492 K, is in the cesium chloride ($B2$) structure, with rubidium atoms on the cesium sites and orientationally disordered NO_3 ions on the chlorine sites. Phase II, stable

up to 564 K, is possibly a body-centered cubic structure with up to eight formula units in the conventional unit cell. Phase I, stable up to the melting point, is thought to be a sodium chloride ($B1$) structure, with NO_3 again playing the role of chlorine.

- The Phase IV structure also exists in the enantiomorphic space group $P3_2$ #145. (Shamsuzzoha, 1982)
- We used the data taken at room temperature, 298 K.

Trigonal Hexagonal primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$\tfrac{1}{2}(x_1+y_1)\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}(-x_1+y_1)\,a\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	$(3a)$	N I
$\mathbf{B}_2$	=	$-y_1\,\mathbf{a}_1 + (x_1-y_1)\,\mathbf{a}_2 + \left(\tfrac{1}{3}+z_1\right)\mathbf{a}_3$	=	$\left(\tfrac{1}{2}x_1 - y_1\right)a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}x_1 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{3}+z_1\right)c\,\hat{\mathbf{z}}$	$(3a)$	N I
$\mathbf{B}_3$	=	$(-x_1+y_1)\,\mathbf{a}_1 - x_1\,\mathbf{a}_2 + \left(\tfrac{2}{3}+z_1\right)\mathbf{a}_3$	=	$\left(-x_1 + \tfrac{1}{2}y_1\right)a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}y_1 a\,\hat{\mathbf{y}} + \left(\tfrac{2}{3}+z_1\right)c\,\hat{\mathbf{z}}$	$(3a)$	N I
$\mathbf{B}_4$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\tfrac{1}{2}(x_2+y_2)\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}(-x_2+y_2)\,a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(3a)$	N II
$\mathbf{B}_5$	=	$-y_2\,\mathbf{a}_1 + (x_2-y_2)\,\mathbf{a}_2 + \left(\tfrac{1}{3}+z_2\right)\mathbf{a}_3$	=	$\left(\tfrac{1}{2}x_2 - y_2\right)a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}x_2 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{3}+z_2\right)c\,\hat{\mathbf{z}}$	$(3a)$	N II
$\mathbf{B}_6$	=	$(-x_2+y_2)\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 + \left(\tfrac{2}{3}+z_2\right)\mathbf{a}_3$	=	$\left(-x_2 + \tfrac{1}{2}y_2\right)a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}y_2 a\,\hat{\mathbf{y}} + \left(\tfrac{2}{3}+z_2\right)c\,\hat{\mathbf{z}}$	$(3a)$	N II
$\mathbf{B}_7$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\tfrac{1}{2}(x_3+y_3)\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}(-x_3+y_3)\,a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(3a)$	N III
$\mathbf{B}_8$	=	$-y_3\,\mathbf{a}_1 + (x_3-y_3)\,\mathbf{a}_2 + \left(\tfrac{1}{3}+z_3\right)\mathbf{a}_3$	=	$\left(\tfrac{1}{2}x_3 - y_3\right)a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{3}+z_3\right)c\,\hat{\mathbf{z}}$	$(3a)$	N III
$\mathbf{B}_9$	=	$(-x_3+y_3)\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\tfrac{2}{3}+z_3\right)\mathbf{a}_3$	=	$\left(-x_3 + \tfrac{1}{2}y_3\right)a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}y_3 a\,\hat{\mathbf{y}} + \left(\tfrac{2}{3}+z_3\right)c\,\hat{\mathbf{z}}$	$(3a)$	N III
$\mathbf{B}_{10}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\tfrac{1}{2}(x_4+y_4)\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}(-x_4+y_4)\,a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(3a)$	O I
$\mathbf{B}_{11}$	=	$-y_4\,\mathbf{a}_1 + (x_4-y_4)\,\mathbf{a}_2 + \left(\tfrac{1}{3}+z_4\right)\mathbf{a}_3$	=	$\left(\tfrac{1}{2}x_4 - y_4\right)a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{3}+z_4\right)c\,\hat{\mathbf{z}}$	$(3a)$	O I
$\mathbf{B}_{12}$	=	$(-x_4+y_4)\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \left(\tfrac{2}{3}+z_4\right)\mathbf{a}_3$	=	$\left(-x_4 + \tfrac{1}{2}y_4\right)a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}y_4 a\,\hat{\mathbf{y}} + \left(\tfrac{2}{3}+z_4\right)c\,\hat{\mathbf{z}}$	$(3a)$	O I
$\mathbf{B}_{13}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\tfrac{1}{2}(x_5+y_5)\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}(-x_5+y_5)\,a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(3a)$	O II

$$\mathbf{B}_{14} = -y_5\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + \left(\tfrac{1}{3} + z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}x_5 - y_5\right)a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{3} + z_5\right)c\,\hat{\mathbf{z}} \quad (3a) \quad \text{O II}$$

$$\mathbf{B}_{15} = (-x_5 + y_5)\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \left(\tfrac{2}{3} + z_5\right)\mathbf{a}_3 = \left(-x_5 + \tfrac{1}{2}y_5\right)a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}y_5 a\,\hat{\mathbf{y}} + \left(\tfrac{2}{3} + z_5\right)c\,\hat{\mathbf{z}} \quad (3a) \quad \text{O II}$$

$$\mathbf{B}_{16} = x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 = \tfrac{1}{2}(x_6 + y_6)\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}(-x_6 + y_6)\,a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}} \quad (3a) \quad \text{O III}$$

$$\mathbf{B}_{17} = -y_6\,\mathbf{a}_1 + (x_6 - y_6)\,\mathbf{a}_2 + \left(\tfrac{1}{3} + z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}x_6 - y_6\right)a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{3} + z_6\right)c\,\hat{\mathbf{z}} \quad (3a) \quad \text{O III}$$

$$\mathbf{B}_{18} = (-x_6 + y_6)\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 + \left(\tfrac{2}{3} + z_6\right)\mathbf{a}_3 = \left(-x_6 + \tfrac{1}{2}y_6\right)a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}y_6 a\,\hat{\mathbf{y}} + \left(\tfrac{2}{3} + z_6\right)c\,\hat{\mathbf{z}} \quad (3a) \quad \text{O III}$$

$$\mathbf{B}_{19} = x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 = \tfrac{1}{2}(x_7 + y_7)\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}(-x_7 + y_7)\,a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}} \quad (3a) \quad \text{O IV}$$

$$\mathbf{B}_{20} = -y_7\,\mathbf{a}_1 + (x_7 - y_7)\,\mathbf{a}_2 + \left(\tfrac{1}{3} + z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}x_7 - y_7\right)a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}x_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{3} + z_7\right)c\,\hat{\mathbf{z}} \quad (3a) \quad \text{O IV}$$

$$\mathbf{B}_{21} = (-x_7 + y_7)\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 + \left(\tfrac{2}{3} + z_7\right)\mathbf{a}_3 = \left(-x_7 + \tfrac{1}{2}y_7\right)a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}y_7 a\,\hat{\mathbf{y}} + \left(\tfrac{2}{3} + z_7\right)c\,\hat{\mathbf{z}} \quad (3a) \quad \text{O IV}$$

$$\mathbf{B}_{22} = x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 = \tfrac{1}{2}(x_8 + y_8)\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}(-x_8 + y_8)\,a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}} \quad (3a) \quad \text{O V}$$

$$\mathbf{B}_{23} = -y_8\,\mathbf{a}_1 + (x_8 - y_8)\,\mathbf{a}_2 + \left(\tfrac{1}{3} + z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}x_8 - y_8\right)a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}x_8 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{3} + z_8\right)c\,\hat{\mathbf{z}} \quad (3a) \quad \text{O V}$$

$$\mathbf{B}_{24} = (-x_8 + y_8)\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 + \left(\tfrac{2}{3} + z_8\right)\mathbf{a}_3 = \left(-x_8 + \tfrac{1}{2}y_8\right)a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}y_8 a\,\hat{\mathbf{y}} + \left(\tfrac{2}{3} + z_8\right)c\,\hat{\mathbf{z}} \quad (3a) \quad \text{O V}$$

$$\mathbf{B}_{25} = x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 = \tfrac{1}{2}(x_9 + y_9)\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}(-x_9 + y_9)\,a\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}} \quad (3a) \quad \text{O VI}$$

$$\mathbf{B}_{26} = -y_9\,\mathbf{a}_1 + (x_9 - y_9)\,\mathbf{a}_2 + \left(\tfrac{1}{3} + z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}x_9 - y_9\right)a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}x_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{3} + z_9\right)c\,\hat{\mathbf{z}} \quad (3a) \quad \text{O VI}$$

$$\mathbf{B}_{27} = (-x_9 + y_9)\,\mathbf{a}_1 - x_9\,\mathbf{a}_2 + \left(\tfrac{2}{3} + z_9\right)\mathbf{a}_3 = \left(-x_9 + \tfrac{1}{2}y_9\right)a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}y_9 a\,\hat{\mathbf{y}} + \left(\tfrac{2}{3} + z_9\right)c\,\hat{\mathbf{z}} \quad (3a) \quad \text{O VI}$$

$$\mathbf{B}_{28} = x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 = \tfrac{1}{2}(x_{10} + y_{10})\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}(-x_{10} + y_{10})\,a\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}} \quad (3a) \quad \text{O VII}$$

$$\mathbf{B}_{29} = -y_{10}\,\mathbf{a}_1 + (x_{10} - y_{10})\,\mathbf{a}_2 + \left(\tfrac{1}{3} + z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}x_{10} - y_{10}\right)a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}x_{10} a\,\hat{\mathbf{y}} + \left(\tfrac{1}{3} + z_{10}\right)c\,\hat{\mathbf{z}} \quad (3a) \quad \text{O VII}$$

$$\mathbf{B}_{30} = (-x_{10} + y_{10})\,\mathbf{a}_1 - x_{10}\,\mathbf{a}_2 + \left(\tfrac{2}{3} + z_{10}\right)\mathbf{a}_3 = \left(-x_{10} + \tfrac{1}{2}y_{10}\right)a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}y_{10} a\,\hat{\mathbf{y}} + \left(\tfrac{2}{3} + z_{10}\right)c\,\hat{\mathbf{z}} \quad (3a) \quad \text{O VII}$$

$$\mathbf{B}_{31} = x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 = \tfrac{1}{2}(x_{11} + y_{11})\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}(-x_{11} + y_{11})\,a\,\hat{\mathbf{y}} + z_{11} c\,\hat{\mathbf{z}} \quad (3a) \quad \text{O VIII}$$

$$\mathbf{B}_{32} = -y_{11}\,\mathbf{a}_1 + (x_{11} - y_{11})\,\mathbf{a}_2 + \left(\tfrac{1}{3} + z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}x_{11} - y_{11}\right)a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}x_{11} a\,\hat{\mathbf{y}} + \left(\tfrac{1}{3} + z_{11}\right)c\,\hat{\mathbf{z}} \quad (3a) \quad \text{O VIII}$$

$$\mathbf{B}_{33} = (-x_{11} + y_{11})\,\mathbf{a}_1 - x_{11}\,\mathbf{a}_2 + \left(\tfrac{2}{3} + z_{11}\right)\mathbf{a}_3 = \left(-x_{11} + \tfrac{1}{2}y_{11}\right)a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}y_{11} a\,\hat{\mathbf{y}} + \left(\tfrac{2}{3} + z_{11}\right)c\,\hat{\mathbf{z}} \quad (3a) \quad \text{O VIII}$$

$\mathbf{B}_{34}$	=	$x_{12}\,\mathbf{a}_1 + y_{12}\,\mathbf{a}_2 + z_{12}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_{12} + y_{12})\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_{12} + y_{12})\,a\,\hat{\mathbf{y}} + z_{12}c\,\hat{\mathbf{z}}$	(3*a*)	O IX
$\mathbf{B}_{35}$	=	$-y_{12}\,\mathbf{a}_1 + (x_{12} - y_{12})\,\mathbf{a}_2 + \left(\frac{1}{3} + z_{12}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}x_{12} - y_{12}\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_{12}a\,\hat{\mathbf{y}} + \left(\frac{1}{3} + z_{12}\right)c\,\hat{\mathbf{z}}$	(3*a*)	O IX
$\mathbf{B}_{36}$	=	$(-x_{12} + y_{12})\,\mathbf{a}_1 - x_{12}\,\mathbf{a}_2 + \left(\frac{2}{3} + z_{12}\right)\mathbf{a}_3$	=	$\left(-x_{12} + \frac{1}{2}y_{12}\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_{12}a\,\hat{\mathbf{y}} + \left(\frac{2}{3} + z_{12}\right)c\,\hat{\mathbf{z}}$	(3*a*)	O IX
$\mathbf{B}_{37}$	=	$x_{13}\,\mathbf{a}_1 + y_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_{13} + y_{13})\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_{13} + y_{13})\,a\,\hat{\mathbf{y}} + z_{13}c\,\hat{\mathbf{z}}$	(3*a*)	Rb I
$\mathbf{B}_{38}$	=	$-y_{13}\,\mathbf{a}_1 + (x_{13} - y_{13})\,\mathbf{a}_2 + \left(\frac{1}{3} + z_{13}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}x_{13} - y_{13}\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_{13}a\,\hat{\mathbf{y}} + \left(\frac{1}{3} + z_{13}\right)c\,\hat{\mathbf{z}}$	(3*a*)	Rb I
$\mathbf{B}_{39}$	=	$(-x_{13} + y_{13})\,\mathbf{a}_1 - x_{13}\,\mathbf{a}_2 + \left(\frac{2}{3} + z_{13}\right)\mathbf{a}_3$	=	$\left(-x_{13} + \frac{1}{2}y_{13}\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_{13}a\,\hat{\mathbf{y}} + \left(\frac{2}{3} + z_{13}\right)c\,\hat{\mathbf{z}}$	(3*a*)	Rb I
$\mathbf{B}_{40}$	=	$x_{14}\,\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_{14} + y_{14})\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_{14} + y_{14})\,a\,\hat{\mathbf{y}} + z_{14}c\,\hat{\mathbf{z}}$	(3*a*)	Rb II
$\mathbf{B}_{41}$	=	$-y_{14}\,\mathbf{a}_1 + (x_{14} - y_{14})\,\mathbf{a}_2 + \left(\frac{1}{3} + z_{14}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}x_{14} - y_{14}\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_{14}a\,\hat{\mathbf{y}} + \left(\frac{1}{3} + z_{14}\right)c\,\hat{\mathbf{z}}$	(3*a*)	Rb II
$\mathbf{B}_{42}$	=	$(-x_{14} + y_{14})\,\mathbf{a}_1 - x_{14}\,\mathbf{a}_2 + \left(\frac{2}{3} + z_{14}\right)\mathbf{a}_3$	=	$\left(-x_{14} + \frac{1}{2}y_{14}\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_{14}a\,\hat{\mathbf{y}} + \left(\frac{2}{3} + z_{14}\right)c\,\hat{\mathbf{z}}$	(3*a*)	Rb II
$\mathbf{B}_{43}$	=	$x_{15}\,\mathbf{a}_1 + y_{15}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_{15} + y_{15})\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_{15} + y_{15})\,a\,\hat{\mathbf{y}} + z_{15}c\,\hat{\mathbf{z}}$	(3*a*)	Rb III
$\mathbf{B}_{44}$	=	$-y_{15}\,\mathbf{a}_1 + (x_{15} - y_{15})\,\mathbf{a}_2 + \left(\frac{1}{3} + z_{15}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}x_{15} - y_{15}\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_{15}a\,\hat{\mathbf{y}} + \left(\frac{1}{3} + z_{15}\right)c\,\hat{\mathbf{z}}$	(3*a*)	Rb III
$\mathbf{B}_{45}$	=	$(-x_{15} + y_{15})\,\mathbf{a}_1 - x_{15}\,\mathbf{a}_2 + \left(\frac{2}{3} + z_{15}\right)\mathbf{a}_3$	=	$\left(-x_{15} + \frac{1}{2}y_{15}\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_{15}a\,\hat{\mathbf{y}} + \left(\frac{2}{3} + z_{15}\right)c\,\hat{\mathbf{z}}$	(3*a*)	Rb III

References:

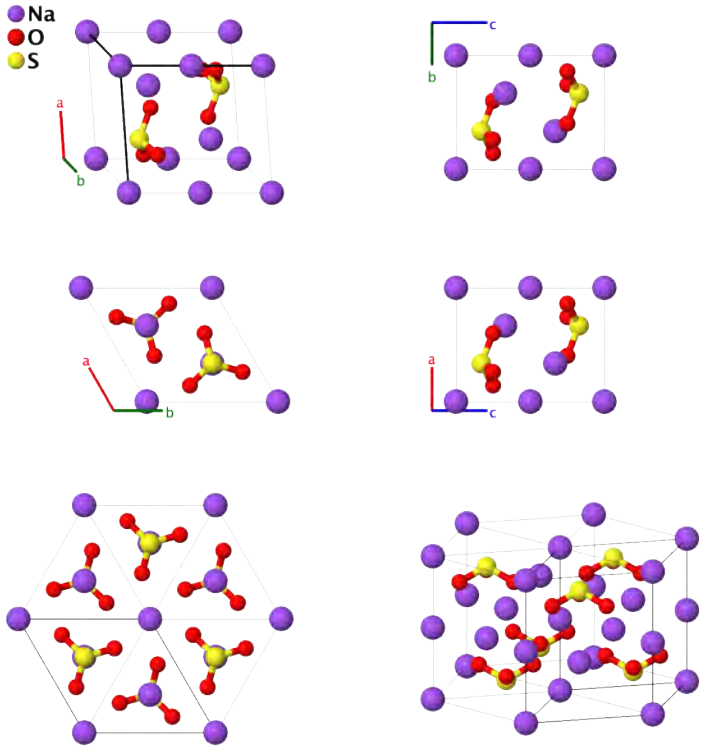
- M. Shamsuzzoha and B. W. Lucas, *Structure (neutron) of phase IV rubidium nitrate at 298 and 403 K*, Acta Crystallogr. Sect. B Struct. Sci. **38**, 2353–2357 (1982), doi:10.1107/S0567740882008772.
- M. Shamsuzzoha and B. W. Lucas, *Polymorphs of rubidium nitrate and their crystallographic relationships*, Can. J. Chem. **66**, 819–823 (1988), doi:10.1139/v88-142.

Geometry files:

- CIF: pp. 1719
- POSCAR: pp. 1719

Na_2SO_3 ($G3_2$) Structure: A2B3C_hP12_147_abd_g_d

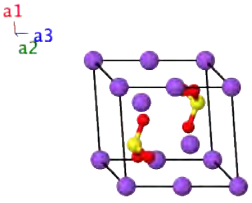
http://aflow.org/prototype-encyclopedia/A2B3C_hP12_147_abd_g_d



Prototype	:	Na_2O_3S
AFLOW prototype label	:	A2B3C_hP12_147_abd_g_d
***Strukturbericht* designation**	:	$G3_2$
Pearson symbol	:	hP12
Space group number	:	147
Space group symbol	:	$P\bar{3}$
AFLOW prototype command	:	`aflow --proto=A2B3C_hP12_147_abd_g_d --params=`$a, c/a, z_3, z_4, x_5, y_5, z_5$

Trigonal Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$ =	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($1a$)	Na I

$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	($1b$)	Na II
$\mathbf{B}_3$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	($2d$)	Na III
$\mathbf{B}_4$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	($2d$)	Na III
$\mathbf{B}_5$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	($2d$)	S
$\mathbf{B}_6$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	($2d$)	S
$\mathbf{B}_7$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_5 + y_5)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_5 + y_5)\,a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	($6g$)	O
$\mathbf{B}_8$	=	$-y_5\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_5 - y_5\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	($6g$)	O
$\mathbf{B}_9$	=	$(-x_5 + y_5)\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\left(-x_5 + \frac{1}{2}y_5\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	($6g$)	O
$\mathbf{B}_{10}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-\frac{1}{2}\,(x_5 + y_5)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(x_5 - y_5)\,a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	($6g$)	O
$\mathbf{B}_{11}$	=	$y_5\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_5 + y_5\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	($6g$)	O
$\mathbf{B}_{12}$	=	$(x_5 - y_5)\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\left(x_5 - \frac{1}{2}y_5\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	($6g$)	O

References:

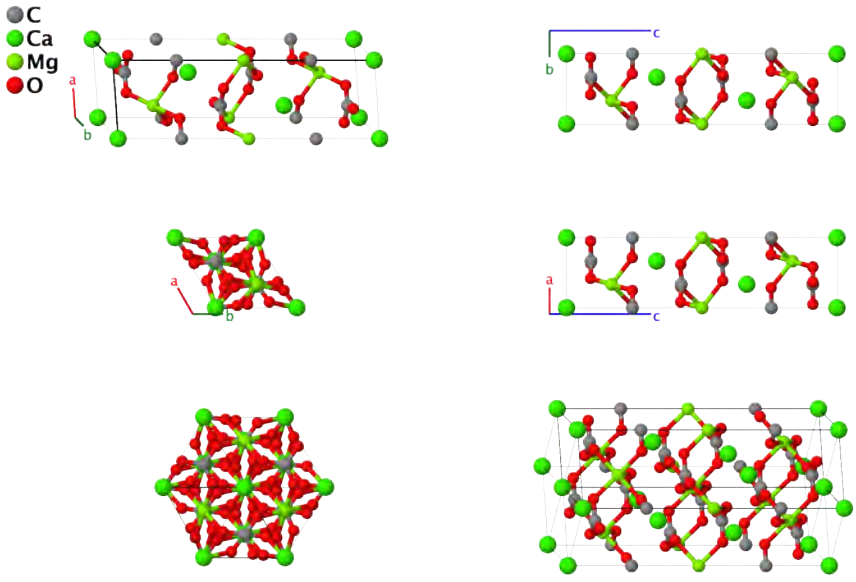
- W. H. Zachariasen and H. E. Buckley, *The Crystal Lattice of Anhydrous Sodium Sulphite, Na_2SO_3*, Phys. Rev. **37**, 1295–1305 (1931), doi:10.1103/PhysRev.37.1295.

Geometry files:

- CIF: pp. 1720
- POSCAR: pp. 1720

Dolomite [$MgCa(CO_3)_2$, $G1_1$] Structure: A2BCD6_hR10_148_c_a_b_f

http://aflow.org/prototype-encyclopedia/A2BCD6_hR10_148_c_a_b_f

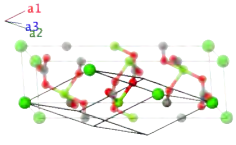


Prototype	:	C_2CaMgO_6
AFLOW prototype label	:	A2BCD6_hR10_148_c_a_b_f
***Strukturbericht* designation**	:	$G1_1$
Pearson symbol	:	hR10
Space group number	:	148
Space group symbol	:	$R\bar{3}$
AFLOW prototype command	:	`aflow --proto=A2BCD6_hR10_148_c_a_b_f [--hex] --params=`$a, c/a, x_3, x_4, y_4, z_4$

- Data was taken at 24 °C.

Rhombohedral primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = -\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($1a$)	Ca
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	($1b$)	Mg
$\mathbf{B}_3$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3 c\,\hat{\mathbf{z}}$	($2c$)	C

$\mathbf{B}_4$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$-x_3 c\,\hat{\mathbf{z}}$	$(2c)$	C
$\mathbf{B}_5$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\frac{1}{2}(x_4 - z_4)\,a\,\hat{\mathbf{x}} + \left(-\frac{1}{2\sqrt{3}}x_4 + \frac{1}{\sqrt{3}}y_4 - \frac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} + \frac{1}{3}(x_4 + y_4 + z_4)\,c\,\hat{\mathbf{z}}$	$(6f)$	O
$\mathbf{B}_6$	=	$z_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + y_4\,\mathbf{a}_3$	=	$\frac{1}{2}(-y_4 + z_4)\,a\,\hat{\mathbf{x}} + \left(\frac{1}{\sqrt{3}}x_4 - \frac{1}{2\sqrt{3}}y_4 - \frac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} + \frac{1}{3}(x_4 + y_4 + z_4)\,c\,\hat{\mathbf{z}}$	$(6f)$	O
$\mathbf{B}_7$	=	$y_4\,\mathbf{a}_1 + z_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$\frac{1}{2}(-x_4 + y_4)\,a\,\hat{\mathbf{x}} + \left(-\frac{1}{2\sqrt{3}}x_4 - \frac{1}{2\sqrt{3}}y_4 + \frac{1}{\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} + \frac{1}{3}(x_4 + y_4 + z_4)\,c\,\hat{\mathbf{z}}$	$(6f)$	O
$\mathbf{B}_8$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\frac{1}{2}(-x_4 + z_4)\,a\,\hat{\mathbf{x}} + \left(\frac{1}{2\sqrt{3}}x_4 - \frac{1}{\sqrt{3}}y_4 + \frac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} - \frac{1}{3}(x_4 + y_4 + z_4)\,c\,\hat{\mathbf{z}}$	$(6f)$	O
$\mathbf{B}_9$	=	$-z_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - y_4\,\mathbf{a}_3$	=	$\frac{1}{2}(y_4 - z_4)\,a\,\hat{\mathbf{x}} + \left(-\frac{1}{\sqrt{3}}x_4 + \frac{1}{2\sqrt{3}}y_4 + \frac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} - \frac{1}{3}(x_4 + y_4 + z_4)\,c\,\hat{\mathbf{z}}$	$(6f)$	O
$\mathbf{B}_{10}$	=	$-y_4\,\mathbf{a}_1 - z_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$\frac{1}{2}(x_4 - y_4)\,a\,\hat{\mathbf{x}} + \left(\frac{1}{2\sqrt{3}}x_4 + \frac{1}{2\sqrt{3}}y_4 - \frac{1}{\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} - \frac{1}{3}(x_4 + y_4 + z_4)\,c\,\hat{\mathbf{z}}$	$(6f)$	O

References:

- R. J. Reeder and S. A. Markgraf, *High-temperature crystal chemistry of dolomite*, Am. Mineral. **71**, 795–804 (1986).

Found in:

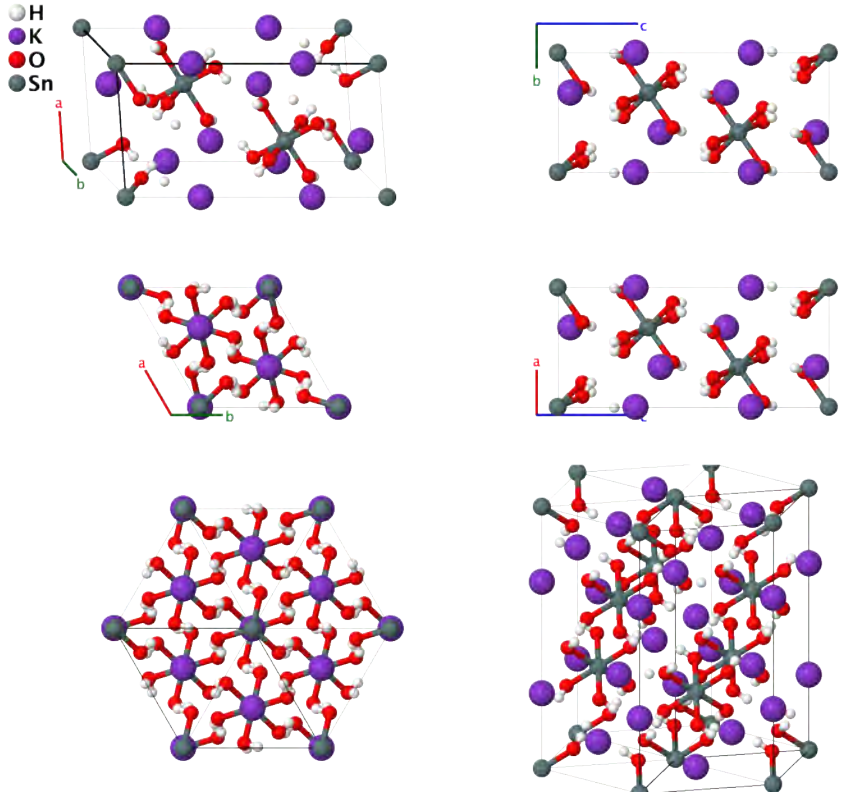
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1720
- POSCAR: pp. 1721

$K_2Sn(OH)_6$ ($H6_2$) Structure: A6B2C6D_hR15_148_f_c_f_a

http://aflow.org/prototype-encyclopedia/A6B2C6D_hR15_148_f_c_f_a



Prototype	:	$H_6K_2O_6Sn$
AFLOW prototype label	:	A6B2C6D_hR15_148_f_c_f_a
***Strukturbericht* designation**	:	$H6_2$
Pearson symbol	:	hR15
Space group number	:	148
Space group symbol	:	$R\bar{3}$
AFLOW prototype command	:	`aflow --proto=A6B2C6D_hR15_148_f_c_f_a [--hex] --params=`$a, c/a, x_2, x_3, y_3, z_3, x_4, y_4, z_4$

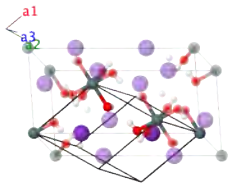
Other compounds with this structure

- $K_2Pt(OH)_6$ and $Na_2Sn(OH)_6$

- (Wyckoff, 1928) found the positions of the potassium and tin atoms in this structure. He could not locate the oxygen atoms, and so was unable to give an unequivocal determination of the space group, beyond noting that it was rhombohedral. We use the modern structure of (Jacobs, 2000), which agrees with Wyckoff in the positioning of the potassium and tin atoms, and finds both the hydrogen and oxygen locations.
- (Ewald, 1931) gave this the *Strukturbericht* designation $H6_2$. (Hermann, 1937) began relabeling the $H6$ structures as $I1$, and (Gottfried, 1937) relabeled them as $J1$, but neither they nor successor volumes revisited this structure, so it was never formally relabeled $J1_2$. We will therefore leave it as $H6_2$.

Rhombohedral primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2\sqrt{3}}\,a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,c\,\hat{\mathbf{z}}\\ \mathbf{a}_2 &= \tfrac{1}{\sqrt{3}}\,a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,c\,\hat{\mathbf{z}}\\ \mathbf{a}_3 &= -\tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2\sqrt{3}}\,a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(1a)$	Sn
$\mathbf{B}_2$	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	$x_2 c\,\hat{\mathbf{z}}$	$(2c)$	K
$\mathbf{B}_3$	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 - x_2\,\mathbf{a}_3$	$-x_2 c\,\hat{\mathbf{z}}$	$(2c)$	K
$\mathbf{B}_4$	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$\frac{1}{2}(x_3 - z_3)\,a\,\hat{\mathbf{x}} + \left(-\frac{1}{2\sqrt{3}}x_3 + \frac{1}{\sqrt{3}}y_3 - \frac{1}{2\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}} + \frac{1}{3}(x_3 + y_3 + z_3)\,c\,\hat{\mathbf{z}}$	$(6f)$	H
$\mathbf{B}_5$	$z_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + y_3\,\mathbf{a}_3$	$\frac{1}{2}(-y_3 + z_3)\,a\,\hat{\mathbf{x}} + \left(\frac{1}{\sqrt{3}}x_3 - \frac{1}{2\sqrt{3}}y_3 - \frac{1}{2\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}} + \frac{1}{3}(x_3 + y_3 + z_3)\,c\,\hat{\mathbf{z}}$	$(6f)$	H
$\mathbf{B}_6$	$y_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	$\frac{1}{2}(-x_3 + y_3)\,a\,\hat{\mathbf{x}} + \left(-\frac{1}{2\sqrt{3}}x_3 - \frac{1}{2\sqrt{3}}y_3 + \frac{1}{\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}} + \frac{1}{3}(x_3 + y_3 + z_3)\,c\,\hat{\mathbf{z}}$	$(6f)$	H
$\mathbf{B}_7$	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	$\frac{1}{2}(-x_3 + z_3)\,a\,\hat{\mathbf{x}} + \left(\frac{1}{2\sqrt{3}}x_3 - \frac{1}{\sqrt{3}}y_3 + \frac{1}{2\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}} - \frac{1}{3}(x_3 + y_3 + z_3)\,c\,\hat{\mathbf{z}}$	$(6f)$	H
$\mathbf{B}_8$	$-z_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - y_3\,\mathbf{a}_3$	$\frac{1}{2}(y_3 - z_3)\,a\,\hat{\mathbf{x}} + \left(-\frac{1}{\sqrt{3}}x_3 + \frac{1}{2\sqrt{3}}y_3 + \frac{1}{2\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}} - \frac{1}{3}(x_3 + y_3 + z_3)\,c\,\hat{\mathbf{z}}$	$(6f)$	H
$\mathbf{B}_9$	$-y_3\,\mathbf{a}_1 - z_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	$\frac{1}{2}(x_3 - y_3)\,a\,\hat{\mathbf{x}} + \left(\frac{1}{2\sqrt{3}}x_3 + \frac{1}{2\sqrt{3}}y_3 - \frac{1}{\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}} - \frac{1}{3}(x_3 + y_3 + z_3)\,c\,\hat{\mathbf{z}}$	$(6f)$	H
$\mathbf{B}_{10}$	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$\frac{1}{2}(x_4 - z_4)\,a\,\hat{\mathbf{x}} + \left(-\frac{1}{2\sqrt{3}}x_4 + \frac{1}{\sqrt{3}}y_4 - \frac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} + \frac{1}{3}(x_4 + y_4 + z_4)\,c\,\hat{\mathbf{z}}$	$(6f)$	O
$\mathbf{B}_{11}$	$z_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + y_4\,\mathbf{a}_3$	$\frac{1}{2}(-y_4 + z_4)\,a\,\hat{\mathbf{x}} + \left(\frac{1}{\sqrt{3}}x_4 - \frac{1}{2\sqrt{3}}y_4 - \frac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} + \frac{1}{3}(x_4 + y_4 + z_4)\,c\,\hat{\mathbf{z}}$	$(6f)$	O
$\mathbf{B}_{12}$	$y_4\,\mathbf{a}_1 + z_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	$\frac{1}{2}(-x_4 + y_4)\,a\,\hat{\mathbf{x}} + \left(-\frac{1}{2\sqrt{3}}x_4 - \frac{1}{2\sqrt{3}}y_4 + \frac{1}{\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} + \frac{1}{3}(x_4 + y_4 + z_4)\,c\,\hat{\mathbf{z}}$	$(6f)$	O
$\mathbf{B}_{13}$	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	$\frac{1}{2}(-x_4 + z_4)\,a\,\hat{\mathbf{x}} + \left(\frac{1}{2\sqrt{3}}x_4 - \frac{1}{\sqrt{3}}y_4 + \frac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} - \frac{1}{3}(x_4 + y_4 + z_4)\,c\,\hat{\mathbf{z}}$	$(6f)$	O
$\mathbf{B}_{14}$	$-z_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - y_4\,\mathbf{a}_3$	$\frac{1}{2}(y_4 - z_4)\,a\,\hat{\mathbf{x}} + \left(-\frac{1}{\sqrt{3}}x_4 + \frac{1}{2\sqrt{3}}y_4 + \frac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} - \frac{1}{3}(x_4 + y_4 + z_4)\,c\,\hat{\mathbf{z}}$	$(6f)$	O
$\mathbf{B}_{15}$	$-y_4\,\mathbf{a}_1 - z_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	$\frac{1}{2}(x_4 - y_4)\,a\,\hat{\mathbf{x}} + \left(\frac{1}{2\sqrt{3}}x_4 + \frac{1}{2\sqrt{3}}y_4 - \frac{1}{\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} - \frac{1}{3}(x_4 + y_4 + z_4)\,c\,\hat{\mathbf{z}}$	$(6f)$	O

References:

- H. Jacobs and R. Stahl, *Neubestimmung der Kristallstrukturen der Hexahydroxometallate $Na_2Sn(OH)_6$, $K_2Sn(OH)_6$ und*

$K_2Pb(OH)_6$, Z. Anorg. Allg. Chem. **626**, 1863–1866 (2000), doi:10.1002/1521-3749(200009)626:9%3C1863::AID-ZAAC1863%3E3.0.CO;2-M.
- R. W. G. Wyckoff, *The Crystal Structure of Potassium Hydroxystannate, $K_2Sn(OH)_6$*, Am. J. Sci. **s5-15**, 297–302 (1928), doi:10.2475/ajs.s5-15.88.297.
- P. P. Ewald and C. Hermann, eds., *Strukturbericht 1913-1928* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1931).
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).
- C. Gottfried and F. Schossberger, eds., *Strukturbericht Band III 1933-1935* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Found in:

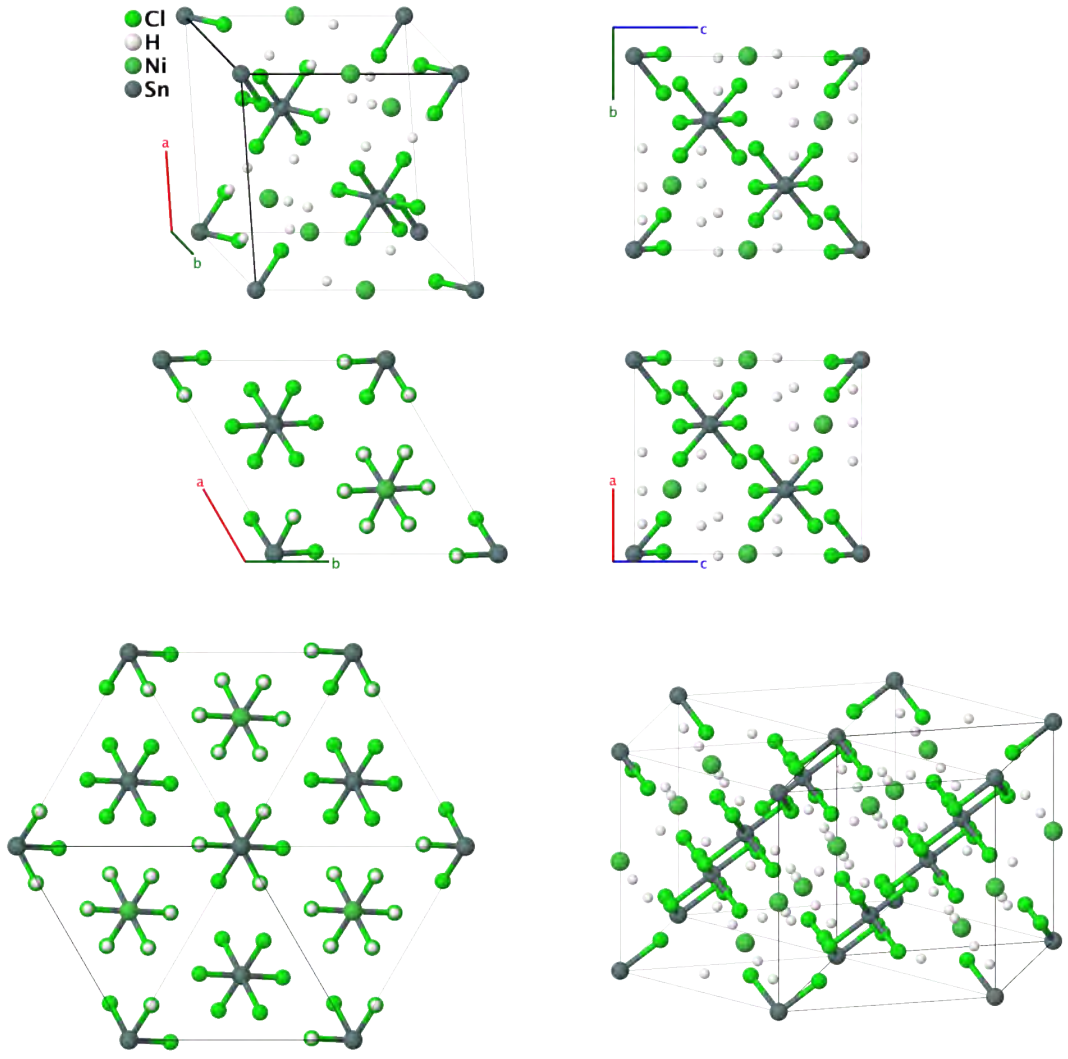
- P. Villars (Chief Editor), *$K_2Sn(OH)_6$ ($K_2Sn[OH]_6$) Crystal Structure*, http://materials.springer.com/isp/crystallographic/docs/sd_1703654 (2016). PAULING FILE in: Inorganic Solid Phases, SpringerMaterials (online database), Springer, Heidelberg (ed.) Springer Materials.

Geometry files:

- CIF: pp. 1721
- POSCAR: pp. 1721

$Ni(H_2O)_6SnCl_6$ ($I6_1$) Structure: A6B6CD_hR14_148_f_f_b_a

http://aflow.org/prototype-encyclopedia/A6B6CD_hR14_148_f_f_b_a



Prototype	:	$Cl_6(H_2O)_6NiSn$
AFLOW prototype label	:	A6B6CD_hR14_148_f_f_b_a
***Strukturbericht* designation**	:	$I6_1$
Pearson symbol	:	hR14
Space group number	:	148
Space group symbol	:	$R\bar{3}$
AFLOW prototype command	:	`aflow --proto=A6B6CD_hR14_148_f_f_b_a [--hex] --params=`$a, c/a, x_3, y_3, z_3, x_4, y_4, z_4$

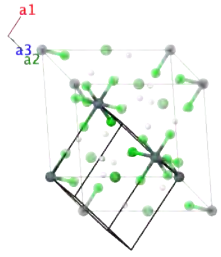
Other compounds with this structure

- $Ca(H_2O)_6SnF_6$, $Co(H_2O)_6PtF_6$, $Co(H_2O)_6SiF_6$, $Co(NH_3)_6Co(CN)_6$, $Fe(H_2O)_6SiF_6$, $Mg(H_2O)_6SiF_6$, $Mg(H_2O)_6SnF_6$, $Mg(H_2O)_6TiF_6$, $Mn(H_2O)_6SiF_6$, $Ni(H_2O)_6SiF_6$, $Zn(H_2O)_6SiF_6$, $Zn(H_2O)_6SnF_6$, $Zn(H_2O)_6TiF_6$, and $Zn(H_2O)_6ZrF_6$

- $Ni(H_2O)_6SnCl_6$ is the prototype for a large class of molecular crystals with the form MG_6LR_6, where MG_6 is a cation and LR_6 is an anion. (Hermann, 1937) gave this the *Strukturbericht* designation $I6_1$. (Gottfried, 1937) changed the I designations to J, so this should have become $J6_1$, but it was never referenced in any form in later volumes of *Strukturbericht*. Since $I6_1$ is the only designation for this structure in the literature, we use it rather than $J6_1$.
- The positions of the hydrogen atoms in the water molecules were not determined, so we only provide the positions of the oxygen atoms (labeled as H_2O).

Rhombohedral primitive vectors:

$$\begin{aligned}
\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2\sqrt{3}}\,a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,c\,\hat{\mathbf{z}} \\
\mathbf{a}_2 &= \tfrac{1}{\sqrt{3}}\,a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,c\,\hat{\mathbf{z}} \\
\mathbf{a}_3 &= -\tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2\sqrt{3}}\,a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,c\,\hat{\mathbf{z}}
\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(1a)$	Sn
$\mathbf{B}_2$	=	$\tfrac{1}{2}\,\mathbf{a}_1 + \tfrac{1}{2}\,\mathbf{a}_2 + \tfrac{1}{2}\,\mathbf{a}_3$	=	$\tfrac{1}{2}c\,\hat{\mathbf{z}}$	$(1b)$	Ni
$\mathbf{B}_3$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\tfrac{1}{2}\,(x_3 - z_3)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_3 + \tfrac{1}{\sqrt{3}}y_3 - \tfrac{1}{2\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,(x_3 + y_3 + z_3)\,c\,\hat{\mathbf{z}}$	$(6f)$	Cl
$\mathbf{B}_4$	=	$z_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + y_3\,\mathbf{a}_3$	=	$\tfrac{1}{2}\,(-y_3 + z_3)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{\sqrt{3}}x_3 - \tfrac{1}{2\sqrt{3}}y_3 - \tfrac{1}{2\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,(x_3 + y_3 + z_3)\,c\,\hat{\mathbf{z}}$	$(6f)$	Cl
$\mathbf{B}_5$	=	$y_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$\tfrac{1}{2}\,(-x_3 + y_3)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_3 - \tfrac{1}{2\sqrt{3}}y_3 + \tfrac{1}{\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,(x_3 + y_3 + z_3)\,c\,\hat{\mathbf{z}}$	$(6f)$	Cl
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\tfrac{1}{2}\,(-x_3 + z_3)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2\sqrt{3}}x_3 - \tfrac{1}{\sqrt{3}}y_3 + \tfrac{1}{2\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}} - \tfrac{1}{3}\,(x_3 + y_3 + z_3)\,c\,\hat{\mathbf{z}}$	$(6f)$	Cl
$\mathbf{B}_7$	=	$-z_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - y_3\,\mathbf{a}_3$	=	$\tfrac{1}{2}\,(y_3 - z_3)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{\sqrt{3}}x_3 + \tfrac{1}{2\sqrt{3}}y_3 + \tfrac{1}{2\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}} - \tfrac{1}{3}\,(x_3 + y_3 + z_3)\,c\,\hat{\mathbf{z}}$	$(6f)$	Cl
$\mathbf{B}_8$	=	$-y_3\,\mathbf{a}_1 - z_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$\tfrac{1}{2}\,(x_3 - y_3)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2\sqrt{3}}x_3 + \tfrac{1}{2\sqrt{3}}y_3 - \tfrac{1}{\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}} - \tfrac{1}{3}\,(x_3 + y_3 + z_3)\,c\,\hat{\mathbf{z}}$	$(6f)$	Cl
$\mathbf{B}_9$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\tfrac{1}{2}\,(x_4 - z_4)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_4 + \tfrac{1}{\sqrt{3}}y_4 - \tfrac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,(x_4 + y_4 + z_4)\,c\,\hat{\mathbf{z}}$	$(6f)$	H_2O
$\mathbf{B}_{10}$	=	$z_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + y_4\,\mathbf{a}_3$	=	$\tfrac{1}{2}\,(-y_4 + z_4)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{\sqrt{3}}x_4 - \tfrac{1}{2\sqrt{3}}y_4 - \tfrac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,(x_4 + y_4 + z_4)\,c\,\hat{\mathbf{z}}$	$(6f)$	H_2O
$\mathbf{B}_{11}$	=	$y_4\,\mathbf{a}_1 + z_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$\tfrac{1}{2}\,(-x_4 + y_4)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_4 - \tfrac{1}{2\sqrt{3}}y_4 + \tfrac{1}{\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,(x_4 + y_4 + z_4)\,c\,\hat{\mathbf{z}}$	$(6f)$	H_2O
$\mathbf{B}_{12}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\tfrac{1}{2}\,(-x_4 + z_4)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2\sqrt{3}}x_4 - \tfrac{1}{\sqrt{3}}y_4 + \tfrac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} - \tfrac{1}{3}\,(x_4 + y_4 + z_4)\,c\,\hat{\mathbf{z}}$	$(6f)$	H_2O

$$\mathbf{B}_{13} = -z_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - y_4\,\mathbf{a}_3 = \tfrac{1}{2}\,(y_4 - z_4)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{\sqrt{3}}x_4 + \tfrac{1}{2\sqrt{3}}y_4 + \tfrac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} - \tfrac{1}{3}\,(x_4 + y_4 + z_4)\,c\,\hat{\mathbf{z}} \quad (6f) \quad H_2O$$

$$\mathbf{B}_{14} = -y_4\,\mathbf{a}_1 - z_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3 = \tfrac{1}{2}\,(x_4 - y_4)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2\sqrt{3}}x_4 + \tfrac{1}{2\sqrt{3}}y_4 - \tfrac{1}{\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} - \tfrac{1}{3}\,(x_4 + y_4 + z_4)\,c\,\hat{\mathbf{z}} \quad (6f) \quad H_2O$$

References:

- L. Pauling, *On the crystal structure of nickel chlorostannate hexahydrate*, Zeitschrift für Kristallographie - Crystalline Materials **72**, 482–492 (1930), doi:10.1524/zkri.1930.72.1.482.
- C. Gottfried and F. Schossberger, eds., *Strukturbericht Band III 1933-1935* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Found in:

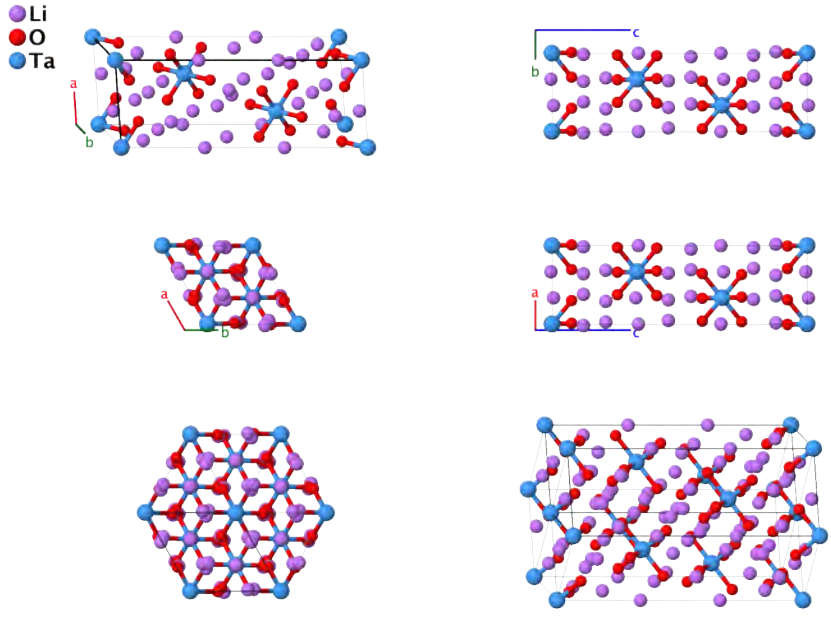
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1721
- POSCAR: pp. 1722

Li_7TaO_6 Structure: A8B6C_hR15_148_cf_f_a

http://aflow.org/prototype-encyclopedia/A8B6C_hR15_148_cf_f_a

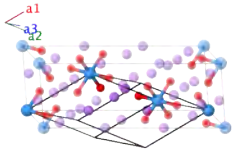


Prototype	:	Li_7O_6Ta
AFLOW prototype label	:	A8B6C_hR15_148_cf_f_a
***Strukturbericht* designation**	:	None
Pearson symbol	:	hR15
Space group number	:	148
Space group symbol	:	$R\bar{3}$
AFLOW prototype command	:	`aflow --proto=A8B6C_hR15_148_cf_f_a [--hex] --params=`$a, c/a, x_2, x_3, y_3, z_3, x_4, y_4, z_4$

- The Li-I ($2c$) site is half-occupied.

Rhombohedral primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = -\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$= 0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	$= 0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($1a$)	Ta
$\mathbf{B}_2$	$= x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	$= x_2 c\,\hat{\mathbf{z}}$	($2c$)	Li I
$\mathbf{B}_3$	$= -x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 - x_2\,\mathbf{a}_3$	$= -x_2 c\,\hat{\mathbf{z}}$	($2c$)	Li I

$$\mathbf{B}_4 = x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3 = \tfrac{1}{2}(x_3 - z_3)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_3 + \tfrac{1}{\sqrt{3}}y_3 - \tfrac{1}{2\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}(x_3 + y_3 + z_3)\,c\,\hat{\mathbf{z}} \quad (6f) \quad \text{Li II}$$

$$\mathbf{B}_5 = z_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + y_3\,\mathbf{a}_3 = \tfrac{1}{2}(-y_3 + z_3)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{\sqrt{3}}x_3 - \tfrac{1}{2\sqrt{3}}y_3 - \tfrac{1}{2\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}(x_3 + y_3 + z_3)\,c\,\hat{\mathbf{z}} \quad (6f) \quad \text{Li II}$$

$$\mathbf{B}_6 = y_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3 = \tfrac{1}{2}(-x_3 + y_3)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_3 - \tfrac{1}{2\sqrt{3}}y_3 + \tfrac{1}{\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}(x_3 + y_3 + z_3)\,c\,\hat{\mathbf{z}} \quad (6f) \quad \text{Li II}$$

$$\mathbf{B}_7 = -x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3 = \tfrac{1}{2}(-x_3 + z_3)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2\sqrt{3}}x_3 - \tfrac{1}{\sqrt{3}}y_3 + \tfrac{1}{2\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}} - \tfrac{1}{3}(x_3 + y_3 + z_3)\,c\,\hat{\mathbf{z}} \quad (6f) \quad \text{Li II}$$

$$\mathbf{B}_8 = -z_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - y_3\,\mathbf{a}_3 = \tfrac{1}{2}(y_3 - z_3)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{\sqrt{3}}x_3 + \tfrac{1}{2\sqrt{3}}y_3 + \tfrac{1}{2\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}} - \tfrac{1}{3}(x_3 + y_3 + z_3)\,c\,\hat{\mathbf{z}} \quad (6f) \quad \text{Li II}$$

$$\mathbf{B}_9 = -y_3\,\mathbf{a}_1 - z_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3 = \tfrac{1}{2}(x_3 - y_3)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2\sqrt{3}}x_3 + \tfrac{1}{2\sqrt{3}}y_3 - \tfrac{1}{\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}} - \tfrac{1}{3}(x_3 + y_3 + z_3)\,c\,\hat{\mathbf{z}} \quad (6f) \quad \text{Li II}$$

$$\mathbf{B}_{10} = x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3 = \tfrac{1}{2}(x_4 - z_4)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_4 + \tfrac{1}{\sqrt{3}}y_4 - \tfrac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}(x_4 + y_4 + z_4)\,c\,\hat{\mathbf{z}} \quad (6f) \quad \text{O}$$

$$\mathbf{B}_{11} = z_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + y_4\,\mathbf{a}_3 = \tfrac{1}{2}(-y_4 + z_4)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{\sqrt{3}}x_4 - \tfrac{1}{2\sqrt{3}}y_4 - \tfrac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}(x_4 + y_4 + z_4)\,c\,\hat{\mathbf{z}} \quad (6f) \quad \text{O}$$

$$\mathbf{B}_{12} = y_4\,\mathbf{a}_1 + z_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3 = \tfrac{1}{2}(-x_4 + y_4)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_4 - \tfrac{1}{2\sqrt{3}}y_4 + \tfrac{1}{\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}(x_4 + y_4 + z_4)\,c\,\hat{\mathbf{z}} \quad (6f) \quad \text{O}$$

$$\mathbf{B}_{13} = -x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3 = \tfrac{1}{2}(-x_4 + z_4)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2\sqrt{3}}x_4 - \tfrac{1}{\sqrt{3}}y_4 + \tfrac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} - \tfrac{1}{3}(x_4 + y_4 + z_4)\,c\,\hat{\mathbf{z}} \quad (6f) \quad \text{O}$$

$$\mathbf{B}_{14} = -z_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - y_4\,\mathbf{a}_3 = \tfrac{1}{2}(y_4 - z_4)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{\sqrt{3}}x_4 + \tfrac{1}{2\sqrt{3}}y_4 + \tfrac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} - \tfrac{1}{3}(x_4 + y_4 + z_4)\,c\,\hat{\mathbf{z}} \quad (6f) \quad \text{O}$$

$$\mathbf{B}_{15} = -y_4\,\mathbf{a}_1 - z_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3 = \tfrac{1}{2}(x_4 - y_4)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2\sqrt{3}}x_4 + \tfrac{1}{2\sqrt{3}}y_4 - \tfrac{1}{\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} - \tfrac{1}{3}(x_4 + y_4 + z_4)\,c\,\hat{\mathbf{z}} \quad (6f) \quad \text{O}$$

References:

- G. Wehrum and R. Hoppe, *Zur Kenntnis 'Kationen-reicher' Tantalate Über $Li_7[TaO_6]$*, Z. Anorg. Allg. Chem. **620**, 659–664 (1994), doi:10.1002/zaac.19946200414.

Found in:

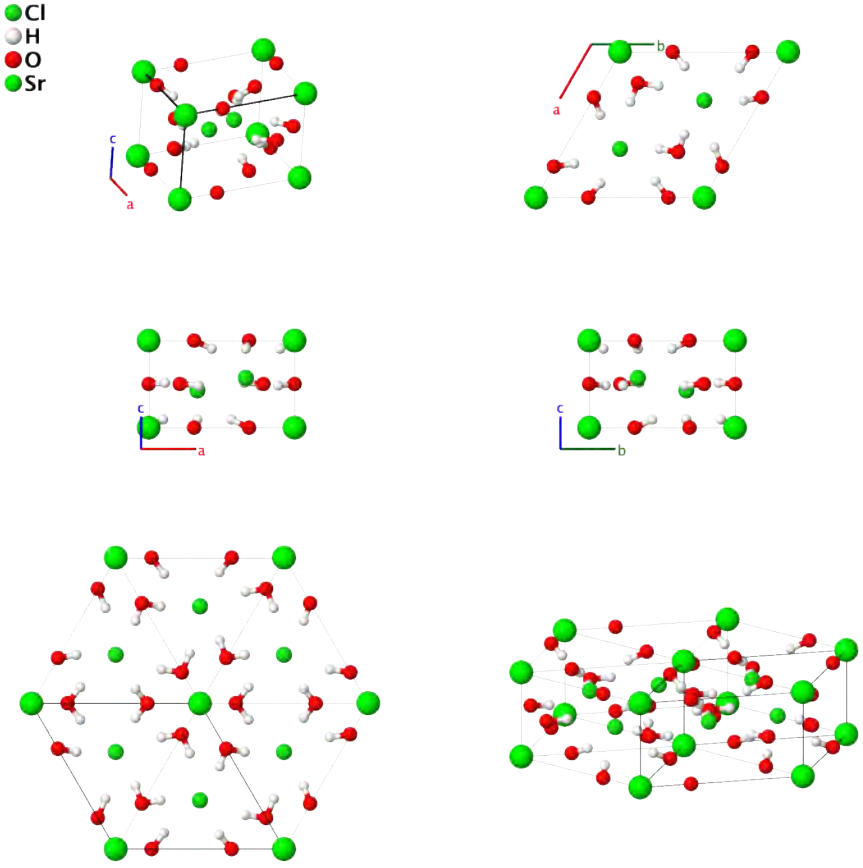
- L. Kahle, X. Cheng, T. Binninger, S. D. Lacey, A. Marcolongo, F. Zipoli, E. Gilardi, C. Villevieille, M. El Kazzi, N. Marzari, and D. Pergolesi, *The solid-state Li-ion conductor Li_7TaO_6: A combined computational and experimental study*, http://arxiv.org/abs/1910.11079 (2019). ArXiv:1910.11079 [cond-mat.mtrl-sci].

Geometry files:

- CIF: pp. 1722
- POSCAR: pp. 1722

$SrCl_2 \cdot (H_2O)_6$ Structure: A2B12C6D_hP21_150_d_2g_ef_a

http://aflow.org/prototype-encyclopedia/A2B12C6D_hP21_150_d_2g_ef_a



Prototype	:	$Cl_2H_{12}O_6Sr$
AFLOW prototype label	:	A2B12C6D_hP21_150_d_2g_ef_a
***Strukturbericht* designation**	:	None
Pearson symbol	:	hP21
Space group number	:	150
Space group symbol	:	$P321$
AFLOW prototype command	:	`aflow --proto=A2B12C6D_hP21_150_d_2g_ef_a --params=`$a, c/a, z_2, x_3, x_4, x_5, y_5, z_5, x_6, y_6, z_6$

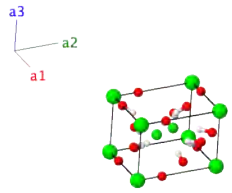
Other compounds with this structure

- $CaCl_2 \cdot (H_2O)_6$ and $CaBr_2 \cdot (H_2O)_6$

- This is a redetermination of the $I1_3$ structure. In addition to locating the hydrogen atoms in the water molecule, (Agron, 1986) show that the space group is $P321$ #150, not $P\bar{3}$ #147.

Trigonal Hexagonal primitive vectors:

$$\mathbf{a}_1 = \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(1a)$	Sr
$\mathbf{B}_2$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(2d)$	Cl
$\mathbf{B}_3$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	$(2d)$	Cl
$\mathbf{B}_4$	=	$x_3\,\mathbf{a}_1$	=	$\frac{1}{2}x_3 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}}$	$(3e)$	O I
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_2$	=	$\frac{1}{2}x_3 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}}$	$(3e)$	O I
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2$	=	$-x_3 a\,\hat{\mathbf{x}}$	$(3e)$	O I
$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}x_4 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(3f)$	O II
$\mathbf{B}_8$	=	$x_4\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}x_4 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(3f)$	O II
$\mathbf{B}_9$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(3f)$	O II
$\mathbf{B}_{10}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\frac{1}{2}(x_5 + y_5)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(-x_5 + y_5)\,a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(6g)$	H I
$\mathbf{B}_{11}$	=	$-y_5\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_5 - y_5\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(6g)$	H I
$\mathbf{B}_{12}$	=	$(-x_5 + y_5)\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\left(-x_5 + \frac{1}{2}y_5\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(6g)$	H I
$\mathbf{B}_{13}$	=	$y_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\frac{1}{2}(x_5 + y_5)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(x_5 - y_5)\,a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	$(6g)$	H I
$\mathbf{B}_{14}$	=	$(x_5 - y_5)\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_5 - y_5\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	$(6g)$	H I
$\mathbf{B}_{15}$	=	$-x_5\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\left(-x_5 + \frac{1}{2}y_5\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	$(6g)$	H I
$\mathbf{B}_{16}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\frac{1}{2}(x_6 + y_6)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(-x_6 + y_6)\,a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(6g)$	H II
$\mathbf{B}_{17}$	=	$-y_6\,\mathbf{a}_1 + (x_6 - y_6)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_6 - y_6\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(6g)$	H II
$\mathbf{B}_{18}$	=	$(-x_6 + y_6)\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\left(-x_6 + \frac{1}{2}y_6\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(6g)$	H II
$\mathbf{B}_{19}$	=	$y_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$\frac{1}{2}(x_6 + y_6)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(x_6 - y_6)\,a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	$(6g)$	H II
$\mathbf{B}_{20}$	=	$(x_6 - y_6)\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_6 - y_6\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	$(6g)$	H II
$\mathbf{B}_{21}$	=	$-x_6\,\mathbf{a}_1 + (-x_6 + y_6)\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$\left(-x_6 + \frac{1}{2}y_6\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_6 a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	$(6g)$	H II

References:

- P. A. Agron and W. R. Busing, *Calcium and strontium dichloride hexahydrates by neutron diffraction*, Acta Crystallogr. C **42**, 141–143 (1986), doi:10.1107/S0108270186097007.

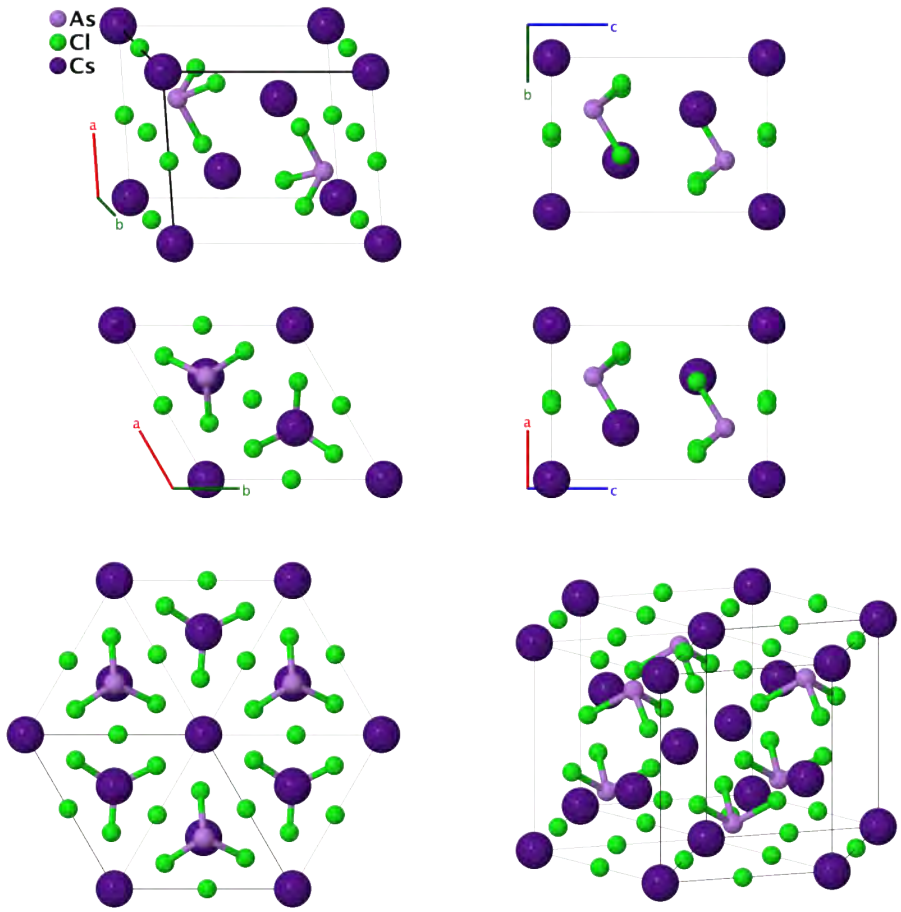
Found in:

- P. Villars (Chief Editor), *PAULING FILE* (2016). In: Inorganic Solid Phases, SpringerMaterials (online database), Springer, Heidelberg SpringerMaterials.

Geometry files:
- CIF: pp. 1722
- POSCAR: pp. 1723

$Cs_3As_2Cl_9$ ($K7_3$) Structure: A2B9C3_hP14_150_d_eg_ad

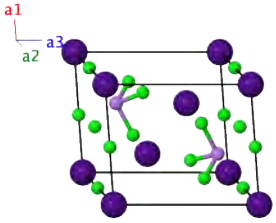
http://aflow.org/prototype-encyclopedia/A2B9C3_hP14_150_d_eg_ad



Prototype	:	$As_2Cl_9Cs_3$
AFLOW prototype label	:	A2B9C3_hP14_150_d_eg_ad
***Strukturbericht* designation**	:	$K7_3$
Pearson symbol	:	hP14
Space group number	:	150
Space group symbol	:	$P321$
AFLOW prototype command	:	`aflow --proto=A2B9C3_hP14_150_d_eg_ad --params=`$a, c/a, z_2, z_3, x_4, x_5, y_5, z_5$

Trigonal Hexagonal primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(1*a*)	Cs I
$\mathbf{B}_2$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(2*d*)	As
$\mathbf{B}_3$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(2*d*)	As
$\mathbf{B}_4$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(2*d*)	Cs II
$\mathbf{B}_5$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(2*d*)	Cs II
$\mathbf{B}_6$	=	$x_4\,\mathbf{a}_1$	=	$\frac{1}{2}x_4 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}}$	(3*e*)	Cl I
$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_2$	=	$\frac{1}{2}x_4 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}}$	(3*e*)	Cl I
$\mathbf{B}_8$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2$	=	$-x_4 a\,\hat{\mathbf{x}}$	(3*e*)	Cl I
$\mathbf{B}_9$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\frac{1}{2}(x_5 + y_5)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(-x_5 + y_5)\,a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(6*g*)	Cl II
$\mathbf{B}_{10}$	=	$-y_5\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_5 - y_5\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(6*g*)	Cl II
$\mathbf{B}_{11}$	=	$(-x_5 + y_5)\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\left(-x_5 + \frac{1}{2}y_5\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(6*g*)	Cl II
$\mathbf{B}_{12}$	=	$y_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\frac{1}{2}(x_5 + y_5)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(x_5 - y_5)\,a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(6*g*)	Cl II
$\mathbf{B}_{13}$	=	$(x_5 - y_5)\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_5 - y_5\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(6*g*)	Cl II
$\mathbf{B}_{14}$	=	$-x_5\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\left(-x_5 + \frac{1}{2}y_5\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(6*g*)	Cl II

References:

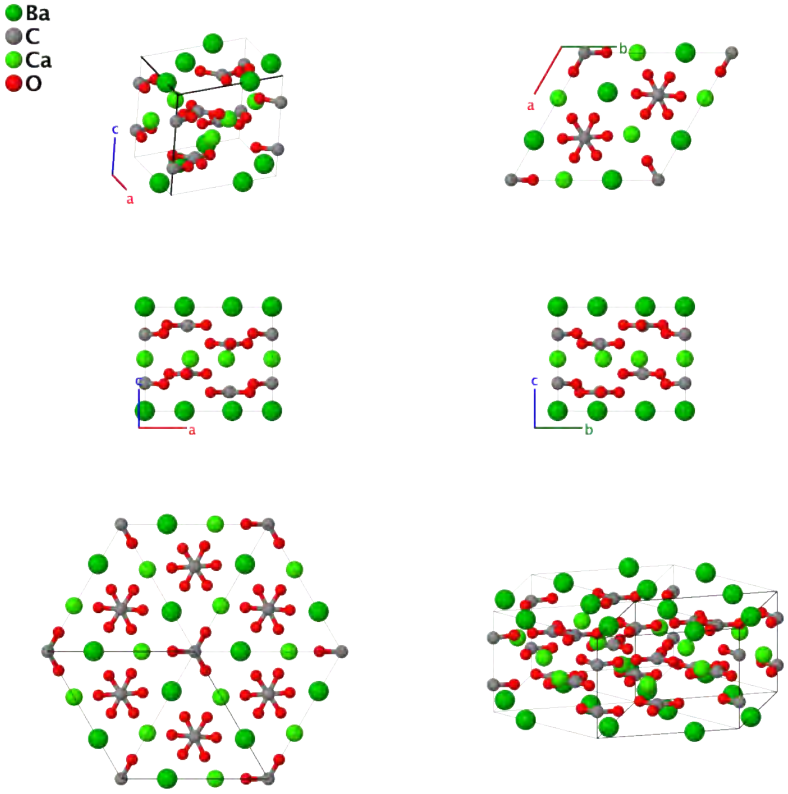
- J. L. Hoard and L. Goldstein, *The Structure of Caesium Enneachlordiarsenite, $Cs_3As_2Cl_9$*, J. Chem. Phys. **3**, 117–122 (1935), doi:10.1063/1.1749606.

Geometry files:

- CIF: pp. 1723
- POSCAR: pp. 1723

Paralstonite ($BaCa(CO_3)_2$) Structure: AB2CD6_hP30_150_e_c2d_f_3g

http://aflow.org/prototype-encyclopedia/AB2CD6_hP30_150_e_c2d_f_3g

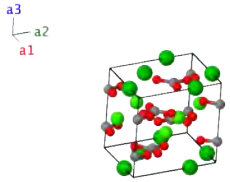


Prototype	:	BaC_2CaO_6
AFLOW prototype label	:	AB2CD6_hP30_150_e_c2d_f_3g
***Strukturbericht* designation**	:	None
Pearson symbol	:	hP30
Space group number	:	150
Space group symbol	:	$P321$
AFLOW prototype command	:	`aflow --proto=AB2CD6_hP30_150_e_c2d_f_3g --params=`$a, c/a, z_1, z_2, z_3, x_4, x_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8$

- $BaCa(CO_3)_2$ comes in a variety of crystal structures (Spahr, 2019):
  - monoclinic barytocalcite, space group $P2_1/m$ #11,
  - trigonal paralstonite, space group $P321$ #150 (the current structure),
  - triclinic alstonite, space group $P1$ #1 or $P\bar{1}$ #2 (Sartori, 1975), and
  - a new monoclinic structure, space group $C2$ #5, synthesized by (Spahr, 2019), and lacking the centrosymmetric character of barytocalcite.
- We were unable to obtain a copy of (Effenberger, 1980), so we use the data provided by (Downs, 2003).

Trigonal Hexagonal primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$z_1\,\mathbf{a}_3$	=	$z_1 c\,\hat{\mathbf{z}}$	$(2c)$	C I
$\mathbf{B}_2$	=	$-z_1\,\mathbf{a}_3$	=	$-z_1 c\,\hat{\mathbf{z}}$	$(2c)$	C I
$\mathbf{B}_3$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(2d)$	C II
$\mathbf{B}_4$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	$(2d)$	C II
$\mathbf{B}_5$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(2d)$	C III
$\mathbf{B}_6$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(2d)$	C III
$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_1$	=	$\frac{1}{2}x_4 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}}$	$(3e)$	Ba
$\mathbf{B}_8$	=	$x_4\,\mathbf{a}_2$	=	$\frac{1}{2}x_4 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}}$	$(3e)$	Ba
$\mathbf{B}_9$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2$	=	$-x_4 a\,\hat{\mathbf{x}}$	$(3e)$	Ba
$\mathbf{B}_{10}$	=	$x_5\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}x_5 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(3f)$	Ca
$\mathbf{B}_{11}$	=	$x_5\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}x_5 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(3f)$	Ca
$\mathbf{B}_{12}$	=	$-x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(3f)$	Ca
$\mathbf{B}_{13}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\frac{1}{2}(x_6 + y_6)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(-x_6 + y_6)\,a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(6g)$	O I
$\mathbf{B}_{14}$	=	$-y_6\,\mathbf{a}_1 + (x_6 - y_6)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_6 - y_6\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(6g)$	O I
$\mathbf{B}_{15}$	=	$(-x_6 + y_6)\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\left(-x_6 + \frac{1}{2}y_6\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(6g)$	O I
$\mathbf{B}_{16}$	=	$y_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$\frac{1}{2}(x_6 + y_6)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(x_6 - y_6)\,a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	$(6g)$	O I
$\mathbf{B}_{17}$	=	$(x_6 - y_6)\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_6 - y_6\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	$(6g)$	O I
$\mathbf{B}_{18}$	=	$-x_6\,\mathbf{a}_1 + (-x_6 + y_6)\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$\left(-x_6 + \frac{1}{2}y_6\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_6 a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	$(6g)$	O I
$\mathbf{B}_{19}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$\frac{1}{2}(x_7 + y_7)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(-x_7 + y_7)\,a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	$(6g)$	O II
$\mathbf{B}_{20}$	=	$-y_7\,\mathbf{a}_1 + (x_7 - y_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_7 - y_7\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_7 a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	$(6g)$	O II
$\mathbf{B}_{21}$	=	$(-x_7 + y_7)\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$\left(-x_7 + \frac{1}{2}y_7\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_7 a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	$(6g)$	O II
$\mathbf{B}_{22}$	=	$y_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$\frac{1}{2}(x_7 + y_7)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(x_7 - y_7)\,a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	$(6g)$	O II
$\mathbf{B}_{23}$	=	$(x_7 - y_7)\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_7 - y_7\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_7 a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	$(6g)$	O II
$\mathbf{B}_{24}$	=	$-x_7\,\mathbf{a}_1 + (-x_7 + y_7)\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$\left(-x_7 + \frac{1}{2}y_7\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_7 a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	$(6g)$	O II
$\mathbf{B}_{25}$	=	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$\frac{1}{2}(x_8 + y_8)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(-x_8 + y_8)\,a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	$(6g)$	O III
$\mathbf{B}_{26}$	=	$-y_8\,\mathbf{a}_1 + (x_8 - y_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_8 - y_8\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_8 a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	$(6g)$	O III
$\mathbf{B}_{27}$	=	$(-x_8 + y_8)\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$\left(-x_8 + \frac{1}{2}y_8\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_8 a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	$(6g)$	O III

$\mathbf{B}_{28}$	=	$y_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_8 + y_8)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(x_8 - y_8)\,a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(6*g*)	O III
$\mathbf{B}_{29}$	=	$(x_8 - y_8)\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_8 - y_8\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_8 a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(6*g*)	O III
$\mathbf{B}_{30}$	=	$-x_8\,\mathbf{a}_1 + (-x_8 + y_8)\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$\left(-x_8 + \frac{1}{2}y_8\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_8 a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(6*g*)	O III

References:

- H. Effenberger, *Die Kristallstruktur des Minerals Paralstonite, $BaCa(CO_3)_2$*, Neues Jahrb. Mineral. Monatsh. **1980**, 353–363 (1980).
- F. Sartori, *New data on alstonite*, Lithos **8**, 199–207 (1975), doi:10.1016/0024-4937(75)90036-5.
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Found in:

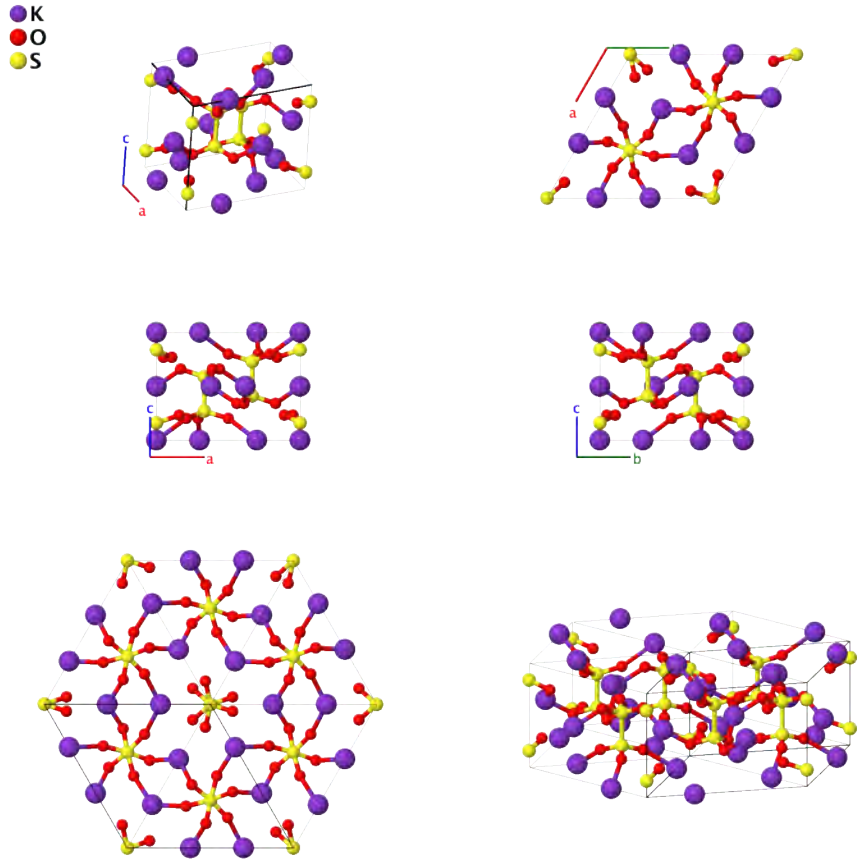
- D. Spahr, L. Bayarjargal, V. Vinograd, R. Luchitskaia, V. Milman, and B. Winkler, *A new $BaCa(CO_3)_2$ polymorph*, Acta Crystallogr. Sect. B Struct. Sci. **75**, 291–300 (2019), doi:10.1107/S2052520619003238.

Geometry files:

- CIF: pp. 1723
- POSCAR: pp. 1724

KSO_3 ($K1_1$) Structure: AB3C_hP30_150_ef_3g_c2d

http://aflow.org/prototype-encyclopedia/AB3C_hP30_150_ef_3g_c2d



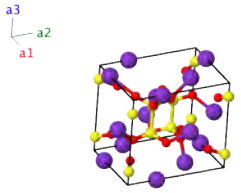
Prototype	:	KO_3S
AFLOW prototype label	:	AB3C_hP30_150_ef_3g_c2d
***Strukturbericht* designation**	:	$K1_1$
Pearson symbol	:	hP30
Space group number	:	150
Space group symbol	:	$P321$
AFLOW prototype command	:	`aflow --proto=AB3C_hP30_150_ef_3g_c2d --params=`$a, c/a, z_1, z_2, z_3, x_4, x_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8$

Other compounds with this structure

- $RbSO_3$

Trigonal Hexagonal primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$z_1\,\mathbf{a}_3$	=	$z_1 c\,\hat{\mathbf{z}}$	$(2c)$	S I
$\mathbf{B}_2$	=	$-z_1\,\mathbf{a}_3$	=	$-z_1 c\,\hat{\mathbf{z}}$	$(2c)$	S I
$\mathbf{B}_3$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(2d)$	S II
$\mathbf{B}_4$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	$(2d)$	S II
$\mathbf{B}_5$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(2d)$	S III
$\mathbf{B}_6$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(2d)$	S III
$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_1$	=	$\frac{1}{2}x_4 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}}$	$(3e)$	K I
$\mathbf{B}_8$	=	$x_4\,\mathbf{a}_2$	=	$\frac{1}{2}x_4 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}}$	$(3e)$	K I
$\mathbf{B}_9$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2$	=	$-x_4 a\,\hat{\mathbf{x}}$	$(3e)$	K I
$\mathbf{B}_{10}$	=	$x_5\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}x_5 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(3f)$	K II
$\mathbf{B}_{11}$	=	$x_5\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}x_5 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(3f)$	K II
$\mathbf{B}_{12}$	=	$-x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(3f)$	K II
$\mathbf{B}_{13}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\frac{1}{2}(x_6 + y_6)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(-x_6 + y_6)\,a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(6g)$	O I
$\mathbf{B}_{14}$	=	$-y_6\,\mathbf{a}_1 + (x_6 - y_6)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_6 - y_6\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(6g)$	O I
$\mathbf{B}_{15}$	=	$(-x_6 + y_6)\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\left(-x_6 + \frac{1}{2}y_6\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(6g)$	O I
$\mathbf{B}_{16}$	=	$y_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$\frac{1}{2}(x_6 + y_6)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(x_6 - y_6)\,a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	$(6g)$	O I
$\mathbf{B}_{17}$	=	$(x_6 - y_6)\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_6 - y_6\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	$(6g)$	O I
$\mathbf{B}_{18}$	=	$-x_6\,\mathbf{a}_1 + (-x_6 + y_6)\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$\left(-x_6 + \frac{1}{2}y_6\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_6 a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	$(6g)$	O I
$\mathbf{B}_{19}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$\frac{1}{2}(x_7 + y_7)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(-x_7 + y_7)\,a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	$(6g)$	O II
$\mathbf{B}_{20}$	=	$-y_7\,\mathbf{a}_1 + (x_7 - y_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_7 - y_7\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_7 a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	$(6g)$	O II
$\mathbf{B}_{21}$	=	$(-x_7 + y_7)\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$\left(-x_7 + \frac{1}{2}y_7\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_7 a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	$(6g)$	O II
$\mathbf{B}_{22}$	=	$y_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$\frac{1}{2}(x_7 + y_7)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(x_7 - y_7)\,a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	$(6g)$	O II
$\mathbf{B}_{23}$	=	$(x_7 - y_7)\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_7 - y_7\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_7 a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	$(6g)$	O II
$\mathbf{B}_{24}$	=	$-x_7\,\mathbf{a}_1 + (-x_7 + y_7)\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$\left(-x_7 + \frac{1}{2}y_7\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_7 a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	$(6g)$	O II
$\mathbf{B}_{25}$	=	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$\frac{1}{2}(x_8 + y_8)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(-x_8 + y_8)\,a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	$(6g)$	O III
$\mathbf{B}_{26}$	=	$-y_8\,\mathbf{a}_1 + (x_8 - y_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_8 - y_8\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_8 a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	$(6g)$	O III
$\mathbf{B}_{27}$	=	$(-x_8 + y_8)\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$\left(-x_8 + \frac{1}{2}y_8\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_8 a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	$(6g)$	O III

$\mathbf{B}_{28}$	=	$y_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_8 + y_8)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(x_8 - y_8)\,a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(6*g*)	O III
$\mathbf{B}_{29}$	=	$(x_8 - y_8)\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_8 - y_8\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_8 a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(6*g*)	O III
$\mathbf{B}_{30}$	=	$-x_8\,\mathbf{a}_1 + (-x_8 + y_8)\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$\left(-x_8 + \frac{1}{2}y_8\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_8 a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	(6*g*)	O III

References:

- M. L. Huggins and G. O. Frank, *The crystal structure of potassium dithionate, $K_2S_2O_6$*, Am. Mineral. **16**, 580–591 (1931).

Found in:

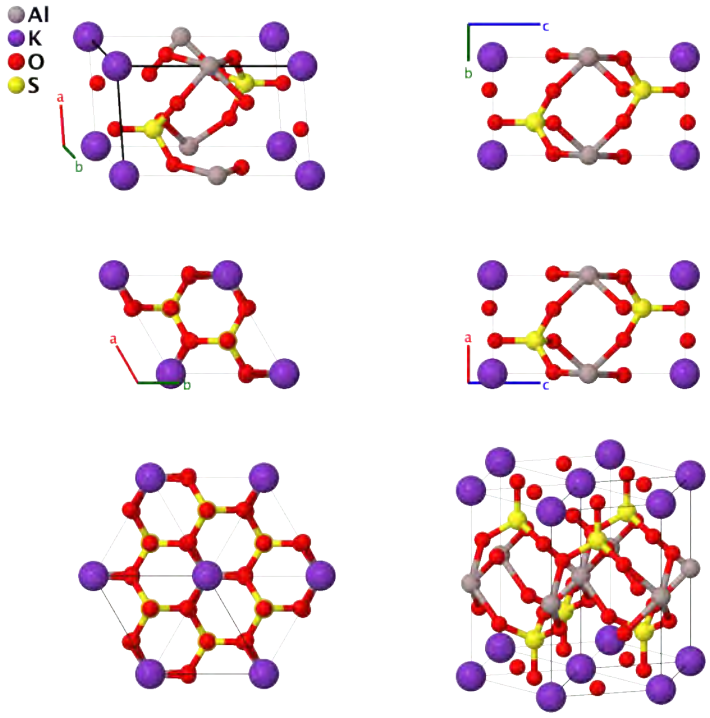
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1724
- POSCAR: pp. 1724

Steklite [$KAl(SO_4)_2$, $H3_2$] Structure: ABC8D2_hP12_150_b_a_dg_d

http://aflow.org/prototype-encyclopedia/ABC8D2_hP12_150_b_a_dg_d



Prototype	:	$AlKO_8S_2$
AFLOW prototype label	:	ABC8D2_hP12_150_b_a_dg_d
***Strukturbericht* designation**	:	$H3_2$
Pearson symbol	:	hP12
Space group number	:	150
Space group symbol	:	$P321$
AFLOW prototype command	:	`aflow --proto=ABC8D2_hP12_150_b_a_dg_d --params=`$a, c/a, z_3, z_4, x_5, y_5, z_5$

Other compounds with this structure

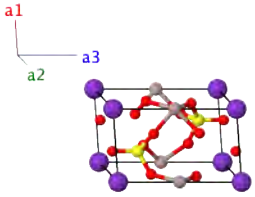
- $NH_4(Al,Fe)(SO_4)_2$ (godovikovite), $KCr(SO_4)_2$, and $RbCr(SO_4)_2$

- This has been a rather difficult structure to follow through the literature. (Villars, 2016) quotes the structure given by (Manoli, 1970), but gives the space group as $P\bar{3}$ #147. (Murashko, 2013) lists the space group as both $P312$ #149 and $P321$ #150, as well as listing obviously incorrect Wyckoff positions.
- (West, 2008) states that the simple structure is in $P\bar{3}$ but that it may be doubled along the c axis and be in space group $P321$.
- After correcting Murashko's results, we find that all of these interpretations yield essentially the same structure in a given layer, and only differ as the structure is reflected through the $z = 0$ plane. As it is not clear which structure is correct, we will use the original $H3_2$ structure given by (Hermann, 1937).

- Steklite is the name of the mineral form of this compound (Murashko, 2013). (Hermann, 1937) simply calls it *Wasserfreier Alaun* (anhydrous alum). For hydrated alum, $KAl(SO_4)_2 \cdot 12H_2O$, see the $H4_{13}$ structure.

Trigonal Hexagonal primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(1*a*)	K
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	(1*b*)	Al
$\mathbf{B}_3$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(2*d*)	O I
$\mathbf{B}_4$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(2*d*)	O I
$\mathbf{B}_5$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(2*d*)	S
$\mathbf{B}_6$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(2*d*)	S
$\mathbf{B}_7$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_5 + y_5)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_5 + y_5)\,a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(6*g*)	O II
$\mathbf{B}_8$	=	$-y_5\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_5 - y_5\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(6*g*)	O II
$\mathbf{B}_9$	=	$(-x_5 + y_5)\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\left(-x_5 + \frac{1}{2}y_5\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(6*g*)	O II
$\mathbf{B}_{10}$	=	$y_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_5 + y_5)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(x_5 - y_5)\,a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(6*g*)	O II
$\mathbf{B}_{11}$	=	$(x_5 - y_5)\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_5 - y_5\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(6*g*)	O II
$\mathbf{B}_{12}$	=	$-x_5\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\left(-x_5 + \frac{1}{2}y_5\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(6*g*)	O II

References:

- L. Vegard and A. Maurstad, *Die Kristallstruktur der wasserfreien Alaune R'R"$(SO_4)_2$*, Zeitschrift für Kristallographie - Crystalline Materials **69**, 519–532 (1929), doi:10.1524/zkri.1929.69.1.519.
- L. Vegard and A. Maurstad, *Die Kristallstruktur der wasserfreien Alaune R'R"$(SO_4)_2$*, Skrifter utgitt av det Norske Videnskaps-Akademi i Oslo pp. 1–24 (1928).
- D. V. West, Q. Huang, H. W. Zandbergen, T. M. McQueen, and R. J. Cava, *Structural disorder, octahedral coordination and two-dimensional ferromagnetism in anhydrous alums*, J. Solid State Chem. **181**, 2768–2775 (2008), doi:10.1016/j.jssc.2008.07.006.
- M. N. Murashko, I. V. Pekov, S. V. Krivovichev, A. P. Chernyatyeva, V. O. Yapaskurt, A. E. Zadov, and M. E. Zelensky, *Steklite, $KAl(SO_4)_2$: A finding at the Tolbachik Volcano, Kamchatka, Russia, validating its status as a mineral species and crystal structure*, Geol. Ore Deposits **55**, 594–600 (2013), doi:10.1134/S1075701513070088.
- P. Villars (Chief Editor), *$KAl(SO_4)_2$ ($KAl[SO_4]_2$ trig1) Crystal Structure*, http://materials.springer.com/isp/crystallographic/docs/sd_0381481 (2016). PAULING FILE in: Inorganic Solid Phases, SpringerMaterials (online database).
- J.-M. Manoli, P. Herpin, and G. Pannetier, *Structure cristalline du sulfate double d'aluminium et de potassium*, Bull. Soc. Chim. France pp. 98–101 (1970).

Found in:

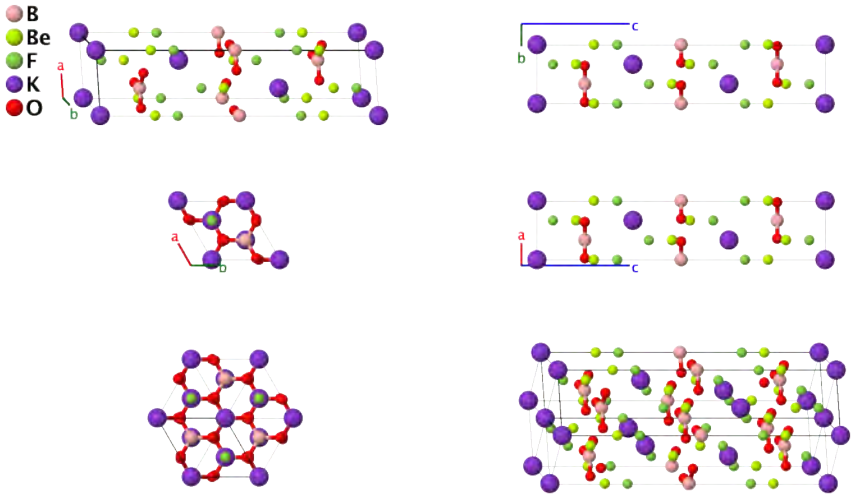
- P. P. Ewald and C. Hermann, eds., *Strukturbericht 1913-1928* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1931).

- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:
- CIF: pp. 1724
- POSCAR: pp. 1725

$KBe_2BO_3F_2$ Structure: AB2C2DE3_hR9_155_b_c_c_a_e

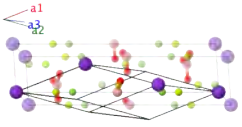
http://aflow.org/prototype-encyclopedia/AB2C2DE3_hR9_155_b_c_c_a_e



Prototype	:	$BBe_2F_2KO_3$
AFLOW prototype label	:	AB2C2DE3_hR9_155_b_c_c_a_e
***Strukturbericht* designation**	:	None
Pearson symbol	:	hR9
Space group number	:	155
Space group symbol	:	$R32$
AFLOW prototype command	:	`aflow --proto=AB2C2DE3_hR9_155_b_c_c_a_e [--hex] --params=`$a, c/a, x_3, x_4, y_5$

Rhombohedral primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = -\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(1a)$	K
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	$(1b)$	B
$\mathbf{B}_3$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3 c\,\hat{\mathbf{z}}$	$(2c)$	Be
$\mathbf{B}_4$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$-x_3 c\,\hat{\mathbf{z}}$	$(2c)$	Be
$\mathbf{B}_5$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4 c\,\hat{\mathbf{z}}$	$(2c)$	F
$\mathbf{B}_6$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$-x_4 c\,\hat{\mathbf{z}}$	$(2c)$	F
$\mathbf{B}_7$	=	$\frac{1}{2}\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 - y_5\,\mathbf{a}_3$	=	$\left(\frac{1}{4} + \frac{1}{2}y_5\right)a\,\hat{\mathbf{x}} + \left(-\frac{1}{4\sqrt{3}} + \frac{\sqrt{3}}{2}y_5\right)a\,\hat{\mathbf{y}} + \frac{1}{6}c\,\hat{\mathbf{z}}$	$(3e)$	O
$\mathbf{B}_8$	=	$-y_5\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + y_5\,\mathbf{a}_3$	=	$-y_5 a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{6}c\,\hat{\mathbf{z}}$	$(3e)$	O

$$\mathbf{B}_9 = y_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \tfrac{1}{2}\,\mathbf{a}_3 = \left(-\tfrac{1}{4} + \tfrac{1}{2}y_5\right) a\,\hat{\mathbf{x}} - \left(\tfrac{\sqrt{3}}{2}y_5 + \tfrac{1}{4\sqrt{3}}\right) a\,\hat{\mathbf{y}} + \tfrac{1}{6}c\,\hat{\mathbf{z}} \quad (3e) \quad \text{O}$$

References:

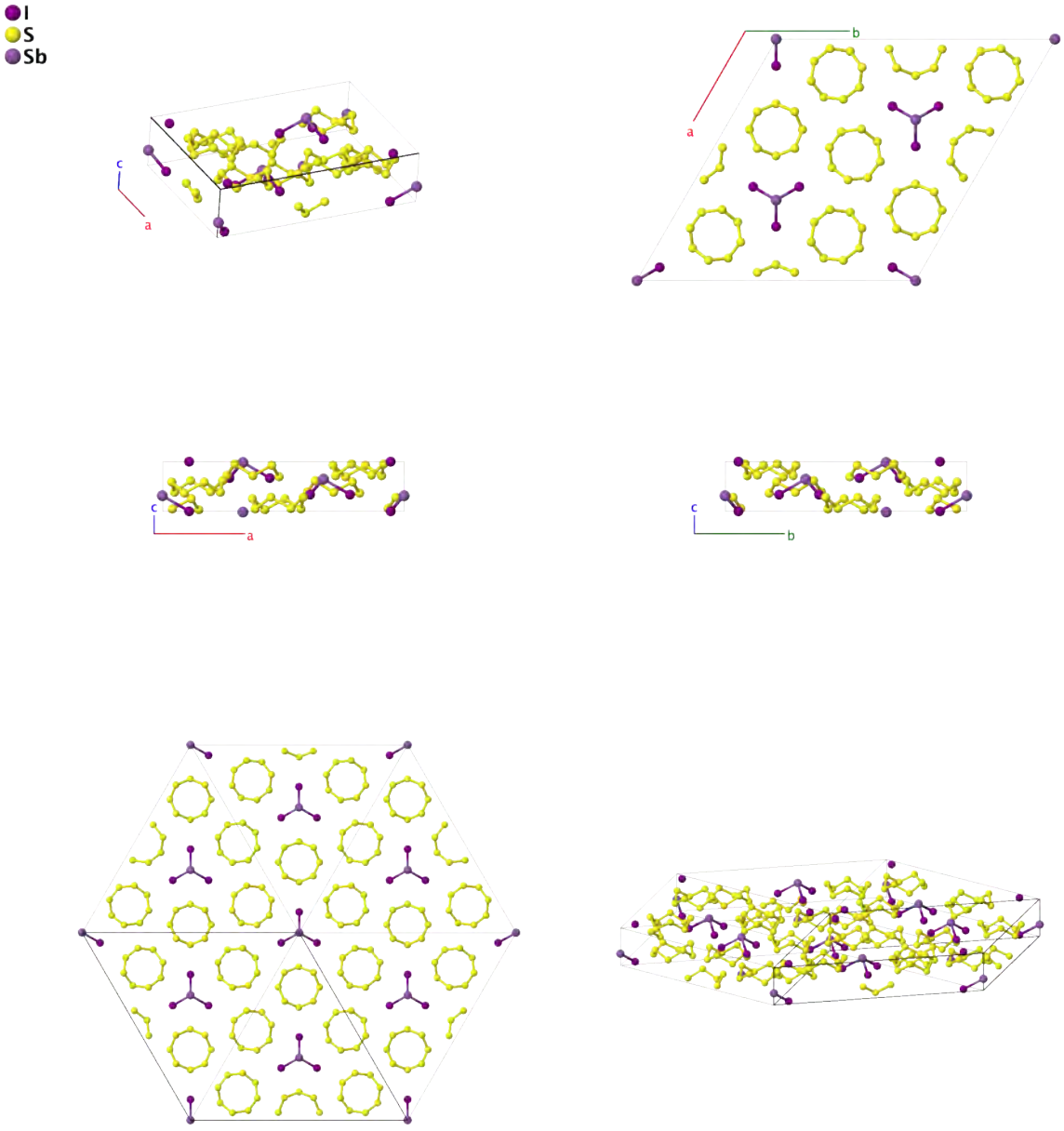
- L. Mei, X. Huang, Y. Wang, Q. Wu, and C. Chen, *Crystal Structure of* $KBe_2BO_3F_2$, Zeitschrift für Kristallographie - Crystalline Materials **210**, 93–95 (1995), doi:10.1524/zkri.1995.210.2.93.

Geometry files:

- CIF: pp. 1725
- POSCAR: pp. 1725

SbI_3S_{24} Structure: A3B24C_hR28_160_b_2b3c_a

http://aflow.org/prototype-encyclopedia/A3B24C_hR28_160_b_2b3c_a

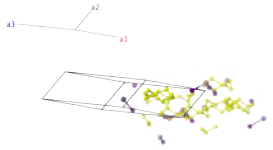


Prototype	:	$I_3S_{24}Sb$
AFLOW prototype label	:	A3B24C_hR28_160_b_2b3c_a
***Strukturbericht* designation**	:	None
Pearson symbol	:	hR28
Space group number	:	160
Space group symbol	:	$R3m$
AFLOW prototype command	:	`aflow --proto=A3B24C_hR28_160_b_2b3c_a [--hex] --params=`$a, c/a, x_1, x_2, z_2, x_3, z_3, x_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7$

- Since there are three S_8 molecules in this structure, (Bjorvatten, 1963) refer to it as SbI_3:3S_8.
- Space group *R*3*m* #160 does not fix the zero of the *z*-axis. Here it is set to coincide with the plane of the iodine atoms.

Rhombohedral primitive vectors:

$$\begin{aligned}
\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2\sqrt{3}}\,a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,c\,\hat{\mathbf{z}} \\
\mathbf{a}_2 &= \tfrac{1}{\sqrt{3}}\,a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,c\,\hat{\mathbf{z}} \\
\mathbf{a}_3 &= -\tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2\sqrt{3}}\,a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,c\,\hat{\mathbf{z}}
\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + x_1\,\mathbf{a}_2 + x_1\,\mathbf{a}_3$	=	$x_1 c\,\hat{\mathbf{z}}$	(1*a*)	Sb
$\mathbf{B}_2$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_2 - z_2)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}\,(x_2 - z_2)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_2 + \frac{1}{3}z_2\right)c\,\hat{\mathbf{z}}$	(3*b*)	I
$\mathbf{B}_3$	=	$z_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$\frac{1}{2}\,(-x_2 + z_2)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}\,(x_2 - z_2)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_2 + \frac{1}{3}z_2\right)c\,\hat{\mathbf{z}}$	(3*b*)	I
$\mathbf{B}_4$	=	$x_2\,\mathbf{a}_1 + z_2\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$\frac{1}{\sqrt{3}}\,(-x_2 + z_2)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_2 + \frac{1}{3}z_2\right)c\,\hat{\mathbf{z}}$	(3*b*)	I
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_3 - z_3)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}\,(x_3 - z_3)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_3 + \frac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	(3*b*)	S I
$\mathbf{B}_6$	=	$z_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$\frac{1}{2}\,(-x_3 + z_3)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}\,(x_3 - z_3)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_3 + \frac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	(3*b*)	S I
$\mathbf{B}_7$	=	$x_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$\frac{1}{\sqrt{3}}\,(-x_3 + z_3)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_3 + \frac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	(3*b*)	S I
$\mathbf{B}_8$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_4 - z_4)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}\,(x_4 - z_4)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_4 + \frac{1}{3}z_4\right)c\,\hat{\mathbf{z}}$	(3*b*)	S II
$\mathbf{B}_9$	=	$z_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$\frac{1}{2}\,(-x_4 + z_4)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}\,(x_4 - z_4)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_4 + \frac{1}{3}z_4\right)c\,\hat{\mathbf{z}}$	(3*b*)	S II
$\mathbf{B}_{10}$	=	$x_4\,\mathbf{a}_1 + z_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$\frac{1}{\sqrt{3}}\,(-x_4 + z_4)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_4 + \frac{1}{3}z_4\right)c\,\hat{\mathbf{z}}$	(3*b*)	S II
$\mathbf{B}_{11}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_5 - z_5)\,a\,\hat{\mathbf{x}} + \left(-\frac{1}{2\sqrt{3}}x_5 + \frac{1}{\sqrt{3}}y_5 - \frac{1}{2\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}} + \frac{1}{3}\,(x_5 + y_5 + z_5)\,c\,\hat{\mathbf{z}}$	(6*c*)	S III
$\mathbf{B}_{12}$	=	$z_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + y_5\,\mathbf{a}_3$	=	$\frac{1}{2}\,(-y_5 + z_5)\,a\,\hat{\mathbf{x}} + \left(\frac{1}{\sqrt{3}}x_5 - \frac{1}{2\sqrt{3}}y_5 - \frac{1}{2\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}} + \frac{1}{3}\,(x_5 + y_5 + z_5)\,c\,\hat{\mathbf{z}}$	(6*c*)	S III
$\mathbf{B}_{13}$	=	$y_5\,\mathbf{a}_1 + z_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$\frac{1}{2}\,(-x_5 + y_5)\,a\,\hat{\mathbf{x}} + \left(-\frac{1}{2\sqrt{3}}x_5 - \frac{1}{2\sqrt{3}}y_5 + \frac{1}{\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}} + \frac{1}{3}\,(x_5 + y_5 + z_5)\,c\,\hat{\mathbf{z}}$	(6*c*)	S III
$\mathbf{B}_{14}$	=	$z_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$\frac{1}{2}\,(-x_5 + z_5)\,a\,\hat{\mathbf{x}} + \left(-\frac{1}{2\sqrt{3}}x_5 + \frac{1}{\sqrt{3}}y_5 - \frac{1}{2\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}} + \frac{1}{3}\,(x_5 + y_5 + z_5)\,c\,\hat{\mathbf{z}}$	(6*c*)	S III
$\mathbf{B}_{15}$	=	$y_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\frac{1}{2}\,(y_5 - z_5)\,a\,\hat{\mathbf{x}} + \left(\frac{1}{\sqrt{3}}x_5 - \frac{1}{2\sqrt{3}}y_5 - \frac{1}{2\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}} + \frac{1}{3}\,(x_5 + y_5 + z_5)\,c\,\hat{\mathbf{z}}$	(6*c*)	S III
$\mathbf{B}_{16}$	=	$x_5\,\mathbf{a}_1 + z_5\,\mathbf{a}_2 + y_5\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_5 - y_5)\,a\,\hat{\mathbf{x}} + \left(-\frac{1}{2\sqrt{3}}x_5 - \frac{1}{2\sqrt{3}}y_5 + \frac{1}{\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}} + \frac{1}{3}\,(x_5 + y_5 + z_5)\,c\,\hat{\mathbf{z}}$	(6*c*)	S III
$\mathbf{B}_{17}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_6 - z_6)\,a\,\hat{\mathbf{x}} + \left(-\frac{1}{2\sqrt{3}}x_6 + \frac{1}{\sqrt{3}}y_6 - \frac{1}{2\sqrt{3}}z_6\right)a\,\hat{\mathbf{y}} + \frac{1}{3}\,(x_6 + y_6 + z_6)\,c\,\hat{\mathbf{z}}$	(6*c*)	S IV

$$\mathbf{B}_{18} = z_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + y_6\,\mathbf{a}_3 = \tfrac{1}{2}(-y_6 + z_6)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{\sqrt{3}}x_6 - \tfrac{1}{2\sqrt{3}}y_6 - \tfrac{1}{2\sqrt{3}}z_6\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}(x_6 + y_6 + z_6)\,c\,\hat{\mathbf{z}} \quad (6c) \quad \text{S IV}$$

$$\mathbf{B}_{19} = y_6\,\mathbf{a}_1 + z_6\,\mathbf{a}_2 + x_6\,\mathbf{a}_3 = \tfrac{1}{2}(-x_6 + y_6)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_6 - \tfrac{1}{2\sqrt{3}}y_6 + \tfrac{1}{\sqrt{3}}z_6\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}(x_6 + y_6 + z_6)\,c\,\hat{\mathbf{z}} \quad (6c) \quad \text{S IV}$$

$$\mathbf{B}_{20} = z_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + x_6\,\mathbf{a}_3 = \tfrac{1}{2}(-x_6 + z_6)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_6 + \tfrac{1}{\sqrt{3}}y_6 - \tfrac{1}{2\sqrt{3}}z_6\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}(x_6 + y_6 + z_6)\,c\,\hat{\mathbf{z}} \quad (6c) \quad \text{S IV}$$

$$\mathbf{B}_{21} = y_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 = \tfrac{1}{2}(y_6 - z_6)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{\sqrt{3}}x_6 - \tfrac{1}{2\sqrt{3}}y_6 - \tfrac{1}{2\sqrt{3}}z_6\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}(x_6 + y_6 + z_6)\,c\,\hat{\mathbf{z}} \quad (6c) \quad \text{S IV}$$

$$\mathbf{B}_{22} = x_6\,\mathbf{a}_1 + z_6\,\mathbf{a}_2 + y_6\,\mathbf{a}_3 = \tfrac{1}{2}(x_6 - y_6)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_6 - \tfrac{1}{2\sqrt{3}}y_6 + \tfrac{1}{\sqrt{3}}z_6\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}(x_6 + y_6 + z_6)\,c\,\hat{\mathbf{z}} \quad (6c) \quad \text{S IV}$$

$$\mathbf{B}_{23} = x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 = \tfrac{1}{2}(x_7 - z_7)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_7 + \tfrac{1}{\sqrt{3}}y_7 - \tfrac{1}{2\sqrt{3}}z_7\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}(x_7 + y_7 + z_7)\,c\,\hat{\mathbf{z}} \quad (6c) \quad \text{S V}$$

$$\mathbf{B}_{24} = z_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + y_7\,\mathbf{a}_3 = \tfrac{1}{2}(-y_7 + z_7)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{\sqrt{3}}x_7 - \tfrac{1}{2\sqrt{3}}y_7 - \tfrac{1}{2\sqrt{3}}z_7\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}(x_7 + y_7 + z_7)\,c\,\hat{\mathbf{z}} \quad (6c) \quad \text{S V}$$

$$\mathbf{B}_{25} = y_7\,\mathbf{a}_1 + z_7\,\mathbf{a}_2 + x_7\,\mathbf{a}_3 = \tfrac{1}{2}(-x_7 + y_7)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_7 - \tfrac{1}{2\sqrt{3}}y_7 + \tfrac{1}{\sqrt{3}}z_7\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}(x_7 + y_7 + z_7)\,c\,\hat{\mathbf{z}} \quad (6c) \quad \text{S V}$$

$$\mathbf{B}_{26} = z_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + x_7\,\mathbf{a}_3 = \tfrac{1}{2}(-x_7 + z_7)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_7 + \tfrac{1}{\sqrt{3}}y_7 - \tfrac{1}{2\sqrt{3}}z_7\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}(x_7 + y_7 + z_7)\,c\,\hat{\mathbf{z}} \quad (6c) \quad \text{S V}$$

$$\mathbf{B}_{27} = y_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 = \tfrac{1}{2}(y_7 - z_7)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{\sqrt{3}}x_7 - \tfrac{1}{2\sqrt{3}}y_7 - \tfrac{1}{2\sqrt{3}}z_7\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}(x_7 + y_7 + z_7)\,c\,\hat{\mathbf{z}} \quad (6c) \quad \text{S V}$$

$$\mathbf{B}_{28} = x_7\,\mathbf{a}_1 + z_7\,\mathbf{a}_2 + y_7\,\mathbf{a}_3 = \tfrac{1}{2}(x_7 - y_7)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_7 - \tfrac{1}{2\sqrt{3}}y_7 + \tfrac{1}{\sqrt{3}}z_7\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}(x_7 + y_7 + z_7)\,c\,\hat{\mathbf{z}} \quad (6c) \quad \text{S V}$$

References:

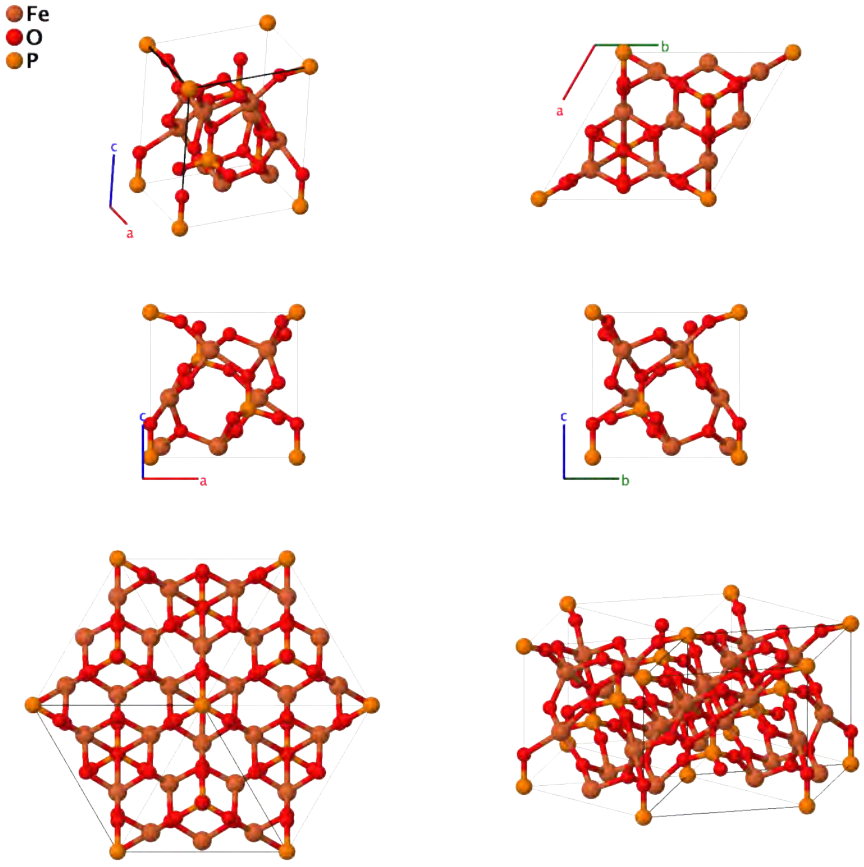
- T. Bjorvatten, O. Hassel, and A. Lindheim, *Crystal Structure of the Addition Compound SbI_3:3S_8*, Acta Chem. Scand. **17**, 689–702 (1963), doi:10.3891/acta.chem.scand.17-0689.

Geometry files:

- CIF: pp. 1725
- POSCAR: pp. 1726

Fe_3PO_7 Structure: A3B7C_hR11_160_b_a2b_a

http://aflow.org/prototype-encyclopedia/A3B7C_hR11_160_b_a2b_a

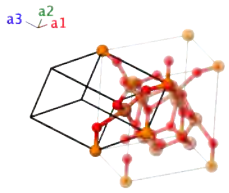


Prototype	:	Fe_3O_7P
AFLOW prototype label	:	A3B7C_hR11_160_b_a2b_a
***Strukturbericht* designation**	:	None
Pearson symbol	:	hR11
Space group number	:	160
Space group symbol	:	*R*3*m*
AFLOW prototype command	:	`aflow --proto=A3B7C_hR11_160_b_a2b_a [--hex] --params=`$a, c/a, x_1, x_2, x_3, z_3, x_4, z_4, x_5, z_5$

- Space group *R*3*m* #160 allows an arbitary definition of the zero of the z-axis. Here we select $z_1 = 0$, putting the phosphorus atom at the origin.

Rhombohedral primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}\,a\,\hat{\mathbf{y}} + \frac{1}{3}\,c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{\sqrt{3}}\,a\,\hat{\mathbf{y}} + \frac{1}{3}\,c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = -\frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}\,a\,\hat{\mathbf{y}} + \frac{1}{3}\,c\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$= x_1\,\mathbf{a}_1 + x_1\,\mathbf{a}_2 + x_1\,\mathbf{a}_3$	$= x_1 c\,\hat{\mathbf{z}}$	(1*a*)	O I
$\mathbf{B}_2$	$= x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	$= x_2 c\,\hat{\mathbf{z}}$	(1*a*)	P
$\mathbf{B}_3$	$= x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$= \frac{1}{2}\,(x_3 - z_3)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}\,(x_3 - z_3)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_3 + \frac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	(3*b*)	Fe
$\mathbf{B}_4$	$= z_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	$= \frac{1}{2}\,(-x_3 + z_3)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}\,(x_3 - z_3)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_3 + \frac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	(3*b*)	Fe
$\mathbf{B}_5$	$= x_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	$= \frac{1}{\sqrt{3}}\,(-x_3 + z_3)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_3 + \frac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	(3*b*)	Fe
$\mathbf{B}_6$	$= x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$= \frac{1}{2}\,(x_4 - z_4)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}\,(x_4 - z_4)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_4 + \frac{1}{3}z_4\right)c\,\hat{\mathbf{z}}$	(3*b*)	O II
$\mathbf{B}_7$	$= z_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	$= \frac{1}{2}\,(-x_4 + z_4)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}\,(x_4 - z_4)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_4 + \frac{1}{3}z_4\right)c\,\hat{\mathbf{z}}$	(3*b*)	O II
$\mathbf{B}_8$	$= x_4\,\mathbf{a}_1 + z_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	$= \frac{1}{\sqrt{3}}\,(-x_4 + z_4)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_4 + \frac{1}{3}z_4\right)c\,\hat{\mathbf{z}}$	(3*b*)	O II
$\mathbf{B}_9$	$= x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	$= \frac{1}{2}\,(x_5 - z_5)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}\,(x_5 - z_5)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_5 + \frac{1}{3}z_5\right)c\,\hat{\mathbf{z}}$	(3*b*)	O III
$\mathbf{B}_{10}$	$= z_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	$= \frac{1}{2}\,(-x_5 + z_5)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}\,(x_5 - z_5)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_5 + \frac{1}{3}z_5\right)c\,\hat{\mathbf{z}}$	(3*b*)	O III
$\mathbf{B}_{11}$	$= x_5\,\mathbf{a}_1 + z_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	$= \frac{1}{\sqrt{3}}\,(-x_5 + z_5)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_5 + \frac{1}{3}z_5\right)c\,\hat{\mathbf{z}}$	(3*b*)	O III

References:

- A. Modaressi, A. Courtois, R. Gerardin, B. Malaman, and C. Gleitzer, *Fe_3PO_7, Un cas de coordinence 5 du fer trivalent, etude structurale et magnetique*, J. Solid State Chem. **47**, 245–255 (1983), doi:10.1016/0022-4596(83)90016-6.

Found in:

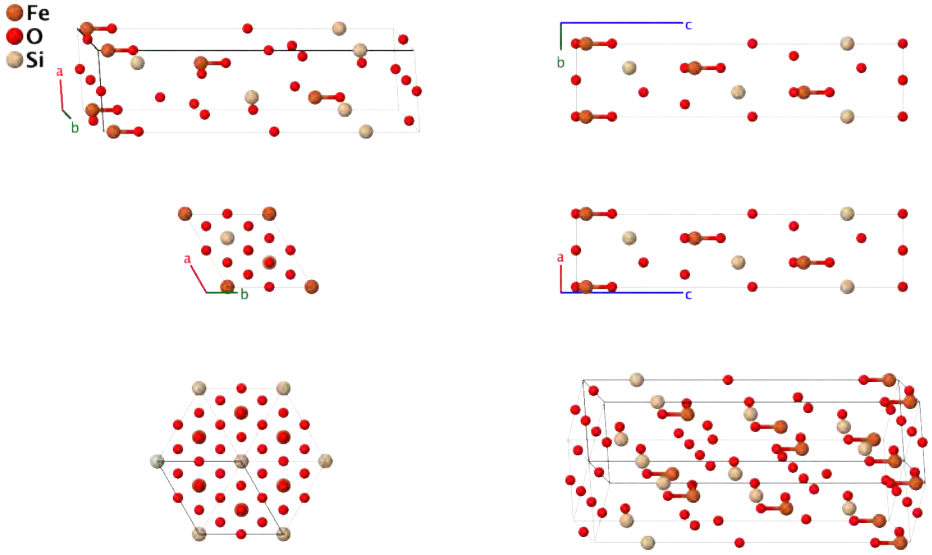
- C. L. Sarkis, M. J. Tarne, J. R. Neilson, H. B. Cao, E. Coldren, M. P. Gelfand, and K. A. Ross, *Partial antiferromagnetic helical order in single-crystal $Fe_3PO_4O_3$*, Phys. Rev. B **101**, 184417 (2020), doi:10.1103/PhysRevB.101.184417.

Geometry files:

- CIF: pp. 1726
- POSCAR: pp. 1726

Cronstedtite {Fe(Fe,Si)[(OH)$_2$,O]O$_3$, $S5_7$} Structure: AB3C2D_hR7_160_a_b_2a_a

http://aflow.org/prototype-encyclopedia/AB3C2D_hR7_160_a_b_2a_a

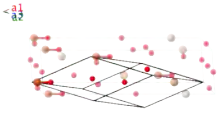


Prototype	:	$FeO_3(OH)_2Si$
AFLOW prototype label	:	AB3C2D_hR7_160_a_b_2a_a
***Strukturbericht* designation**	:	$S5_7$
Pearson symbol	:	hR7
Space group number	:	160
Space group symbol	:	$R3m$
AFLOW prototype command	:	`aflow --proto=AB3C2D_hR7_160_a_b_2a_a [--hex] --params=`$a, c/a, x_1, x_2, x_3, x_4, x_5, z_5$

- This particular structure of cronstedtite (also spelled cronstedite) is technically cronstedtite-3R, as there are three layers in the hexagonal cell, repeated in rhombohedral order. It is characterized by a high degree of disorder. The site we have labeled Si is better written as (Si,Fe), and the site labeled OH-I is actually [(OH)$_2$,O]. In addition the pure iron site, Fe ($1a$), is only occupied 90% of the time.
- As noted by (Frondel, 1962), there are a large number of varieties of cronstedtite, with each type having a different stacking sequence in the hexagonal cell.
- (Sterling, 1939) and (Hermann, 1943) put this structure in space group $R3$ #146, but the coordinates given by Sterling resolve to space group $R3m$ #160, and so we list this structure there.
- This is the final structure given a *Strukturbericht* designation.

Rhombohedral primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = -\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + x_1\,\mathbf{a}_2 + x_1\,\mathbf{a}_3$	=	$x_1 c\,\hat{\mathbf{z}}$	($1a$)	Fe
$\mathbf{B}_2$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$x_2 c\,\hat{\mathbf{z}}$	($1a$)	OH I
$\mathbf{B}_3$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3 c\,\hat{\mathbf{z}}$	($1a$)	OH II
$\mathbf{B}_4$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4 c\,\hat{\mathbf{z}}$	($1a$)	Si
$\mathbf{B}_5$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_5 - z_5)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}\,(x_5 - z_5)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_5 + \frac{1}{3}z_5\right) c\,\hat{\mathbf{z}}$	($3b$)	O
$\mathbf{B}_6$	=	$z_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$\frac{1}{2}\,(-x_5 + z_5)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}\,(x_5 - z_5)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_5 + \frac{1}{3}z_5\right) c\,\hat{\mathbf{z}}$	($3b$)	O
$\mathbf{B}_7$	=	$x_5\,\mathbf{a}_1 + z_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$\frac{1}{\sqrt{3}}\,(-x_5 + z_5)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_5 + \frac{1}{3}z_5\right) c\,\hat{\mathbf{z}}$	($3b$)	O

References:

- S. B. Hendricks, *Random structures of layer minerals as illustrated by cronstedite ($2FeO{\cdot}Fe_2O_3{\cdot}SiO_2{\cdot}2H_2O$). Possible iron content of kaolin*, Am. Mineral. **24**, 529–539 (1939).
- C. Frondel, *Polytypism in Cronstedtite*, Am. Mineral. **47**, 781–783 (1962).

Found in:

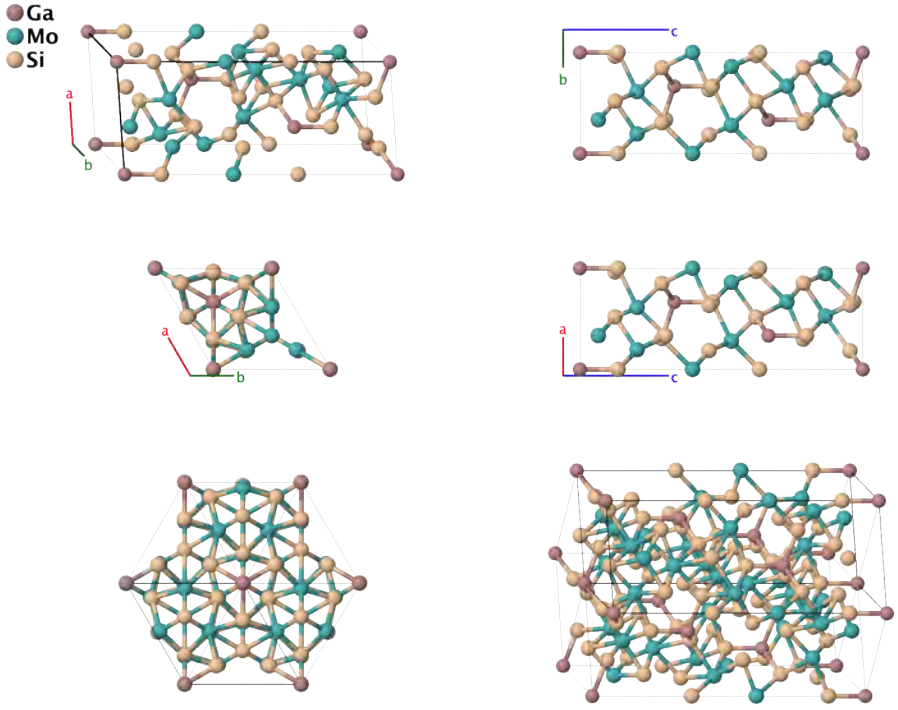
- K. Herrmann, ed., *Strukturbericht Band VII 1939* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1943).

Geometry files:

- CIF: pp. 1727
- POSCAR: pp. 1727

Low-Temperature $GaMo_4S_8$ Structure: AB4C8_hR13_160_a_ab_2a2b

http://aflow.org/prototype-encyclopedia/AB4C8_hR13_160_a_ab_2a2b

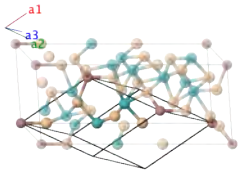


Prototype	:	$GaMo_4S_8$
AFLOW prototype label	:	AB4C8_hR13_160_a_ab_2a2b
***Strukturbericht* designation**	:	None
Pearson symbol	:	hR13
Space group number	:	160
Space group symbol	:	*R*3*m*
AFLOW prototype command	:	`aflow --proto=AB4C8_hR13_160_a_ab_2a2b [--hex] --params=`$a, c/a, x_1, x_2, x_3, x_4, x_5, z_5, x_6, z_6, x_7, z_7$

- At temperatures below 45 K $GaMo_4S_8$ transforms from its high-temperature cubic structure to this rhombohedral structure.
- We use the data from (François, 1991) at 8 K.
- Space group *R*3*m* #160 allows an arbitrary choice of the zero of the z-axis. Here we chose to place the gallium atom at the origin ($z_1 = 0$).

Rhombohedral primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}\,a\,\hat{\mathbf{y}} + \frac{1}{3}\,c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{\sqrt{3}}\,a\,\hat{\mathbf{y}} + \frac{1}{3}\,c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = -\frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}\,a\,\hat{\mathbf{y}} + \frac{1}{3}\,c\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$= x_1\,\mathbf{a}_1 + x_1\,\mathbf{a}_2 + x_1\,\mathbf{a}_3$	$= x_1 c\,\hat{\mathbf{z}}$	$(1a)$	Ga
$\mathbf{B}_2$	$= x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	$= x_2 c\,\hat{\mathbf{z}}$	$(1a)$	Mo I
$\mathbf{B}_3$	$= x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	$= x_3 c\,\hat{\mathbf{z}}$	$(1a)$	Si I
$\mathbf{B}_4$	$= x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	$= x_4 c\,\hat{\mathbf{z}}$	$(1a)$	Si II
$\mathbf{B}_5$	$= x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	$= \frac{1}{2}(x_5 - z_5)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}(x_5 - z_5)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_5 + \frac{1}{3}z_5\right)c\,\hat{\mathbf{z}}$	$(3b)$	Mo II
$\mathbf{B}_6$	$= z_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	$= \frac{1}{2}(-x_5 + z_5)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}(x_5 - z_5)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_5 + \frac{1}{3}z_5\right)c\,\hat{\mathbf{z}}$	$(3b)$	Mo II
$\mathbf{B}_7$	$= x_5\,\mathbf{a}_1 + z_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	$= \frac{1}{\sqrt{3}}(-x_5 + z_5)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_5 + \frac{1}{3}z_5\right)c\,\hat{\mathbf{z}}$	$(3b)$	Mo II
$\mathbf{B}_8$	$= x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	$= \frac{1}{2}(x_6 - z_6)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}(x_6 - z_6)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_6 + \frac{1}{3}z_6\right)c\,\hat{\mathbf{z}}$	$(3b)$	Si III
$\mathbf{B}_9$	$= z_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + x_6\,\mathbf{a}_3$	$= \frac{1}{2}(-x_6 + z_6)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}(x_6 - z_6)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_6 + \frac{1}{3}z_6\right)c\,\hat{\mathbf{z}}$	$(3b)$	Si III
$\mathbf{B}_{10}$	$= x_6\,\mathbf{a}_1 + z_6\,\mathbf{a}_2 + x_6\,\mathbf{a}_3$	$= \frac{1}{\sqrt{3}}(-x_6 + z_6)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_6 + \frac{1}{3}z_6\right)c\,\hat{\mathbf{z}}$	$(3b)$	Si III
$\mathbf{B}_{11}$	$= x_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	$= \frac{1}{2}(x_7 - z_7)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}(x_7 - z_7)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_7 + \frac{1}{3}z_7\right)c\,\hat{\mathbf{z}}$	$(3b)$	Si IV
$\mathbf{B}_{12}$	$= z_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + x_7\,\mathbf{a}_3$	$= \frac{1}{2}(-x_7 + z_7)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}(x_7 - z_7)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_7 + \frac{1}{3}z_7\right)c\,\hat{\mathbf{z}}$	$(3b)$	Si IV
$\mathbf{B}_{13}$	$= x_7\,\mathbf{a}_1 + z_7\,\mathbf{a}_2 + x_7\,\mathbf{a}_3$	$= \frac{1}{\sqrt{3}}(-x_7 + z_7)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_7 + \frac{1}{3}z_7\right)c\,\hat{\mathbf{z}}$	$(3b)$	Si IV

References:

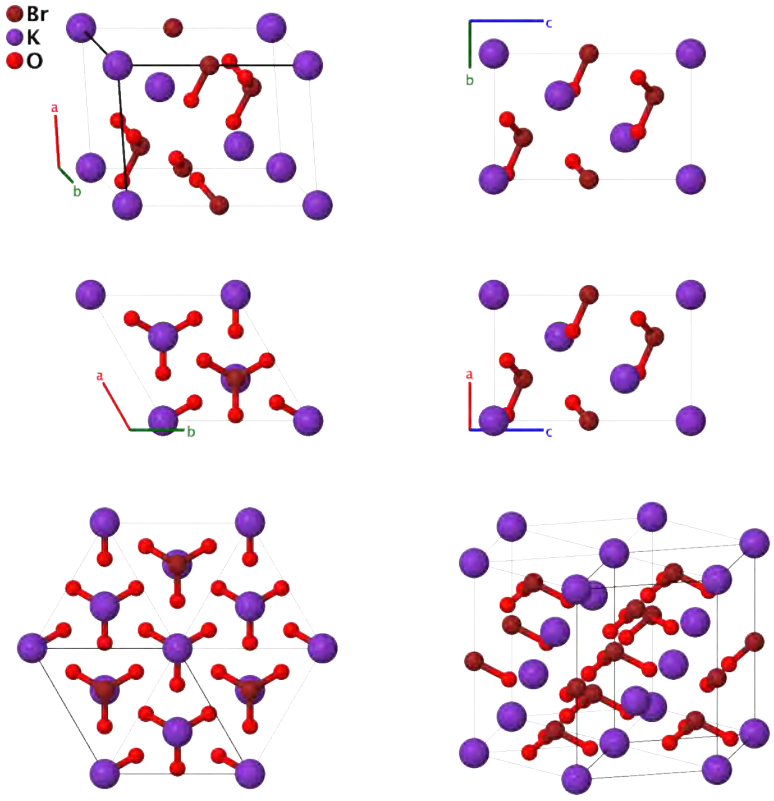
- M. François, W. Lengauer, K. Yvon, M. Sergent, M. Potel, P. Gougeon, and H. Ben Yaich-Aerrache, *Structural phase transition in $GaMo_4S_8$ by X-ray powder diffraction*, Zeitschrift für Kristallographie - Crystalline Materials **196**, 111–128 (1991), doi:10.1524/zkri.1991.196.14.111.

Geometry files:

- CIF: pp. 1727
- POSCAR: pp. 1728

$KBrO_3$ ($G0_7$) Structure: ABC3_hR5_160_a_a_b

http://aflow.org/prototype-encyclopedia/ABC3_hR5_160_a_a_b.KBrO3



Prototype	:	$BrKO_3$
AFLOW prototype label	:	ABC3_hR5_160_a_a_b
***Strukturbericht* designation**	:	$G0_7$
Pearson symbol	:	hR5
Space group number	:	160
Space group symbol	:	$R3m$
AFLOW prototype command	:	`aflow --proto=ABC3_hR5_160_a_a_b [--hex] --params=`$a, c/a, x_1, x_2, x_3, z_3$

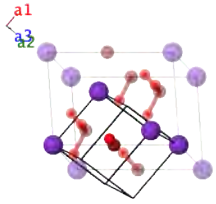
Other compounds with this structure

- KNO_3, γ-KNO_3, and $RbNO_3$

- γ-KNO_3 and $KBrO_3$ ($G0_7$) have the same AFLOW prototype label, ABC3_hR5_160_a_a_b. They are generated by the same symmetry operations with different sets of parameters (`--params`) specified in their corresponding CIF files.

Rhombohedral primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}\,a\,\hat{\mathbf{y}} + \frac{1}{3}\,c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{\sqrt{3}}\,a\,\hat{\mathbf{y}} + \frac{1}{3}\,c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = -\frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}\,a\,\hat{\mathbf{y}} + \frac{1}{3}\,c\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$x_1\,\mathbf{a}_1 + x_1\,\mathbf{a}_2 + x_1\,\mathbf{a}_3$	$x_1 c\,\hat{\mathbf{z}}$	(1*a*)	Br
$\mathbf{B}_2$	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	$x_2 c\,\hat{\mathbf{z}}$	(1*a*)	K
$\mathbf{B}_3$	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$\frac{1}{2}\,(x_3 - z_3)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}\,(x_3 - z_3)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_3 + \frac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	(3*b*)	O
$\mathbf{B}_4$	$z_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	$\frac{1}{2}\,(-x_3 + z_3)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}\,(x_3 - z_3)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_3 + \frac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	(3*b*)	O
$\mathbf{B}_5$	$x_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	$\frac{1}{\sqrt{3}}\,(-x_3 + z_3)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_3 + \frac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	(3*b*)	O

References:

- D. H. Templeton and L. K. Templeton, *Tensor X-ray optical properties of the bromate ion*, Acta Crystallogr. Sect. A pp. 133–142 (1985), doi:10.1107/S0108767385000277.

Found in:

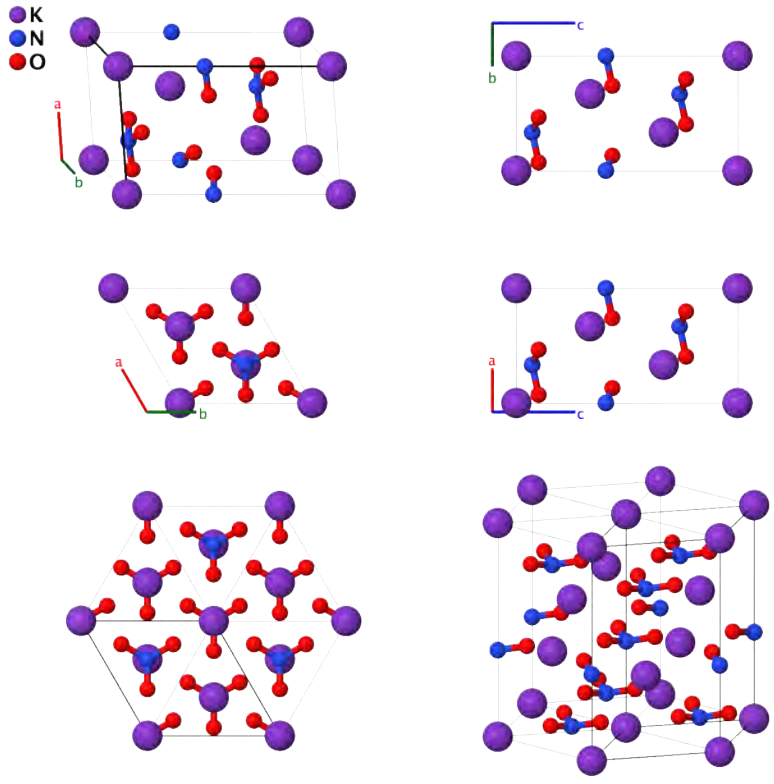
- D. Santamaría-Pérez, R. Chulia-Jordan, P. Rodríguez-Hernández, and A. Muñoz, *Crystal behavior of potassium bromate under compression*, Acta Crystallogr. Sect. B Struct. Sci. **71**, 798–804 (2015), doi:10.1107/S2052520615018156.

Geometry files:

- CIF: pp. 1728
- POSCAR: pp. 1728

γ-Potassium Nitrate (KNO_3) Structure: ABC3_hR5_160_a_a_b

http://aflow.org/prototype-encyclopedia/ABC3_hR5_160_a_a_b



Prototype	:	KNO_3
AFLOW prototype label	:	ABC3_hR5_160_a_a_b
***Strukturbericht* designation**	:	None
Pearson symbol	:	hR5
Space group number	:	160
Space group symbol	:	$R3m$
AFLOW prototype command	:	`aflow --proto=ABC3_hR5_160_a_a_b [--hex] --params=`$a, c/a, x_1, x_2, x_3, z_3$

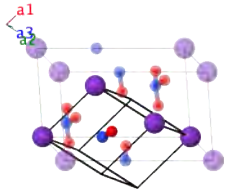
Other compounds with this structure

- NH_4ClO_4

- On heating, α-KNO_3 (either Structure I or Structure II) transforms into β-KNO_3 at 128 °C. When heated above 200 °C and then cooled, the β-phase transforms into the metastable ferroelectric γ-KNO_3 phase, which can remain down to room temperature.
- (Nimmo, 1976) give the data for γ-KNO_3 taken at 91 °C.
- Although this is isostructural with the $KBrO_3$ ($G0_7$) structure, we have included it here to facilitate the comparison of the various KNO_3 phases. γ-KNO_3 and $KBrO_3$ ($G0_7$) have the same AFLOW prototype label, ABC3_hR5_160_a_a_b. They are generated by the same symmetry operations with different sets of parameters (`--params`) specified in their corresponding CIF files.

Rhombohedral primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}\,a\,\hat{\mathbf{y}} + \frac{1}{3}\,c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{\sqrt{3}}\,a\,\hat{\mathbf{y}} + \frac{1}{3}\,c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = -\frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}\,a\,\hat{\mathbf{y}} + \frac{1}{3}\,c\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$x_1\,\mathbf{a}_1 + x_1\,\mathbf{a}_2 + x_1\,\mathbf{a}_3$	$x_1 c\,\hat{\mathbf{z}}$	($1a$)	K
$\mathbf{B}_2$	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	$x_2 c\,\hat{\mathbf{z}}$	($1a$)	N
$\mathbf{B}_3$	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$\frac{1}{2}\,(x_3 - z_3)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}\,(x_3 - z_3)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_3 + \frac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	($3b$)	O
$\mathbf{B}_4$	$z_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	$\frac{1}{2}\,(-x_3 + z_3)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}\,(x_3 - z_3)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_3 + \frac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	($3b$)	O
$\mathbf{B}_5$	$x_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	$\frac{1}{\sqrt{3}}\,(-x_3 + z_3)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_3 + \frac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	($3b$)	O

References:

- J. K. Nimmo and B. W. Lucas, *The crystal structures of γ-and β-KNO_3 and the α-β-γ phase transformations*, Acta Crystallogr. Sect. B Struct. Sci. **32**, 1968–1971 (1976), doi:10.1107/S0567740876006894.

Found in:

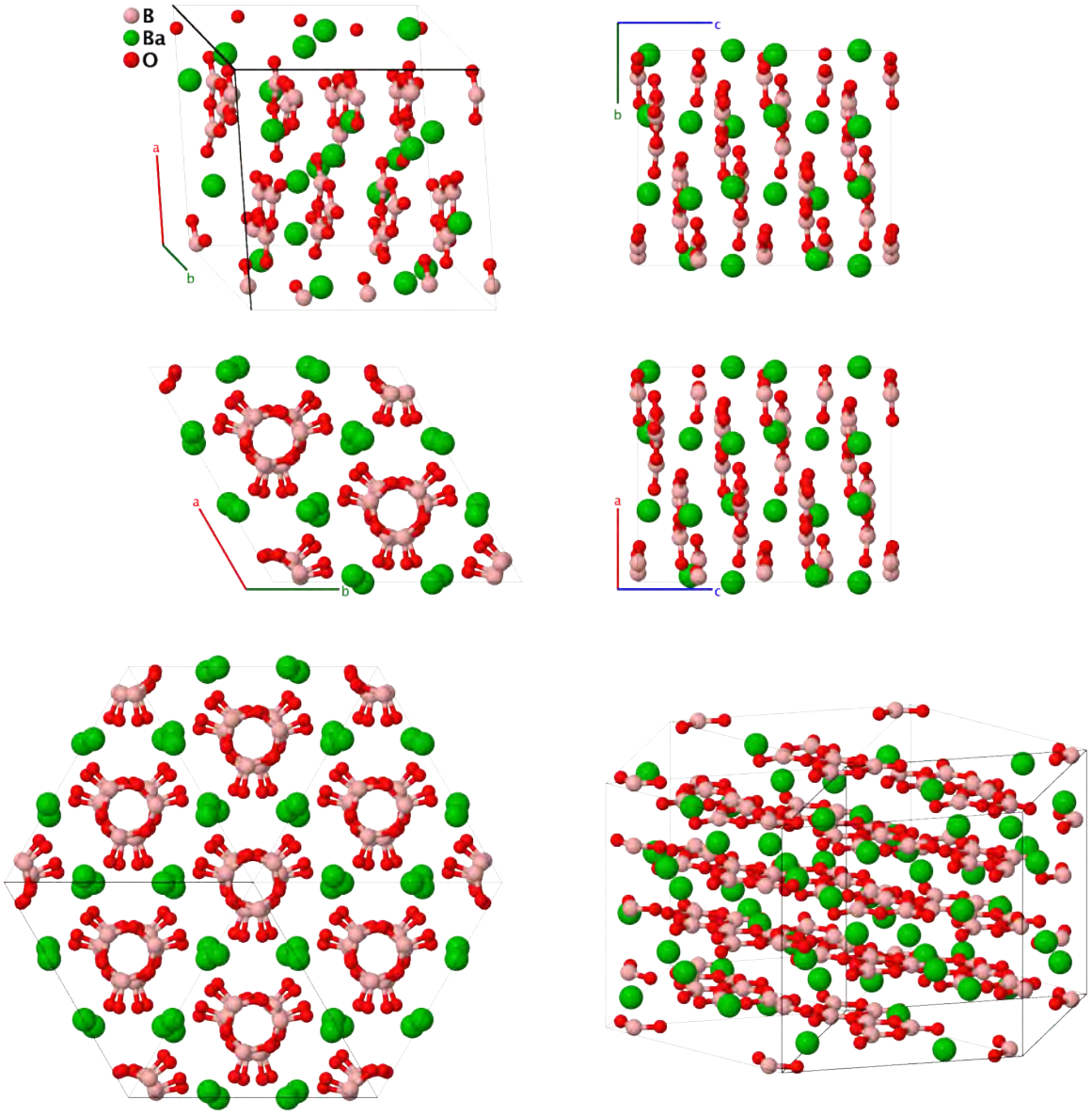
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1728
- POSCAR: pp. 1728

α-BaB_2O_4 (Low-Temperature) Structure: A2BC4_hR42_161_2b_b_4b

http://aflow.org/prototype-encyclopedia/A2BC4_hR42_161_2b_b_4b

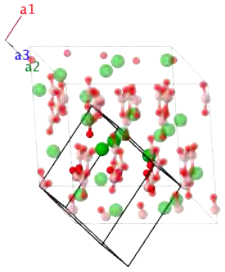


Prototype	:	B_2BaO_4
AFLOW prototype label	:	A2BC4_hR42_161_2b_b_4b
***Strukturbericht* designation**	:	None
Pearson symbol	:	hR42
Space group number	:	161
Space group symbol	:	$R3c$
AFLOW prototype command	:	`aflow --proto=A2BC4_hR42_161_2b_b_4b [--hex] --params=`$a, c/a, x_1, y_1, z_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7$

- This is the low-temperature structure. Heating to temperatures between 100-400 °C transforms it into β-BaB_2O_4. The principle difference between the two forms is the lack of inversion symmetry in the low-temperature structure.

Rhombohedral primitive vectors:

$$\begin{aligned}
\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2\sqrt{3}}\,a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,c\,\hat{\mathbf{z}} \\
\mathbf{a}_2 &= \tfrac{1}{\sqrt{3}}\,a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,c\,\hat{\mathbf{z}} \\
\mathbf{a}_3 &= -\tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2\sqrt{3}}\,a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,c\,\hat{\mathbf{z}}
\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + y_1\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$\tfrac{1}{2}\,(x_1 - z_1)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_1 + \tfrac{1}{\sqrt{3}}y_1 - \tfrac{1}{2\sqrt{3}}z_1\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,(x_1 + y_1 + z_1)\,c\,\hat{\mathbf{z}}$	(6*b*)	B I
$\mathbf{B}_2$	=	$z_1\,\mathbf{a}_1 + x_1\,\mathbf{a}_2 + y_1\,\mathbf{a}_3$	=	$\tfrac{1}{2}\,(-y_1 + z_1)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{\sqrt{3}}x_1 - \tfrac{1}{2\sqrt{3}}y_1 - \tfrac{1}{2\sqrt{3}}z_1\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,(x_1 + y_1 + z_1)\,c\,\hat{\mathbf{z}}$	(6*b*)	B I
$\mathbf{B}_3$	=	$y_1\,\mathbf{a}_1 + z_1\,\mathbf{a}_2 + x_1\,\mathbf{a}_3$	=	$\tfrac{1}{2}\,(-x_1 + y_1)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_1 - \tfrac{1}{2\sqrt{3}}y_1 + \tfrac{1}{\sqrt{3}}z_1\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,(x_1 + y_1 + z_1)\,c\,\hat{\mathbf{z}}$	(6*b*)	B I
$\mathbf{B}_4$	=	$\left(\tfrac{1}{2} + z_1\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_1\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_1\right)\mathbf{a}_3$	=	$\tfrac{1}{2}\,(-x_1 + z_1)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_1 + \tfrac{1}{\sqrt{3}}y_1 - \tfrac{1}{2\sqrt{3}}z_1\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + \tfrac{1}{3}x_1 + \tfrac{1}{3}y_1 + \tfrac{1}{3}z_1\right)c\,\hat{\mathbf{z}}$	(6*b*)	B I
$\mathbf{B}_5$	=	$\left(\tfrac{1}{2} + y_1\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_1\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\tfrac{1}{2}\,(y_1 - z_1)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{\sqrt{3}}x_1 - \tfrac{1}{2\sqrt{3}}y_1 - \tfrac{1}{2\sqrt{3}}z_1\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + \tfrac{1}{3}x_1 + \tfrac{1}{3}y_1 + \tfrac{1}{3}z_1\right)c\,\hat{\mathbf{z}}$	(6*b*)	B I
$\mathbf{B}_6$	=	$\left(\tfrac{1}{2} + x_1\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + z_1\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + y_1\right)\mathbf{a}_3$	=	$\tfrac{1}{2}\,(x_1 - y_1)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_1 - \tfrac{1}{2\sqrt{3}}y_1 + \tfrac{1}{\sqrt{3}}z_1\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + \tfrac{1}{3}x_1 + \tfrac{1}{3}y_1 + \tfrac{1}{3}z_1\right)c\,\hat{\mathbf{z}}$	(6*b*)	B I
$\mathbf{B}_7$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\tfrac{1}{2}\,(x_2 - z_2)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_2 + \tfrac{1}{\sqrt{3}}y_2 - \tfrac{1}{2\sqrt{3}}z_2\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,(x_2 + y_2 + z_2)\,c\,\hat{\mathbf{z}}$	(6*b*)	B II
$\mathbf{B}_8$	=	$z_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + y_2\,\mathbf{a}_3$	=	$\tfrac{1}{2}\,(-y_2 + z_2)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{\sqrt{3}}x_2 - \tfrac{1}{2\sqrt{3}}y_2 - \tfrac{1}{2\sqrt{3}}z_2\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,(x_2 + y_2 + z_2)\,c\,\hat{\mathbf{z}}$	(6*b*)	B II
$\mathbf{B}_9$	=	$y_2\,\mathbf{a}_1 + z_2\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$\tfrac{1}{2}\,(-x_2 + y_2)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_2 - \tfrac{1}{2\sqrt{3}}y_2 + \tfrac{1}{\sqrt{3}}z_2\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,(x_2 + y_2 + z_2)\,c\,\hat{\mathbf{z}}$	(6*b*)	B II

$$\mathbf{B}_{10} = \left(\tfrac{1}{2}+z_2\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_2\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_2\right)\mathbf{a}_3 = \tfrac{1}{2}(-x_2+z_2)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_2+\tfrac{1}{\sqrt{3}}y_2-\tfrac{1}{2\sqrt{3}}z_2\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+\tfrac{1}{3}x_2+\tfrac{1}{3}y_2+\tfrac{1}{3}z_2\right)c\,\hat{\mathbf{z}} \quad (6b) \quad \text{B II}$$

$$\mathbf{B}_{11} = \left(\tfrac{1}{2}+y_2\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_2\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_2\right)\mathbf{a}_3 = \tfrac{1}{2}(y_2-z_2)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{\sqrt{3}}x_2-\tfrac{1}{2\sqrt{3}}y_2-\tfrac{1}{2\sqrt{3}}z_2\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+\tfrac{1}{3}x_2+\tfrac{1}{3}y_2+\tfrac{1}{3}z_2\right)c\,\hat{\mathbf{z}} \quad (6b) \quad \text{B II}$$

$$\mathbf{B}_{12} = \left(\tfrac{1}{2}+x_2\right)\mathbf{a}_1+\left(\tfrac{1}{2}+z_2\right)\mathbf{a}_2+\left(\tfrac{1}{2}+y_2\right)\mathbf{a}_3 = \tfrac{1}{2}(x_2-y_2)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_2-\tfrac{1}{2\sqrt{3}}y_2+\tfrac{1}{\sqrt{3}}z_2\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+\tfrac{1}{3}x_2+\tfrac{1}{3}y_2+\tfrac{1}{3}z_2\right)c\,\hat{\mathbf{z}} \quad (6b) \quad \text{B II}$$

$$\mathbf{B}_{13} = x_3\,\mathbf{a}_1+y_3\,\mathbf{a}_2+z_3\,\mathbf{a}_3 = \tfrac{1}{2}(x_3-z_3)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_3+\tfrac{1}{\sqrt{3}}y_3-\tfrac{1}{2\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}}+\tfrac{1}{3}(x_3+y_3+z_3)\,c\,\hat{\mathbf{z}} \quad (6b) \quad \text{Ba}$$

$$\mathbf{B}_{14} = z_3\,\mathbf{a}_1+x_3\,\mathbf{a}_2+y_3\,\mathbf{a}_3 = \tfrac{1}{2}(-y_3+z_3)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{\sqrt{3}}x_3-\tfrac{1}{2\sqrt{3}}y_3-\tfrac{1}{2\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}}+\tfrac{1}{3}(x_3+y_3+z_3)\,c\,\hat{\mathbf{z}} \quad (6b) \quad \text{Ba}$$

$$\mathbf{B}_{15} = y_3\,\mathbf{a}_1+z_3\,\mathbf{a}_2+x_3\,\mathbf{a}_3 = \tfrac{1}{2}(-x_3+y_3)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_3-\tfrac{1}{2\sqrt{3}}y_3+\tfrac{1}{\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}}+\tfrac{1}{3}(x_3+y_3+z_3)\,c\,\hat{\mathbf{z}} \quad (6b) \quad \text{Ba}$$

$$\mathbf{B}_{16} = \left(\tfrac{1}{2}+z_3\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_3\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_3\right)\mathbf{a}_3 = \tfrac{1}{2}(-x_3+z_3)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_3+\tfrac{1}{\sqrt{3}}y_3-\tfrac{1}{2\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+\tfrac{1}{3}x_3+\tfrac{1}{3}y_3+\tfrac{1}{3}z_3\right)c\,\hat{\mathbf{z}} \quad (6b) \quad \text{Ba}$$

$$\mathbf{B}_{17} = \left(\tfrac{1}{2}+y_3\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_3\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_3\right)\mathbf{a}_3 = \tfrac{1}{2}(y_3-z_3)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{\sqrt{3}}x_3-\tfrac{1}{2\sqrt{3}}y_3-\tfrac{1}{2\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+\tfrac{1}{3}x_3+\tfrac{1}{3}y_3+\tfrac{1}{3}z_3\right)c\,\hat{\mathbf{z}} \quad (6b) \quad \text{Ba}$$

$$\mathbf{B}_{18} = \left(\tfrac{1}{2}+x_3\right)\mathbf{a}_1+\left(\tfrac{1}{2}+z_3\right)\mathbf{a}_2+\left(\tfrac{1}{2}+y_3\right)\mathbf{a}_3 = \tfrac{1}{2}(x_3-y_3)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_3-\tfrac{1}{2\sqrt{3}}y_3+\tfrac{1}{\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+\tfrac{1}{3}x_3+\tfrac{1}{3}y_3+\tfrac{1}{3}z_3\right)c\,\hat{\mathbf{z}} \quad (6b) \quad \text{Ba}$$

$$\mathbf{B}_{19} = x_4\,\mathbf{a}_1+y_4\,\mathbf{a}_2+z_4\,\mathbf{a}_3 = \tfrac{1}{2}(x_4-z_4)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_4+\tfrac{1}{\sqrt{3}}y_4-\tfrac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}}+\tfrac{1}{3}(x_4+y_4+z_4)\,c\,\hat{\mathbf{z}} \quad (6b) \quad \text{O I}$$

$$\mathbf{B}_{20} = z_4\,\mathbf{a}_1+x_4\,\mathbf{a}_2+y_4\,\mathbf{a}_3 = \tfrac{1}{2}(-y_4+z_4)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{\sqrt{3}}x_4-\tfrac{1}{2\sqrt{3}}y_4-\tfrac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}}+\tfrac{1}{3}(x_4+y_4+z_4)\,c\,\hat{\mathbf{z}} \quad (6b) \quad \text{O I}$$

$$\mathbf{B}_{21} = y_4\,\mathbf{a}_1+z_4\,\mathbf{a}_2+x_4\,\mathbf{a}_3 = \tfrac{1}{2}(-x_4+y_4)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_4-\tfrac{1}{2\sqrt{3}}y_4+\tfrac{1}{\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}}+\tfrac{1}{3}(x_4+y_4+z_4)\,c\,\hat{\mathbf{z}} \quad (6b) \quad \text{O I}$$

$$\mathbf{B}_{22} = \left(\tfrac{1}{2}+z_4\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_4\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_4\right)\mathbf{a}_3 = \tfrac{1}{2}(-x_4+z_4)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_4+\tfrac{1}{\sqrt{3}}y_4-\tfrac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+\tfrac{1}{3}x_4+\tfrac{1}{3}y_4+\tfrac{1}{3}z_4\right)c\,\hat{\mathbf{z}} \quad (6b) \quad \text{O I}$$

$$\mathbf{B}_{23} = \left(\tfrac{1}{2}+y_4\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_4\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_4\right)\mathbf{a}_3 = \tfrac{1}{2}(y_4-z_4)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{\sqrt{3}}x_4-\tfrac{1}{2\sqrt{3}}y_4-\tfrac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+\tfrac{1}{3}x_4+\tfrac{1}{3}y_4+\tfrac{1}{3}z_4\right)c\,\hat{\mathbf{z}} \quad (6b) \quad \text{O I}$$

$$\mathbf{B}_{24} = \left(\tfrac{1}{2}+x_4\right)\mathbf{a}_1+\left(\tfrac{1}{2}+z_4\right)\mathbf{a}_2+\left(\tfrac{1}{2}+y_4\right)\mathbf{a}_3 = \tfrac{1}{2}(x_4-y_4)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_4-\tfrac{1}{2\sqrt{3}}y_4+\tfrac{1}{\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+\tfrac{1}{3}x_4+\tfrac{1}{3}y_4+\tfrac{1}{3}z_4\right)c\,\hat{\mathbf{z}} \quad (6b) \quad \text{O I}$$

$$\mathbf{B}_{25} = x_5\,\mathbf{a}_1+y_5\,\mathbf{a}_2+z_5\,\mathbf{a}_3 = \tfrac{1}{2}(x_5-z_5)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_5+\tfrac{1}{\sqrt{3}}y_5-\tfrac{1}{2\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}}+\tfrac{1}{3}(x_5+y_5+z_5)\,c\,\hat{\mathbf{z}} \quad (6b) \quad \text{O II}$$

$$\mathbf{B}_{26} = z_5\,\mathbf{a}_1+x_5\,\mathbf{a}_2+y_5\,\mathbf{a}_3 = \tfrac{1}{2}(-y_5+z_5)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{\sqrt{3}}x_5-\tfrac{1}{2\sqrt{3}}y_5-\tfrac{1}{2\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}}+\tfrac{1}{3}(x_5+y_5+z_5)\,c\,\hat{\mathbf{z}} \quad (6b) \quad \text{O II}$$

$$\mathbf{B}_{27} = y_5\,\mathbf{a}_1+z_5\,\mathbf{a}_2+x_5\,\mathbf{a}_3 = \tfrac{1}{2}(-x_5+y_5)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_5-\tfrac{1}{2\sqrt{3}}y_5+\tfrac{1}{\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}}+\tfrac{1}{3}(x_5+y_5+z_5)\,c\,\hat{\mathbf{z}} \quad (6b) \quad \text{O II}$$

$$\mathbf{B}_{28} = \left(\tfrac{1}{2}+z_5\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_5\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_5\right)\mathbf{a}_3 = \tfrac{1}{2}(-x_5+z_5)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_5+\tfrac{1}{\sqrt{3}}y_5-\tfrac{1}{2\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+\tfrac{1}{3}x_5+\tfrac{1}{3}y_5+\tfrac{1}{3}z_5\right)c\,\hat{\mathbf{z}} \quad (6b) \quad \text{O II}$$

$$\mathbf{B}_{29} = \left(\tfrac{1}{2}+y_5\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_5\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_5\right)\mathbf{a}_3 = \tfrac{1}{2}(y_5-z_5)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{\sqrt{3}}x_5-\tfrac{1}{2\sqrt{3}}y_5-\tfrac{1}{2\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+\tfrac{1}{3}x_5+\tfrac{1}{3}y_5+\tfrac{1}{3}z_5\right)c\,\hat{\mathbf{z}} \quad (6b) \quad \text{O II}$$

$$\mathbf{B}_{30} = \left(\tfrac{1}{2}+x_5\right)\mathbf{a}_1+\left(\tfrac{1}{2}+z_5\right)\mathbf{a}_2+\left(\tfrac{1}{2}+y_5\right)\mathbf{a}_3 = \tfrac{1}{2}(x_5-y_5)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_5-\tfrac{1}{2\sqrt{3}}y_5+\tfrac{1}{\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+\tfrac{1}{3}x_5+\tfrac{1}{3}y_5+\tfrac{1}{3}z_5\right)c\,\hat{\mathbf{z}} \quad (6b) \quad \text{O II}$$

$$\mathbf{B}_{31} = x_6\,\mathbf{a}_1+y_6\,\mathbf{a}_2+z_6\,\mathbf{a}_3 = \tfrac{1}{2}(x_6-z_6)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_6+\tfrac{1}{\sqrt{3}}y_6-\tfrac{1}{2\sqrt{3}}z_6\right)a\,\hat{\mathbf{y}}+\tfrac{1}{3}(x_6+y_6+z_6)\,c\,\hat{\mathbf{z}} \quad (6b) \quad \text{O III}$$

$$\mathbf{B}_{32} = z_6\,\mathbf{a}_1+x_6\,\mathbf{a}_2+y_6\,\mathbf{a}_3 = \tfrac{1}{2}(-y_6+z_6)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{\sqrt{3}}x_6-\tfrac{1}{2\sqrt{3}}y_6-\tfrac{1}{2\sqrt{3}}z_6\right)a\,\hat{\mathbf{y}}+\tfrac{1}{3}(x_6+y_6+z_6)\,c\,\hat{\mathbf{z}} \quad (6b) \quad \text{O III}$$

$$\mathbf{B}_{33} = y_6\,\mathbf{a}_1+z_6\,\mathbf{a}_2+x_6\,\mathbf{a}_3 = \tfrac{1}{2}(-x_6+y_6)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_6-\tfrac{1}{2\sqrt{3}}y_6+\tfrac{1}{\sqrt{3}}z_6\right)a\,\hat{\mathbf{y}}+\tfrac{1}{3}(x_6+y_6+z_6)\,c\,\hat{\mathbf{z}} \quad (6b) \quad \text{O III}$$

$$\mathbf{B}_{34} = \left(\tfrac{1}{2}+z_6\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_6\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_6\right)\mathbf{a}_3 = \tfrac{1}{2}(-x_6+z_6)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_6+\tfrac{1}{\sqrt{3}}y_6-\tfrac{1}{2\sqrt{3}}z_6\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+\tfrac{1}{3}x_6+\tfrac{1}{3}y_6+\tfrac{1}{3}z_6\right)c\,\hat{\mathbf{z}} \quad (6b) \quad \text{O III}$$

$$\mathbf{B}_{35} = \left(\tfrac{1}{2}+y_6\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_6\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_6\right)\mathbf{a}_3 = \tfrac{1}{2}(y_6-z_6)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{\sqrt{3}}x_6-\tfrac{1}{2\sqrt{3}}y_6-\tfrac{1}{2\sqrt{3}}z_6\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+\tfrac{1}{3}x_6+\tfrac{1}{3}y_6+\tfrac{1}{3}z_6\right)c\,\hat{\mathbf{z}} \quad (6b) \quad \text{O III}$$

$$\mathbf{B}_{36} = \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_1+\left(\tfrac{1}{2}+z_6\right)\mathbf{a}_2+\left(\tfrac{1}{2}+y_6\right)\mathbf{a}_3 = \tfrac{1}{2}(x_6-y_6)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_6-\tfrac{1}{2\sqrt{3}}y_6+\tfrac{1}{\sqrt{3}}z_6\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+\tfrac{1}{3}x_6+\tfrac{1}{3}y_6+\tfrac{1}{3}z_6\right)c\,\hat{\mathbf{z}} \quad (6b) \quad \text{O III}$$

$$\mathbf{B}_{37} = x_7\,\mathbf{a}_1+y_7\,\mathbf{a}_2+z_7\,\mathbf{a}_3 = \tfrac{1}{2}(x_7-z_7)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_7+\tfrac{1}{\sqrt{3}}y_7-\tfrac{1}{2\sqrt{3}}z_7\right)a\,\hat{\mathbf{y}}+\tfrac{1}{3}(x_7+y_7+z_7)\,c\,\hat{\mathbf{z}} \quad (6b) \quad \text{O IV}$$

$$\mathbf{B}_{38} = z_7\,\mathbf{a}_1+x_7\,\mathbf{a}_2+y_7\,\mathbf{a}_3 = \tfrac{1}{2}(-y_7+z_7)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{\sqrt{3}}x_7-\tfrac{1}{2\sqrt{3}}y_7-\tfrac{1}{2\sqrt{3}}z_7\right)a\,\hat{\mathbf{y}}+\tfrac{1}{3}(x_7+y_7+z_7)\,c\,\hat{\mathbf{z}} \quad (6b) \quad \text{O IV}$$

$$\mathbf{B}_{39} = y_7\,\mathbf{a}_1+z_7\,\mathbf{a}_2+x_7\,\mathbf{a}_3 = \tfrac{1}{2}(-x_7+y_7)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_7-\tfrac{1}{2\sqrt{3}}y_7+\tfrac{1}{\sqrt{3}}z_7\right)a\,\hat{\mathbf{y}}+\tfrac{1}{3}(x_7+y_7+z_7)\,c\,\hat{\mathbf{z}} \quad (6b) \quad \text{O IV}$$

$$\mathbf{B}_{40} = \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_7\right)\mathbf{a}_3 = \tfrac{1}{2}(-x_7+z_7)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_7+\tfrac{1}{\sqrt{3}}y_7-\tfrac{1}{2\sqrt{3}}z_7\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+\tfrac{1}{3}x_7+\tfrac{1}{3}y_7+\tfrac{1}{3}z_7\right)c\,\hat{\mathbf{z}} \quad (6b) \quad \text{O IV}$$

$$\mathbf{B}_{41} = \left(\tfrac{1}{2}+y_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 = \tfrac{1}{2}(y_7-z_7)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{\sqrt{3}}x_7-\tfrac{1}{2\sqrt{3}}y_7-\tfrac{1}{2\sqrt{3}}z_7\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+\tfrac{1}{3}x_7+\tfrac{1}{3}y_7+\tfrac{1}{3}z_7\right)c\,\hat{\mathbf{z}} \quad (6b) \quad \text{O IV}$$

$$\mathbf{B}_{42} = \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}+z_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}+y_7\right)\mathbf{a}_3 = \tfrac{1}{2}(x_7-y_7)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_7-\tfrac{1}{2\sqrt{3}}y_7+\tfrac{1}{\sqrt{3}}z_7\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+\tfrac{1}{3}x_7+\tfrac{1}{3}y_7+\tfrac{1}{3}z_7\right)c\,\hat{\mathbf{z}} \quad (6b) \quad \text{O IV}$$

References:

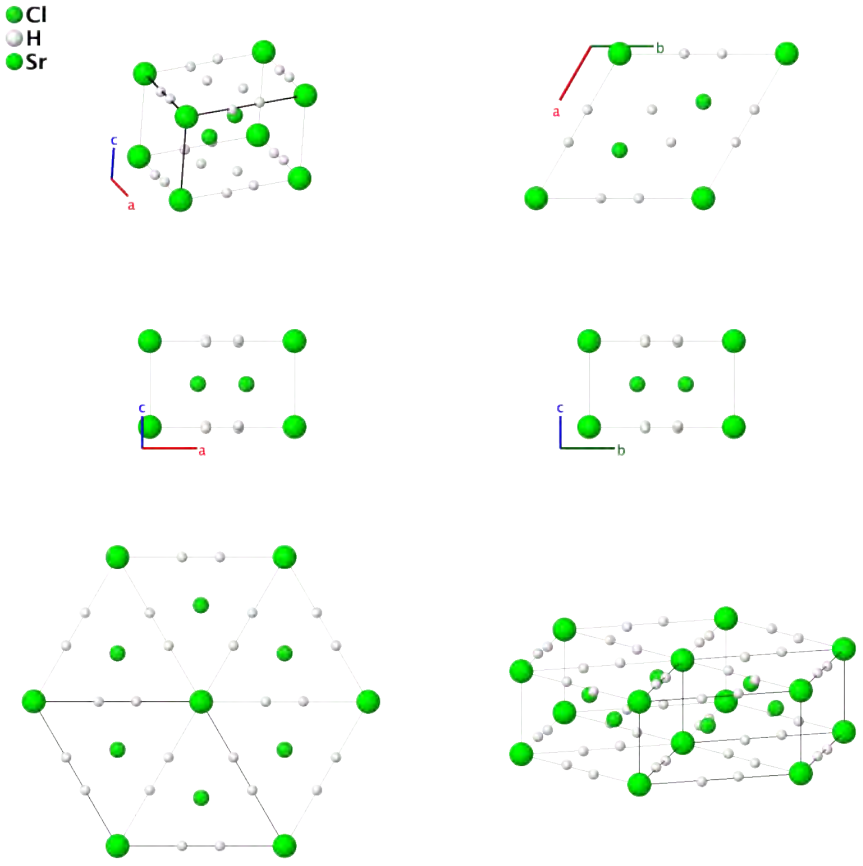
- R. Fröhlich, *Crystal Structure of the low-temperature form of* BaB_2O_4, Zeitschrift für Kristallographie - Crystalline Materials **168**, 109–112 (1984), doi:10.1524/zkri.1984.168.14.109.

Geometry files:

- CIF: pp. 1729
- POSCAR: pp. 1729

$I1_3$ ($SrCl_2{\cdot}(H_2O)_6$) (*obsolete*) Structure: A2B6C_hP9_162_d_k_a

http://aflow.org/prototype-encyclopedia/A2B6C_hP9_162_d_k_a



Prototype	:	$Cl_2(H_2O)_6Sr$
AFLOW prototype label	:	A2B6C_hP9_162_d_k_a
***Strukturbericht* designation**	:	$I1_3$
Pearson symbol	:	hP9
Space group number	:	162
Space group symbol	:	$P\bar{3}1m$
AFLOW prototype command	:	`aflow --proto=A2B6C_hP9_162_d_k_a --params=`$a, c/a, x_3, z_3$

Other compounds with this structure

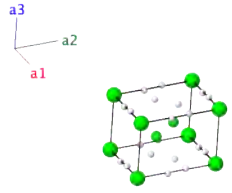
- $CaCl_2 \cdot (H_2O)_6$, $CaBr_2 \cdot (H_2O)_6$, $SrBr_2 \cdot (H_2O)_6$, $CaI_2 \cdot (H_2O)_6$, $SrI_2 \cdot (H_2O)_6$, and $BaI_2 \cdot (H_2O)_6$

- (Hermann, 1937) gives this the *Strukturbericht* designation $I1_3$, but gives the prototype as $K_2Pt(SCN)_6$. As we discussed on the $K_2Pt(SCN)_6$ ($H6_3$) page, the difference between these two structures is significant, so we will use the original $H6_3$ designation for $K_2Pt(SCN)_6$, and $I1_3$ for $SrCl_2 \cdot (H_2O)_6$.
- In any case, (Aagon, 1986) and others have shown that the correct space group of this structure is $P321$ #150. We discuss the corrected structure on the $SrCl_2 \cdot (H_2O)_6$ page.

- Using the notation of (Gottfried, 1937) this could also be designated the $J1_3$ structure. That designation was never used in any *Strukturbericht* volume, so we will use $I1_3$ here.
- The positions of the hydrogen atoms in the water molecules were not determined, so we only provide the positions of the oxygen atoms (labeled as H_2O).

Trigonal Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(1a)$	Sr
$\mathbf{B}_2$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(2d)$	Cl
$\mathbf{B}_3$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(2d)$	Cl
$\mathbf{B}_4$	=	$x_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_3$	=	$\frac{1}{2}x_3a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	$(6k)$	H_2O
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{2}x_3a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	$(6k)$	H_2O
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + z_3c\,\hat{\mathbf{z}}$	$(6k)$	H_2O
$\mathbf{B}_7$	=	$-x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-\frac{1}{2}x_3a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	$(6k)$	H_2O
$\mathbf{B}_8$	=	$-x_3\,\mathbf{a}_1 - z_3\,\mathbf{a}_3$	=	$-\frac{1}{2}x_3a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	$(6k)$	H_2O
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} - z_3c\,\hat{\mathbf{z}}$	$(6k)$	H_2O

References:

- Z. Hermann, *Über die Struktur des Strontiumchlorid-Hexahydrats*, Z. Anorg. Allg. Chem. **187**, 231–236 (1930), doi:10.1002/zaac.19301870121.
- P. A. Agron and W. R. Busing, *Calcium and strontium dichloride hexahydrates by neutron diffraction*, Acta Crystallogr. C **42**, 141–143 (1986), doi:10.1107/S0108270186097007.
- C. Gottfried and F. Schossberger, eds., *Strukturbericht Band III 1933-1935* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Found in:

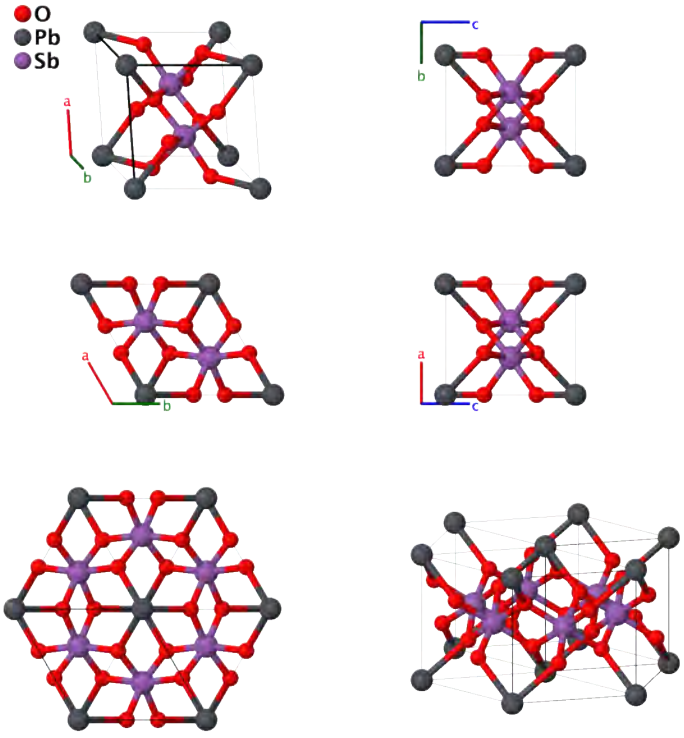
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1729
- POSCAR: pp. 1730

Rosiaite ($PbSb_2O_6$) Structure: A6BC2_hP9_162_k_a_d

http://aflow.org/prototype-encyclopedia/A6BC2_hP9_162_k_a_d



Prototype	:	O_6PbSb_2
AFLOW prototype label	:	A6BC2_hP9_162_k_a_d
***Strukturbericht* designation**	:	None
Pearson symbol	:	hP9
Space group number	:	162
Space group symbol	:	$P\bar{3}1m$
AFLOW prototype command	:	`aflow --proto=A6BC2_hP9_162_k_a_d --params=`$a, c/a, x_3, z_3$

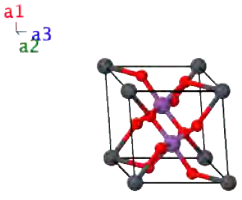
Other compounds with this structure

- $MnSeTeO_6$, $MnSnTeO_6$, and $PbTeGeO_6$

- This is the ternary form of the $L'3_2$ (β-V_2N) structure.

Trigonal Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(1a)$	Pb
$\mathbf{B}_2$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(2d)$	Sb
$\mathbf{B}_3$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(2d)$	Sb
$\mathbf{B}_4$	=	$x_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_3$	=	$\frac{1}{2}x_3a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	$(6k)$	O
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{2}x_3a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	$(6k)$	O
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + z_3c\,\hat{\mathbf{z}}$	$(6k)$	O
$\mathbf{B}_7$	=	$-x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-\frac{1}{2}x_3a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	$(6k)$	O
$\mathbf{B}_8$	=	$-x_3\,\mathbf{a}_1 - z_3\,\mathbf{a}_3$	=	$-\frac{1}{2}x_3a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	$(6k)$	O
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} - z_3c\,\hat{\mathbf{z}}$	$(6k)$	O

References:

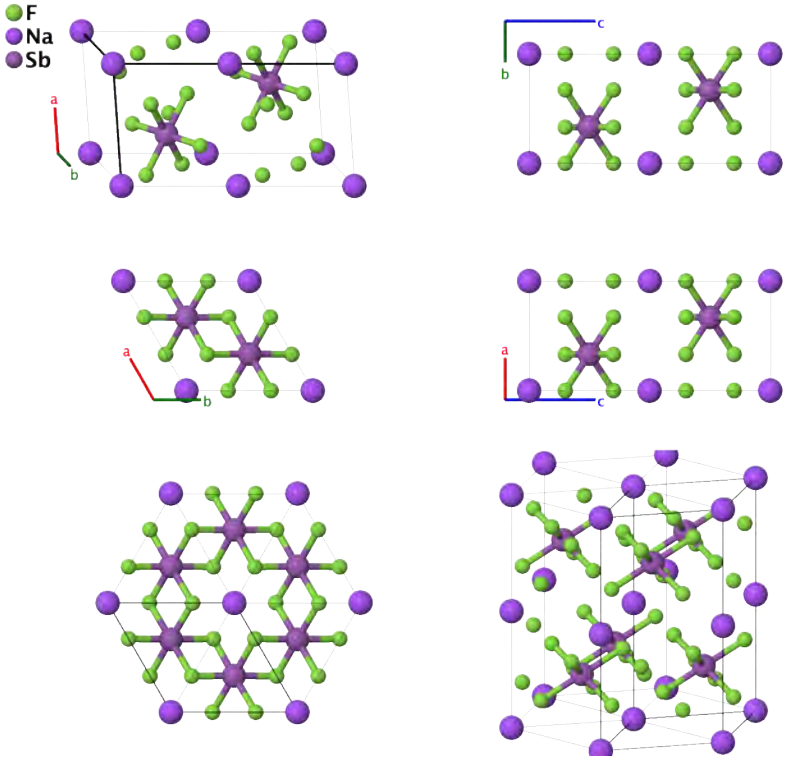
- R. Basso, G. Lucchetti, L. Zefiro, and A. Palenzona, *Rosiaite, $PbSb_2O_6$, a new mineral from the Cetine mine, Siena, Italy*, Eur. J. Mineral. **8**, 487–492 (1996), doi:10.1127/ejm/8/3/0487.

Geometry files:

- CIF: pp. 1730
- POSCAR: pp. 1730

$NaSbF_4(OH)_2$ ($J1_{12}$) Structure: A6BC_hP16_163_i_b_c

http://aflow.org/prototype-encyclopedia/A6BC_hP16_163_i_b_c

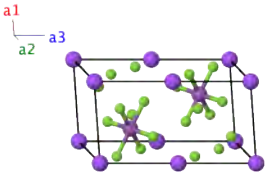


Prototype	:	F_4NaSb
AFLOW prototype label	:	A6BC_hP16_163_i_b_c
***Strukturbericht* designation**	:	$J1_{12}$
Pearson symbol	:	hP16
Space group number	:	163
Space group symbol	:	$P\bar{3}1c$
AFLOW prototype command	:	`aflow --proto=A6BC_hP16_163_i_b_c --params=`$a, c/a, x_3, y_3, z_3$

- The (12*i*) site, which we show as occupied by fluorine, is actually occupied by a random mixture of fluorine atoms (67%) and OH radicals (33%).
- Although the replacement of fluorine by OH does not affect the shape of the Sb-$(F,OH)_6$ ions, it has a profound effect on the structure, as can be seen by looking at $NaSbF_6$ and $NaSb(OH)_6$ ($J1_{11}$).

Trigonal Hexagonal primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(2*b*)	Na
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	(2*b*)	Na
$\mathbf{B}_3$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(2*c*)	Sb
$\mathbf{B}_4$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(2*c*)	Sb
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_3 + y_3)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_3 + y_3)\,a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(12*i*)	F
$\mathbf{B}_6$	=	$-y_3\,\mathbf{a}_1 + (x_3 - y_3)\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_3 - y_3\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(12*i*)	F
$\mathbf{B}_7$	=	$(-x_3 + y_3)\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\left(-x_3 + \frac{1}{2}y_3\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(12*i*)	F
$\mathbf{B}_8$	=	$-y_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$-\frac{1}{2}\,(x_3 + y_3)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_3 + y_3)\,a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(12*i*)	F
$\mathbf{B}_9$	=	$(-x_3 + y_3)\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_3 + y_3\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(12*i*)	F
$\mathbf{B}_{10}$	=	$x_3\,\mathbf{a}_1 + (x_3 - y_3)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(x_3 - \frac{1}{2}y_3\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	(12*i*)	F
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-\frac{1}{2}\,(x_3 + y_3)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(x_3 - y_3)\,a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(12*i*)	F
$\mathbf{B}_{12}$	=	$y_3\,\mathbf{a}_1 + (-x_3 + y_3)\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_3 + y_3\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(12*i*)	F
$\mathbf{B}_{13}$	=	$(x_3 - y_3)\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\left(x_3 - \frac{1}{2}y_3\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_3 a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(12*i*)	F
$\mathbf{B}_{14}$	=	$y_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\frac{1}{2}\,(x_3 + y_3)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(x_3 - y_3)\,a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(12*i*)	F
$\mathbf{B}_{15}$	=	$(x_3 - y_3)\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}x_3 - y_3\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(12*i*)	F
$\mathbf{B}_{16}$	=	$-x_3\,\mathbf{a}_1 + (-x_3 + y_3)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(-x_3 + \frac{1}{2}y_3\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(12*i*)	F

References:

- N. Schrewelius, *Röntgenuntersuchung der Verbindungen $NaSb(OH)_6$, $NaSbF_6$, $NaSbO_3$ und gleichartiger Stoffe*, Z. Anorg. Allg. Chem. **238**, 241–254 (1938), doi:10.1002/zaac.19382380209.

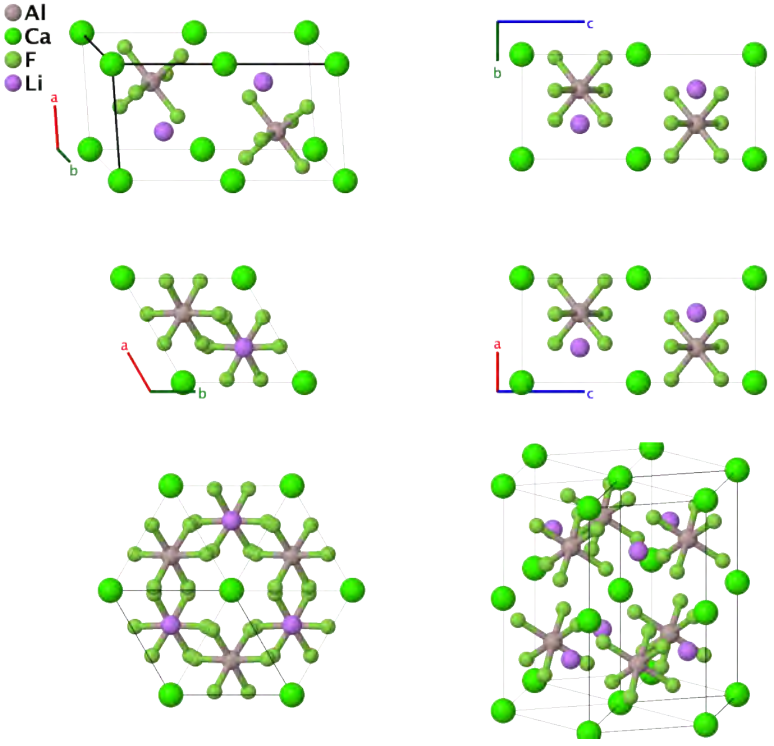
Found in:

- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:
- CIF: pp. 1730
- POSCAR: pp. 1731

Colquiriite ($LiCaAlF_6$) Structure: ABC6D_hP18_163_d_b_i_c

http://aflow.org/prototype-encyclopedia/ABC6D_hP18_163_d_b_i_c



Prototype	:	$AlCaF_6Li$
AFLOW prototype label	:	ABC6D_hP18_163_d_b_i_c
***Strukturbericht* designation**	:	None
Pearson symbol	:	hP18
Space group number	:	163
Space group symbol	:	$P\bar{3}1c$
AFLOW prototype command	:	`aflow --proto=ABC6D_hP18_163_d_b_i_c --params=`$a, c/a, x_4, y_4, z_4$

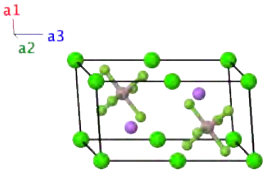
Other compounds with this structure

- $LiSrAlF_6$

- This is the low-temperature standard-pressure structure of $LiCaAlF_6$. We used the data taken at 300 K and ambient pressure.

Trigonal Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(2*b*)	Ca
$\mathbf{B}_2$	$\frac{1}{2}\,\mathbf{a}_3$	$\frac{1}{2}c\,\hat{\mathbf{z}}$	(2*b*)	Ca
$\mathbf{B}_3$	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(2*c*)	Li
$\mathbf{B}_4$	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(2*c*)	Li
$\mathbf{B}_5$	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(2*d*)	Al
$\mathbf{B}_6$	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(2*d*)	Al
$\mathbf{B}_7$	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$\frac{1}{2}(x_4 + y_4)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(-x_4 + y_4)\,a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(12*i*)	F
$\mathbf{B}_8$	$-y_4\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$\left(\frac{1}{2}x_4 - y_4\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(12*i*)	F
$\mathbf{B}_9$	$(-x_4 + y_4)\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$\left(-x_4 + \frac{1}{2}y_4\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_4 a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(12*i*)	F
$\mathbf{B}_{10}$	$-y_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	$-\frac{1}{2}(x_4 + y_4)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(-x_4 + y_4)\,a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(12*i*)	F
$\mathbf{B}_{11}$	$(-x_4 + y_4)\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	$\left(-\frac{1}{2}x_4 + y_4\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(12*i*)	F
$\mathbf{B}_{12}$	$x_4\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	$\left(x_4 - \frac{1}{2}y_4\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_4 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(12*i*)	F
$\mathbf{B}_{13}$	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	$-\frac{1}{2}(x_4 + y_4)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(x_4 - y_4)\,a\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(12*i*)	F
$\mathbf{B}_{14}$	$y_4\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	$\left(-\frac{1}{2}x_4 + y_4\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(12*i*)	F
$\mathbf{B}_{15}$	$(x_4 - y_4)\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	$\left(x_4 - \frac{1}{2}y_4\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_4 a\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(12*i*)	F
$\mathbf{B}_{16}$	$y_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	$\frac{1}{2}(x_4 + y_4)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(x_4 - y_4)\,a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(12*i*)	F
$\mathbf{B}_{17}$	$(x_4 - y_4)\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	$\left(\frac{1}{2}x_4 - y_4\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(12*i*)	F
$\mathbf{B}_{18}$	$-x_4\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	$\left(-x_4 + \frac{1}{2}y_4\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_4 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(12*i*)	F

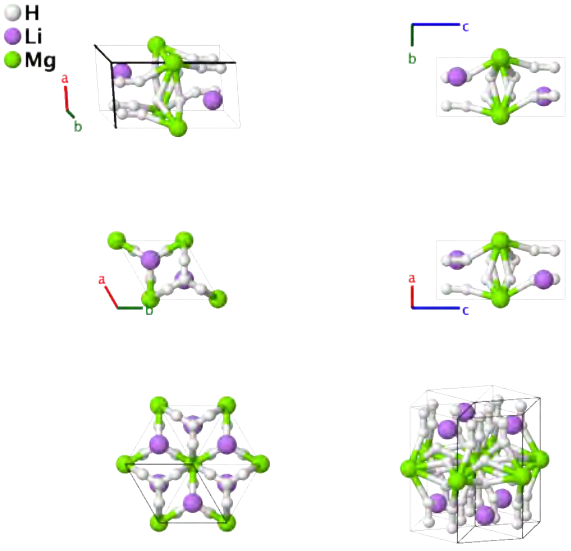
References:

- S. Kuze, D. du Boulay, N. Ishizawa, N. Kodama, M. Yamaga, and B. Henderson, *Structures of $LiCaAlF_6$ and $LiSrAlF_6$ at 120 and 300 K by synchrotron X-ray single-crystal diffraction*, J. Solid State Chem. **177**, 3505–3513 (2004), doi:10.1016/j.jssc.2004.04.039.

Geometry files:
- CIF: pp. 1731
- POSCAR: pp. 1731

Predicted Li_2MgH_{16} 300 GPa Structure: A16B2C_hP19_164_2d2i_d_b

http://aflow.org/prototype-encyclopedia/A16B2C_hP19_164_2d2i_d_b

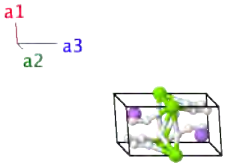


Prototype	:	$H_{16}Li_2Mg$
AFLOW prototype label	:	A16B2C_hP19_164_2d2i_d_b
***Strukturbericht* designation**	:	None
Pearson symbol	:	hP19
Space group number	:	164
Space group symbol	:	$P\bar{3}m1$
AFLOW prototype command	:	`aflow --proto=A16B2C_hP19_164_2d2i_d_b --params=`$a, c/a, z_2, z_3, z_4, x_5, z_5, x_6, z_6$

- This structure is predicted to be the zero-temperature ground state of Li_2MgH_{16} at 300 GPa. It is primarily of interest because a metastable cubic structure with the same composition is predicted to be superconducting at 250 GPa with $T_c = 430 - 473$ K.

Trigonal Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$ =	$\frac{1}{2}\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	$(1b)$	Mg

$\mathbf{B}_2$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(2*d*)	H I
$\mathbf{B}_3$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(2*d*)	H I
$\mathbf{B}_4$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(2*d*)	H II
$\mathbf{B}_5$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(2*d*)	H II
$\mathbf{B}_6$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(2*d*)	Li
$\mathbf{B}_7$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(2*d*)	Li
$\mathbf{B}_8$	=	$x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$-\sqrt{3}x_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(6*i*)	H III
$\mathbf{B}_9$	=	$x_5\,\mathbf{a}_1 + 2x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\frac{3}{2}x_5 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(6*i*)	H III
$\mathbf{B}_{10}$	=	$-2x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$-\frac{3}{2}x_5 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(6*i*)	H III
$\mathbf{B}_{11}$	=	$-x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\sqrt{3}x_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(6*i*)	H III
$\mathbf{B}_{12}$	=	$2x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\frac{3}{2}x_5 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(6*i*)	H III
$\mathbf{B}_{13}$	=	$-x_5\,\mathbf{a}_1 - 2x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-\frac{3}{2}x_5 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(6*i*)	H III
$\mathbf{B}_{14}$	=	$x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$-\sqrt{3}x_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(6*i*)	H IV
$\mathbf{B}_{15}$	=	$x_6\,\mathbf{a}_1 + 2x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\frac{3}{2}x_6 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(6*i*)	H IV
$\mathbf{B}_{16}$	=	$-2x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$-\frac{3}{2}x_6 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(6*i*)	H IV
$\mathbf{B}_{17}$	=	$-x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$\sqrt{3}x_6 a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(6*i*)	H IV
$\mathbf{B}_{18}$	=	$2x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$\frac{3}{2}x_6 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(6*i*)	H IV
$\mathbf{B}_{19}$	=	$-x_6\,\mathbf{a}_1 - 2x_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$-\frac{3}{2}x_6 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(6*i*)	H IV

References:

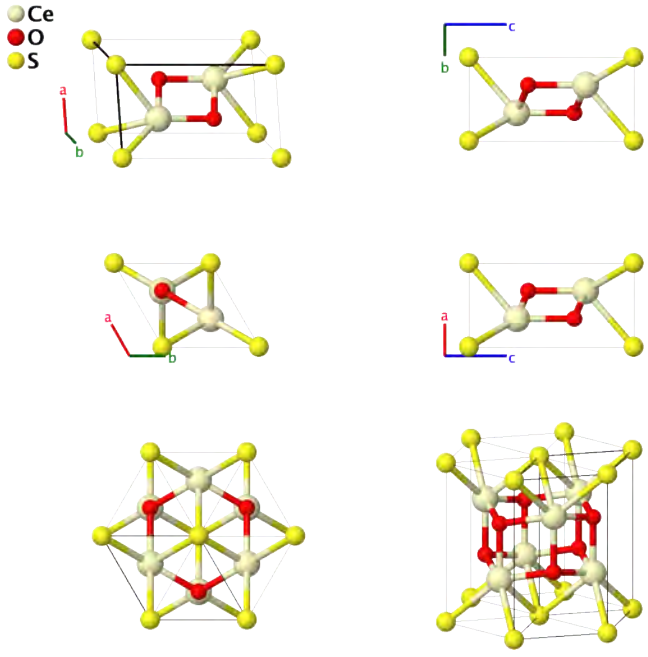
- Y. Sun, J. Lv, Y. Xie, H. Liu, and Y. Ma, *Route to a Superconducting Phase above Room Temperature in Electron-Doped Hydride Compounds under High Pressure*, Phys. Rev. Lett. **123**, 097001 (2019), doi:10.1103/PhysRevLett.123.097001.

Geometry files:

- CIF: pp. 1731
- POSCAR: pp. 1732

Ce_2O_2S Structure: A2B2C_hP5_164_d_d_a

http://aflow.org/prototype-encyclopedia/A2B2C_hP5_164_d_d_a



Prototype	:	Ce_2O_2S
AFLOW prototype label	:	A2B2C_hP5_164_d_d_a
***Strukturbericht* designation**	:	None
Pearson symbol	:	hP5
Space group number	:	164
Space group symbol	:	$P\bar{3}m1$
AFLOW prototype command	:	`aflow --proto=A2B2C_hP5_164_d_d_a --params=`$a, c/a, z_2, z_3$

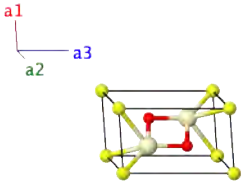
Other compounds with this structure

- Ce_2O_2S, Ce_2Se_2S, La_2O_2S, Pu_2O_2S, $CeCuZnP_2$, $DyCuZnP_2$, $ErCuZnP_2$, $GdCuZnP_2$, $HoCuZnP_2$, $LaCuZnP_2$, $LuCuZnP_2$, $NdCuZnP_2$, $PrCuZnP_2$, $ScCuZnP_2$, $SmCuZnP_2$, $TbCuZnP_2$, $TmCuZnP_2$, $YCuZnP_2$, $YbCuZnP_2$, $CaAl_2Si_2$, $CaAs_2Bi_2$, $CaAs_2Mn_2$, $CaAs_2P_2$, $CaAs_2Sb_2$, $CaBi_2Mn_2$, $SrAs_2Mn_2$, $SrAs_2P_2$, $SrAs_2Sb_2$, and $YbMnCuP_2$

- This is the ternary form of the $D5_2$ La_2O_3 structure and the $D5_{13}$ Al_3Ni_2 structure. We have separated it from the parents because of the large number of compounds involved.
- Authors after (Zachariasen, 1949) often use $CaAl_2Si_2$ as the prototype for this structure.

Trigonal Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(1a)$	S
$\mathbf{B}_2$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(2d)$	Ce
$\mathbf{B}_3$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	$(2d)$	Ce
$\mathbf{B}_4$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(2d)$	O
$\mathbf{B}_5$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(2d)$	O

References:

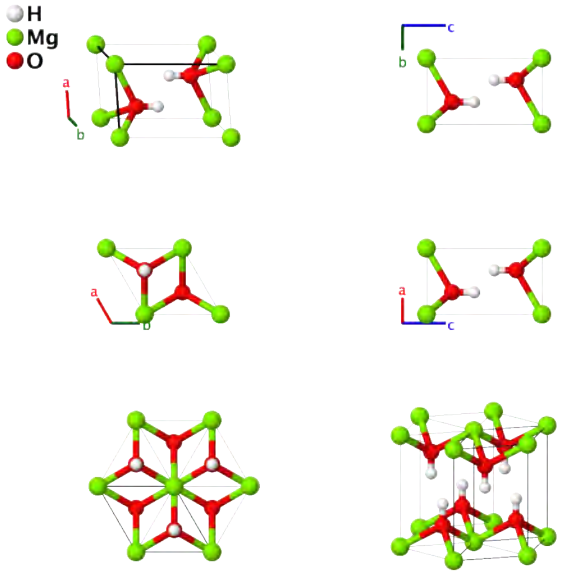
- W. H. Zachariasen, *Crystal chemical studies of the 5f-series of elements. VII. The crystal structure of Ce_2O_2S, La_2O_2S and Pu_2O_2S*, Acta Cryst. **2**, 60–62 (1949), doi:10.1107/S0365110X49000138.

Geometry files:

- CIF: pp. 1732
- POSCAR: pp. 1732

Brucite [$Mg(OH)_2$] Structure: A2BC2_hP5_164_d_a_d

http://aflow.org/prototype-encyclopedia/A2BC2_hP5_164_d_a_d



Prototype	:	H_2MgO_2
AFLOW prototype label	:	A2BC2_hP5_164_d_a_d
***Strukturbericht* designation**	:	None
Pearson symbol	:	hP5
Space group number	:	164
Space group symbol	:	$P\bar{3}m1$
AFLOW prototype command	:	`aflow --proto=A2BC2_hP5_164_d_a_d --params=`$a, c/a, z_2, z_3$

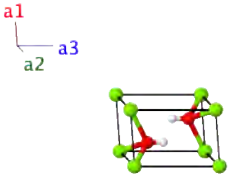
Other compounds with this structure

- $Ca(OH)_2$ (portlandite), $Fe(OH)_2$, $Mn(OH)_2$ (pyrochroite), $Ni(OH)_2$ (theophrastite), and β-$Co(OH)_2$

- We used the data from (Catti, 1995) at ambient pressure. In some brucite-like materials the hydrogen atoms are displaced to the (6*i*) Wyckoff positions $(x, -x, z)$ of space group #164, and these sites are 1/3 occupied.

Trigonal Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(1a)$	Mg
$\mathbf{B}_2$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(2d)$	H
$\mathbf{B}_3$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	$(2d)$	H
$\mathbf{B}_4$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(2d)$	O
$\mathbf{B}_5$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(2d)$	O

References:

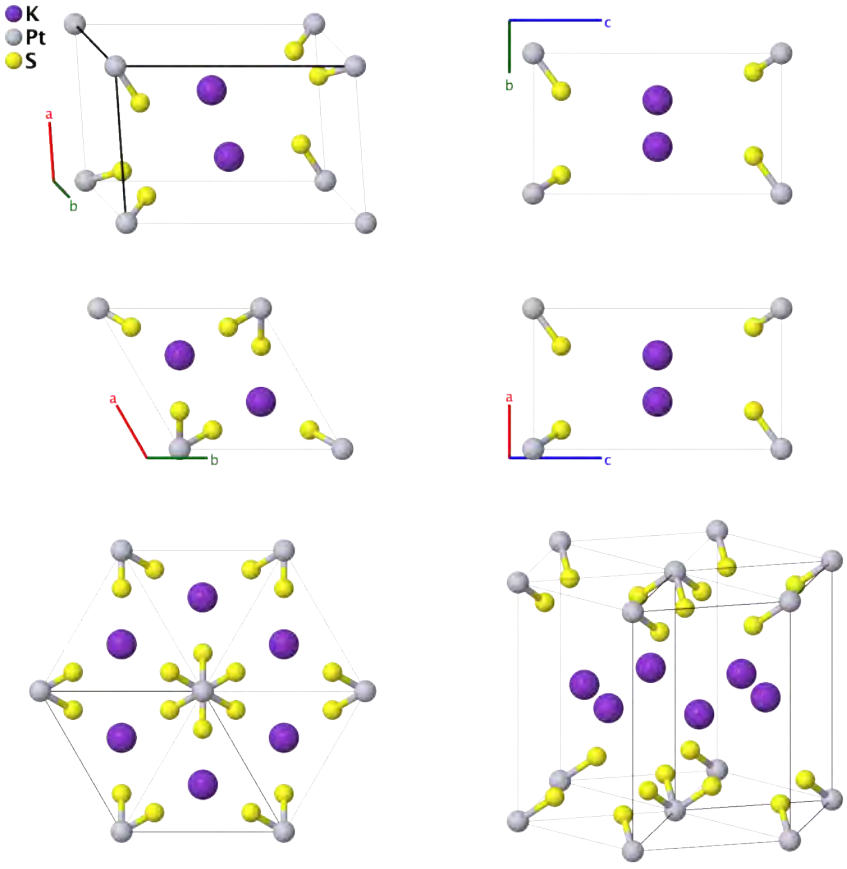
- M. Catti, G. Ferraris, S. Hull, and A. Pavese, *Static compression and H disorder in brucite, Mg(OH)$_2$, to 11 GPa: a powder neutron diffraction study*, Phys. Chem. Miner. **22**, 200–206 (1995), doi:10.1007/BF00202300.

Geometry files:

- CIF: pp. 1732
- POSCAR: pp. 1733

$K_2Pt(SCN)_6$ ($H6_3$) Structure: A2BC6_hP9_164_d_a_i

http://aflow.org/prototype-encyclopedia/A2BC6_hP9_164_d_a_i



Prototype	:	$K_2Pt(SCN)_6$
AFLOW prototype label	:	A2BC6_hP9_164_d_a_i
***Strukturbericht* designation**	:	$H6_3$
Pearson symbol	:	hP9
Space group number	:	164
Space group symbol	:	$P\bar{3}m1$
AFLOW prototype command	:	`aflow --proto=A2BC6_hP9_164_d_a_i --params=`$a, c/a, z_2, x_3, z_3$

Other compounds with this structure

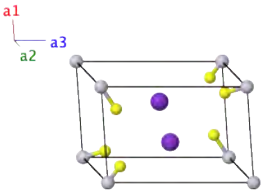
- $(NH4)_2Pt(SCN)_6$ and $Rb_2Pt(SCN)_6$

- The structure here must be regarded at tentative. Even the "official" labeling of the structure is uncertain:
  - (Hendricks, 1928) made the only structural study of this compound that we have been able to find. Unfortunately, they were not able to determine the positions of the carbon and silicon atoms. Presumably $(SCN)^-$ forms a straight-line ionic system making a roughly 105° angle with the Pt-S line, as it does in the hydrated form of this compound $K_2Pt(SCN)_6 \cdot 2H_2O$.
  - (Hendricks, 1928) were also uncertain about the space group, giving two possibilities: $P\bar{3}1m$ #162 and $P\bar{3}m1$ #164. They prefer the latter, so we present this as a $P\bar{3}m1$ #164 structure.

- The positions of the potassium and sulfur ions are not well determined. (Hendricks, 1928) give $z_2 \approx 0.5$, $x_3 = 0.10 - 0.17$, and $z_3 = 0.09 - 0.135$. We used the average values for the S coordinates.
- While (Ewald, 1931) gave this the *Strukturbericht* designation $H6_3$, (Hermann, 1937) lists it as $I1_3$, "previously $H6_3$." However, they then go on to describe structures of the form $SrCl_2(H_2O)_6$, with a very different c/a ratio.
- (Ewald, 1931) originally used the $H6$ *Strukturbericht* category for compounds of the form $B(X)_6$, but (Hermann, 1937) and following volumes use both I and J for $B(X)_6$, sometimes changing a previous $H6$ designation to $I1$ or $J1$. In general we will follow this change in notation, but given the striking difference in structures between $K_2Pt(SCN)_6$ and $SrCl_2(H_2O)_6$ we will continue to use $H6_3$ for $K_2Pt(SCN)_6$ and $I1_3$ for $SrCl_2(H_2O)_6$.

Trigonal Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(1a)$	Pt
$\mathbf{B}_2$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(2d)$	K
$\mathbf{B}_3$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	$(2d)$	K
$\mathbf{B}_4$	=	$x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$-\sqrt{3}x_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(6i)$	S
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + 2x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{3}{2}x_3 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(6i)$	S
$\mathbf{B}_6$	=	$-2x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$-\frac{3}{2}x_3 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(6i)$	S
$\mathbf{B}_7$	=	$-x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\sqrt{3}x_3 a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(6i)$	S
$\mathbf{B}_8$	=	$2x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\frac{3}{2}x_3 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(6i)$	S
$\mathbf{B}_9$	=	$-x_3\,\mathbf{a}_1 - 2x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-\frac{3}{2}x_3 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(6i)$	S

References:

- S. B. Hendricks and H. E. Merwin, *The atomic arrangement in crystals of alkali platini-thiocyanates*, Am. J. Sci. **15**, 487–494 (1928), doi:10.2475/ajs.s5-15.90.487.
- P. P. Ewald and C. Hermann, eds., *Strukturbericht 1913-1928* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1931).
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Found in:

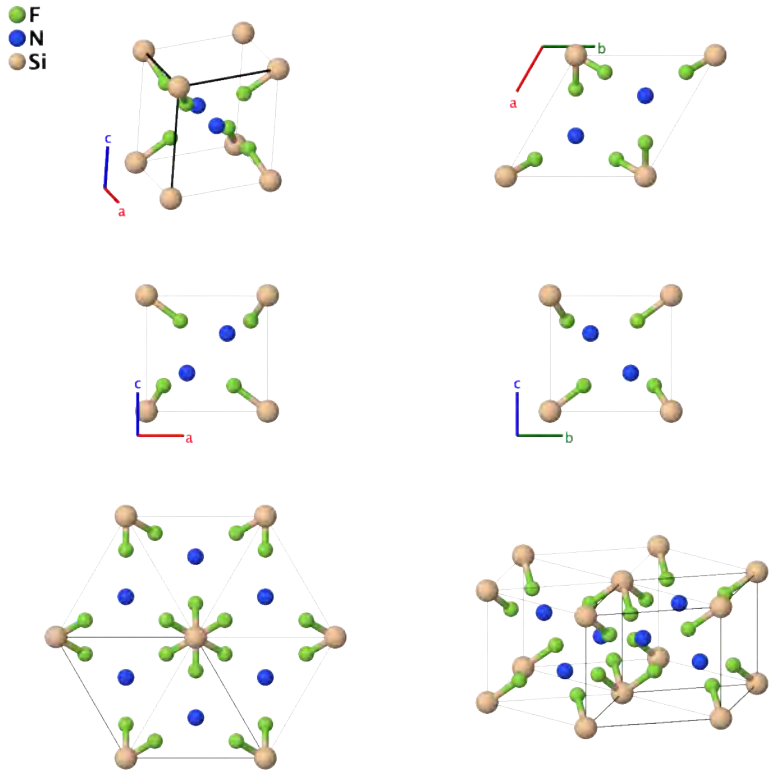
- P. P. Ewald and C. Hermann, eds., *Strukturbericht 1913-1928* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1931).

Geometry files:

- CIF: pp. 1733
- POSCAR: pp. 1733

Bararite (Trigonal $(NH_4)_2SiF_6$, $J1_6$) Structure: A6B2C_hP9_164_i_d_a

http://aflow.org/prototype-encyclopedia/A6B2C_hP9_164_i_d_a

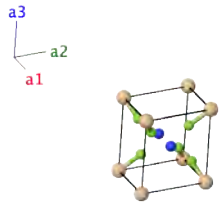


Prototype	:	$F_6(NH_4)_2Si$
AFLOW prototype label	:	A6B2C_hP9_164_i_d_a
***Strukturbericht* designation**	:	$J1_6$
Pearson symbol	:	hP9
Space group number	:	164
Space group symbol	:	$P\bar{3}m1$
AFLOW prototype command	:	`aflow --proto=A6B2C_hP9_164_i_d_a --params=`$a, c/a, z_2, x_3, z_3$

- Bararite is a trigonal form of $(NH_4)_2SiF_6$, metastable at room temperature. The room temperature stable form is cubic cryptohalite, which takes on the $J1_1$ structure. Except for the hydrogen atoms, this structure is very similar to $J1_{13}$, K_2GeF_6. (Schlemper, 1966) state that the hydrogen atoms are on ($2d$) and ($6i$) sites, but were not able to determine the coordinates because of large thermal fluctuations. They were to study the system at 77 K, but we have not found any evidence that this work was ever published.
- The positions of the hydrogen atoms in the NH_4 ions were not determined, so we only provide the positions of the nitrogen atoms (labeled as NH_4).

Trigonal Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(1*a*)	Si
$\mathbf{B}_2$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(2*d*)	NH_4
$\mathbf{B}_3$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(2*d*)	NH_4
$\mathbf{B}_4$	=	$x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$-\sqrt{3}x_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(6*i*)	F
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + 2x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{3}{2}x_3 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(6*i*)	F
$\mathbf{B}_6$	=	$-2x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$-\frac{3}{2}x_3 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(6*i*)	F
$\mathbf{B}_7$	=	$-x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\sqrt{3}x_3 a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(6*i*)	F
$\mathbf{B}_8$	=	$2x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\frac{3}{2}x_3 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(6*i*)	F
$\mathbf{B}_9$	=	$-x_3\,\mathbf{a}_1 - 2x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-\frac{3}{2}x_3 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(6*i*)	F

References:

- E. O. Schlemper and W. C. Hamilton, *On the Structure of Trigonal Ammonium Flourosilicate*, J. Chem. Phys. **45**, 408–409 (1966), doi:10.1063/1.2716548.

Found in:

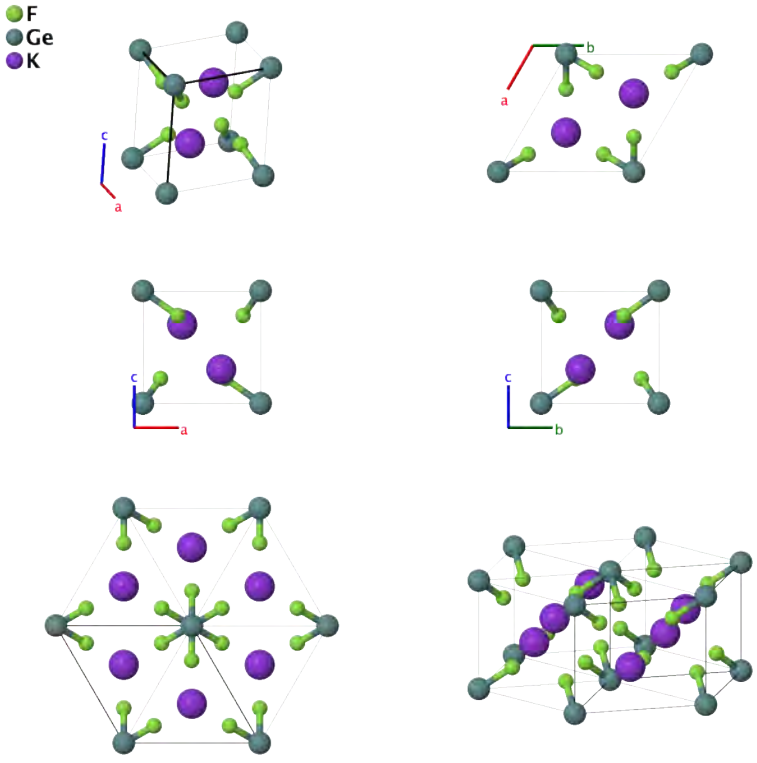
- J. Fábry, J. Chval, and V. Petrícek, *A new modification of diammonium hexafluorosilicate, $(NH_4)_2SiF_6$*, Acta Crystallogr. E **57**, i90–i91 (2001), doi:10.1107/S160053680101501X.

Geometry files:

- CIF: pp. 1733
- POSCAR: pp. 1733

K_2GeF_6 ($J1_{13}$) Structure: A6BC2_hP9_164_i_a_d

http://aflow.org/prototype-encyclopedia/A6BC2_hP9_164_i_a_d



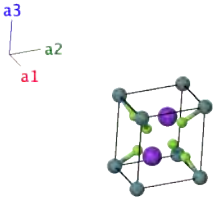
Prototype	:	F_6GeK_2
AFLOW prototype label	:	A6BC2_hP9_164_i_a_d
***Strukturbericht* designation**	:	$J1_{13}$
Pearson symbol	:	hP9
Space group number	:	164
Space group symbol	:	$P\bar{3}m1$
AFLOW prototype command	:	`aflow --proto=A6BC2_hP9_164_i_a_d --params=`$a, c/a, z_2, x_3, z_3$

Other compounds with this structure

- $(NH_4)_2GeF_6$, Cs_2CeCl_6, K_2PtF_6, and K_2SiF_6

Trigonal Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($1a$)	Ge
$\mathbf{B}_2$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	($2d$)	K
$\mathbf{B}_3$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	($2d$)	K
$\mathbf{B}_4$	=	$x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$-\sqrt{3}x_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	($6i$)	F
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + 2x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{3}{2}x_3 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	($6i$)	F
$\mathbf{B}_6$	=	$-2x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$-\frac{3}{2}x_3 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	($6i$)	F
$\mathbf{B}_7$	=	$-x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\sqrt{3}x_3 a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	($6i$)	F
$\mathbf{B}_8$	=	$2x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\frac{3}{2}x_3 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	($6i$)	F
$\mathbf{B}_9$	=	$-x_3\,\mathbf{a}_1 - 2x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-\frac{3}{2}x_3 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	($6i$)	F

References:

- J. L. Hoard and W. B. Vincent, *Structures of Complex Fluorides. Potassium Hexafluogermanate and Ammonium Hexafluogermanate*, J. Am. Chem. Soc. **61**, 2849–2852 (1939), doi:10.1021/ja01265a082.

Found in:

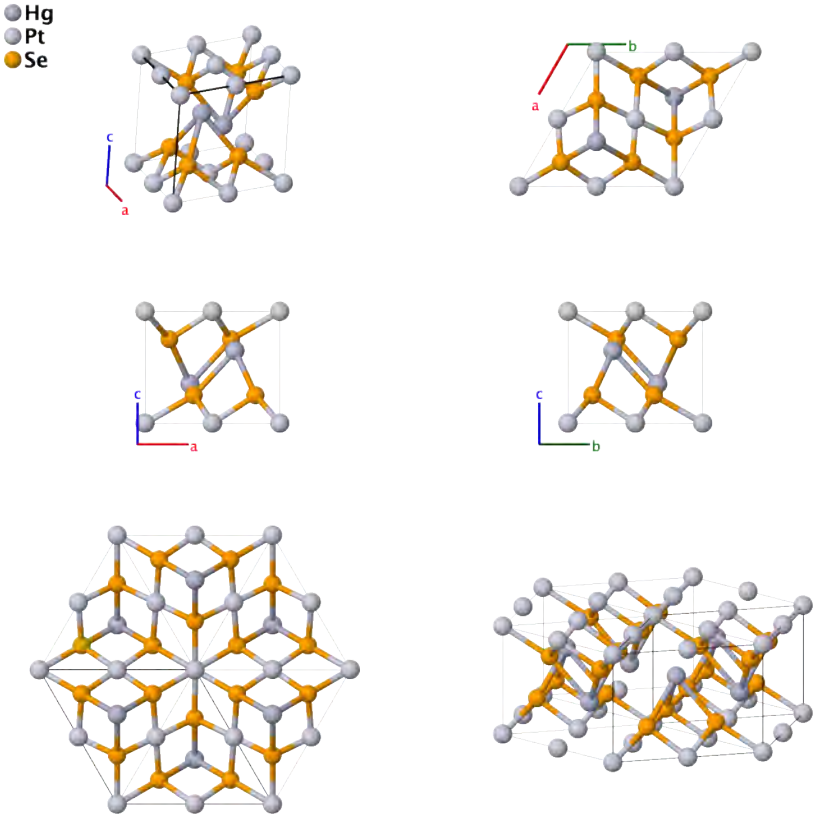
- T. Kaatz and M. Marcovich, *The crystal structure of the compound Cs_2CeCl_6*, Acta Cryst. **21**, 1011 (1966), doi:10.1107/S0365110X66004419.

Geometry files:

- CIF: pp. 1734
- POSCAR: pp. 1734

Jacutingaite (Pt_2HgSe_3) Structure: AB2C3_hP12_164_d_ae_i

http://aflow.org/prototype-encyclopedia/AB2C3_hP12_164_d_ae_i

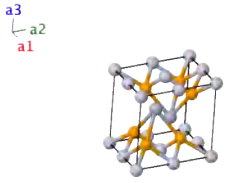


Prototype	:	$HgPt_2Se_3$
AFLOW prototype label	:	AB2C3_hP12_164_d_ae_i
***Strukturbericht* designation**	:	None
Pearson symbol	:	hP12
Space group number	:	164
Space group symbol	:	$P\bar{3}m1$
AFLOW prototype command	:	`aflow --proto=AB2C3_hP12_164_d_ae_i --params=`$a, c/a, z_2, x_4, z_4$

- Table 4 of (Vymazalová, 2012) has a small error: the positions of the Hg(1) atoms should be written $(1/3, 2/3, 0.3507)$ rather than $(1/3, 1/3, 0.3507)$.

Trigonal Hexagonal primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(1*a*)	Pt I
$\mathbf{B}_2$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(2*d*)	Hg
$\mathbf{B}_3$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	(2*d*)	Hg
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{4}a\,\hat{\mathbf{y}}$	(3*e*)	Pt II
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{4}a\,\hat{\mathbf{y}}$	(3*e*)	Pt II
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	(3*e*)	Pt II
$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$-\sqrt{3}x_4 a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(6*i*)	Se
$\mathbf{B}_8$	=	$x_4\,\mathbf{a}_1 + 2x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\frac{3}{2}x_4 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(6*i*)	Se
$\mathbf{B}_9$	=	$-2x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$-\frac{3}{2}x_4 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(6*i*)	Se
$\mathbf{B}_{10}$	=	$-x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\sqrt{3}x_4 a\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(6*i*)	Se
$\mathbf{B}_{11}$	=	$2x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\frac{3}{2}x_4 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(6*i*)	Se
$\mathbf{B}_{12}$	=	$-x_4\,\mathbf{a}_1 - 2x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-\frac{3}{2}x_4 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(6*i*)	Se

References:

- A. Vymazalová, F. Laufek, M. Drábek, A. R. Cabral, J. Haloda, T. Sidorinová, B. Lehmann, H. F. Galbiatti, and J. Drahokoupil, *Jacutingaite, Pt_2HgSe_3, A New Platinum-Group Mineral Species From the Cauê Iron-Ore Deposit, Itabira District, Minas Gerais, Brazil*, Can. Mineral. **50**, 431–440 (2012), doi:10.3749/canmin.50.2.431.

Found in:

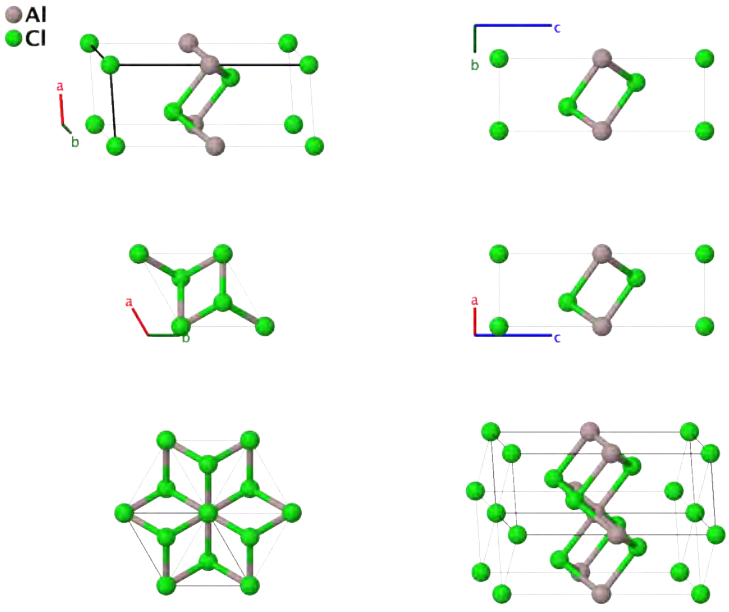
- A. Marrazzo, N. Marzari, and M. Gibertini, *Emergent dual topology in the three-dimensional Kane-Mele Pt_2HgSe_3*, Phys. Rev. Research **2**, 012063 (2020), doi:10.1103/PhysRevResearch.2.012063.

Geometry files:

- CIF: pp. 1734
- POSCAR: pp. 1734

$D0_{13}$ ($AlCl_3$) (*obsolete*) Structure: AB3_hP4_164_b_ad

http://aflow.org/prototype-encyclopedia/AB3_hP4_164_b_ad

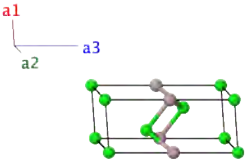


Prototype	:	$AlCl_3$
AFLOW prototype label	:	AB3_hP4_164_b_ad
***Strukturbericht* designation**	:	$D0_{13}$
Pearson symbol	:	hP4
Space group number	:	164
Space group symbol	:	$P\bar{3}m1$
AFLOW prototype command	:	`aflow --proto=AB3_hP4_164_b_ad --params=`$a, c/a, z_3$

- This structure was suggested by (Laschkarew, 1930) and designated as *Strukturbericht* $D0_{13}$ by (Hermann, 1937). However, in the next edition of *Strukturbericht*, (Gottfried, 1937) designated the $AlCl_3$ structure found by (Ketellar, 1935) as $D0_{15}$. This structure, however, has a lower symmetry than the currently accepted structure, which we previously designated $D0_{15}$ and is body-centered orthorhombic, space group $C2/m$ #12.

Trigonal Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(1*a*)	Cl I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	(1*b*)	Al
$\mathbf{B}_3$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(2*d*)	Cl II
$\mathbf{B}_4$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	(2*d*)	Cl II

References:

- W. E. Laschkarew, *Zur Struktur $AlCl_3$*, Z. Anorg. Allg. Chem. **193**, 270–276 (1930), doi:10.1002/zaac.19301930123.
- C. Gottfried and F. Schossberger, eds., *Strukturbericht Band III 1933-1935* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).
- J. A. A. Ketelaar, *Die Kristallstruktur der Aluminiumhalogenide II. Die Kristallstruktur von $AlCl_3$*, Zeitschrift für Kristallographie - Crystalline Materials **90**, 237–255 (1935), doi:10.1524/zkri.1935.90.1.237.

Found in:

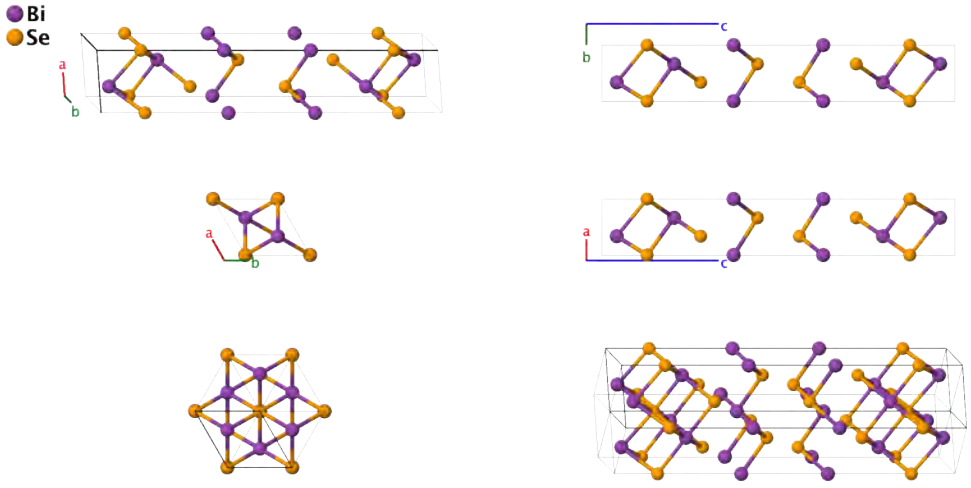
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1735
- POSCAR: pp. 1735

Nevskite (BiSe) Structure: AB_hP12_164_c2d_c2d

http://aflow.org/prototype-encyclopedia/AB_hP12_164_c2d_c2d



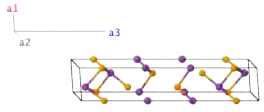
Prototype	:	BiSe
AFLOW prototype label	:	AB_hP12_164_c2d_c2d
***Strukturbericht* designation**	:	None
Pearson symbol	:	hP12
Space group number	:	164
Space group symbol	:	$P\bar{3}m1$
AFLOW prototype command	:	`aflow --proto=AB_hP12_164_c2d_c2d --params=`$a, c/a, z_1, z_2, z_3, z_4, z_5, z_6$

Other compounds with this structure

- BiTe (tsumoite) and $Bi(S_{0.56}Te_{0.44})$ (ingodite)

Trigonal Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$z_1\,\mathbf{a}_3$	=	$z_1 c\,\hat{\mathbf{z}}$	$(2c)$	Bi I
$\mathbf{B}_2$	=	$-z_1\,\mathbf{a}_3$	=	$-z_1 c\,\hat{\mathbf{z}}$	$(2c)$	Bi I
$\mathbf{B}_3$	=	$z_2\,\mathbf{a}_3$	=	$z_2 c\,\hat{\mathbf{z}}$	$(2c)$	Se I
$\mathbf{B}_4$	=	$-z_2\,\mathbf{a}_3$	=	$-z_2 c\,\hat{\mathbf{z}}$	$(2c)$	Se I
$\mathbf{B}_5$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(2d)$	Bi II
$\mathbf{B}_6$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(2d)$	Bi II

$\mathbf{B}_7$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(2*d*)	Bi III
$\mathbf{B}_8$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	(2*d*)	Bi III
$\mathbf{B}_9$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(2*d*)	Se II
$\mathbf{B}_{10}$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(2*d*)	Se II
$\mathbf{B}_{11}$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(2*d*)	Se III
$\mathbf{B}_{12}$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(2*d*)	Se III

References:

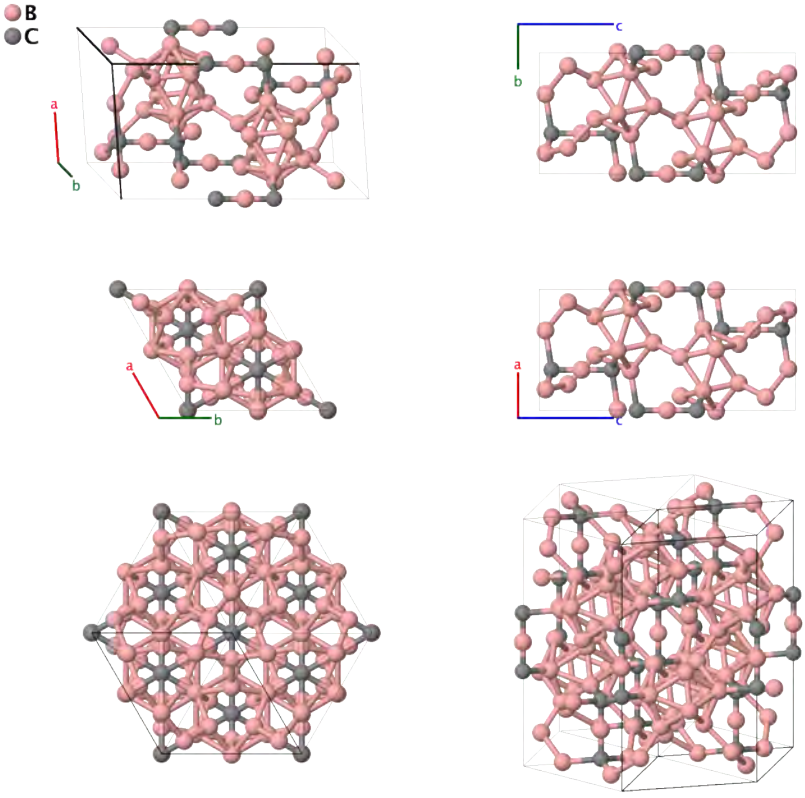
- E. Gaudin, S. Jobic, M. Evain, R. Brec, and J. Rouxel, *Charge balance in some Bi_xSe_y phases through atomic structure determination and band structure calculations*, Mater. Res. Bull. **30**, 549–561 (1995), doi:10.1016/0025-5408(95)00030-5.

Geometry files:

- CIF: pp. 1735
- POSCAR: pp. 1735

$B_{13}C_2$ "B_4C" ($D1_g$) Structure: A13B2_hR15_166_b2h_c

http://aflow.org/prototype-encyclopedia/A13B2_hR15_166_b2h_c



Prototype	:	$B_{13}C_2$
AFLOW prototype label	:	A13B2_hR15_166_b2h_c
***Strukturbericht* designation**	:	$D1_g$
Pearson symbol	:	hR15
Space group number	:	166
Space group symbol	:	$R\bar{3}m$
AFLOW prototype command	:	`aflow --proto=A13B2_hR15_166_b2h_c [--hex] --params=`$a, c/a, x_2, x_3, z_3, x_4, z_4$

Other compounds with this structure

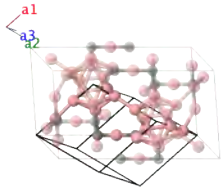
- $B_{1-x}C_x$ ($0.08 < x < 0.20$) and $B_{13}P_2$, B_4Si, B_6O

- This structure has a rather complicated history:
  - It is difficult to determine the species of atoms at a given site because of the similar electronic and nuclear cross sections of ^{11}B and ^{12}C (Domnich, 2011).
  - Early investigations (Clark, 1943) assumed the structure was B_4C, with the extra carbon atom replacing the boron on the (1*b*) site. [Note that Clark has an error in the coordinates of one set of boron atoms, giving a boron-boron distance of less than 1 Å. This error is repeated in (Brandes, 1992) and (Wykcoff, 1964).]

- In reality, concentrations can range from 8-20% carbon (Domnich, 2011).
- (Larson, 1986) states that the (1*b*) site in $B_{13}C_2$ is boron, and as the structure becomes more carbon-rich the carbon atoms replace boron in the icosahedra. We follow this and use the structure determined by (Will, 1976) as our reference.
- (Lazzari, 1999) states that excess electrons go on the "polar" sites of the icosahedron, *i.e.* the sites closest to the carbon atoms on the chains (the B-III atoms in our notation).

Rhombohedral primitive vectors:

$$\begin{aligned}
\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2\sqrt{3}}\,a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,c\,\hat{\mathbf{z}} \\
\mathbf{a}_2 &= \tfrac{1}{\sqrt{3}}\,a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,c\,\hat{\mathbf{z}} \\
\mathbf{a}_3 &= -\tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2\sqrt{3}}\,a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,c\,\hat{\mathbf{z}}
\end{aligned}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$\tfrac{1}{2}\,\mathbf{a}_1 + \tfrac{1}{2}\,\mathbf{a}_2 + \tfrac{1}{2}\,\mathbf{a}_3$	$\tfrac{1}{2}c\,\hat{\mathbf{z}}$	(1*b*)	B I
$\mathbf{B}_2$	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	$x_2 c\,\hat{\mathbf{z}}$	(2*c*)	C
$\mathbf{B}_3$	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 - x_2\,\mathbf{a}_3$	$-x_2 c\,\hat{\mathbf{z}}$	(2*c*)	C
$\mathbf{B}_4$	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$\tfrac{1}{2}\,(x_3 - z_3)\,a\,\hat{\mathbf{x}} + \tfrac{1}{2\sqrt{3}}\,(x_3 - z_3)\,a\,\hat{\mathbf{y}} + \left(\tfrac{2}{3}x_3 + \tfrac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	(6*h*)	B II
$\mathbf{B}_5$	$z_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	$\tfrac{1}{2}\,(-x_3 + z_3)\,a\,\hat{\mathbf{x}} + \tfrac{1}{2\sqrt{3}}\,(x_3 - z_3)\,a\,\hat{\mathbf{y}} + \left(\tfrac{2}{3}x_3 + \tfrac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	(6*h*)	B II
$\mathbf{B}_6$	$x_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	$\tfrac{1}{\sqrt{3}}\,(-x_3 + z_3)\,a\,\hat{\mathbf{y}} + \left(\tfrac{2}{3}x_3 + \tfrac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	(6*h*)	B II
$\mathbf{B}_7$	$-z_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	$\tfrac{1}{2}\,(x_3 - z_3)\,a\,\hat{\mathbf{x}} + \tfrac{1}{2\sqrt{3}}\,(-x_3 + z_3)\,a\,\hat{\mathbf{y}} - \left(\tfrac{2}{3}x_3 + \tfrac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	(6*h*)	B II
$\mathbf{B}_8$	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	$\tfrac{1}{2}\,(-x_3 + z_3)\,a\,\hat{\mathbf{x}} + \tfrac{1}{2\sqrt{3}}\,(-x_3 + z_3)\,a\,\hat{\mathbf{y}} - \left(\tfrac{2}{3}x_3 + \tfrac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	(6*h*)	B II
$\mathbf{B}_9$	$-x_3\,\mathbf{a}_1 - z_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	$\tfrac{1}{\sqrt{3}}\,(x_3 - z_3)\,a\,\hat{\mathbf{y}} - \left(\tfrac{2}{3}x_3 + \tfrac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	(6*h*)	B II
$\mathbf{B}_{10}$	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$\tfrac{1}{2}\,(x_4 - z_4)\,a\,\hat{\mathbf{x}} + \tfrac{1}{2\sqrt{3}}\,(x_4 - z_4)\,a\,\hat{\mathbf{y}} + \left(\tfrac{2}{3}x_4 + \tfrac{1}{3}z_4\right)c\,\hat{\mathbf{z}}$	(6*h*)	B III
$\mathbf{B}_{11}$	$z_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	$\tfrac{1}{2}\,(-x_4 + z_4)\,a\,\hat{\mathbf{x}} + \tfrac{1}{2\sqrt{3}}\,(x_4 - z_4)\,a\,\hat{\mathbf{y}} + \left(\tfrac{2}{3}x_4 + \tfrac{1}{3}z_4\right)c\,\hat{\mathbf{z}}$	(6*h*)	B III
$\mathbf{B}_{12}$	$x_4\,\mathbf{a}_1 + z_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	$\tfrac{1}{\sqrt{3}}\,(-x_4 + z_4)\,a\,\hat{\mathbf{y}} + \left(\tfrac{2}{3}x_4 + \tfrac{1}{3}z_4\right)c\,\hat{\mathbf{z}}$	(6*h*)	B III
$\mathbf{B}_{13}$	$-z_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	$\tfrac{1}{2}\,(x_4 - z_4)\,a\,\hat{\mathbf{x}} + \tfrac{1}{2\sqrt{3}}\,(-x_4 + z_4)\,a\,\hat{\mathbf{y}} - \left(\tfrac{2}{3}x_4 + \tfrac{1}{3}z_4\right)c\,\hat{\mathbf{z}}$	(6*h*)	B III
$\mathbf{B}_{14}$	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	$\tfrac{1}{2}\,(-x_4 + z_4)\,a\,\hat{\mathbf{x}} + \tfrac{1}{2\sqrt{3}}\,(-x_4 + z_4)\,a\,\hat{\mathbf{y}} - \left(\tfrac{2}{3}x_4 + \tfrac{1}{3}z_4\right)c\,\hat{\mathbf{z}}$	(6*h*)	B III
$\mathbf{B}_{15}$	$-x_4\,\mathbf{a}_1 - z_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	$\tfrac{1}{\sqrt{3}}\,(x_4 - z_4)\,a\,\hat{\mathbf{y}} - \left(\tfrac{2}{3}x_4 + \tfrac{1}{3}z_4\right)c\,\hat{\mathbf{z}}$	(6*h*)	B III

References:

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- R. W. G. Wyckoff, *Crystal Structures*, vol. 2 (Interscience (John Wiley & Sons), New York, London, Sydney, 1964), second edn.
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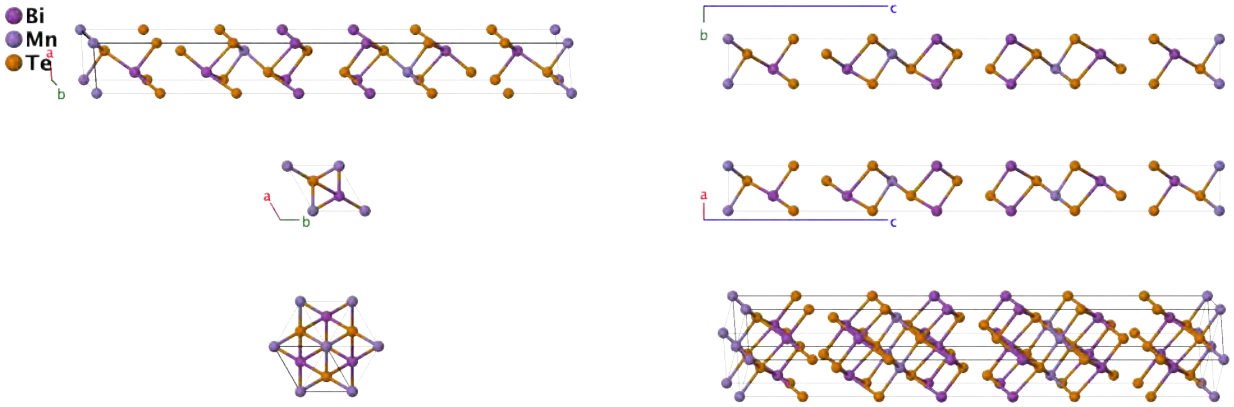
- R. Lazzari, N. Vast, J. M. Besson, S. Baroni, and A. Dal Corso, *Atomic Structure and Vibrational Properties of Icosahedral B_4C Boron Carbide*, Phys. Rev. Lett. **83**, 3230–3233 (1999), doi:10.1103/PhysRevLett.83.3230. *erratum* Phys. Rev. Lett. 85, 4194 (2000).

Geometry files:
- CIF: pp. 1736
- POSCAR: pp. 1736

$MnBi_2Te_4$ Structure: A2BC4_hR7_166_c_a_2c

http://aflow.org/prototype-encyclopedia/A2BC4_hR7_166_c_a_2c

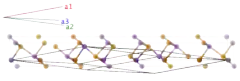


Prototype	:	Bi_2MnTe_4
AFLOW prototype label	:	A2BC4_hR7_166_c_a_2c
***Strukturbericht* designation**	:	None
Pearson symbol	:	hR7
Space group number	:	166
Space group symbol	:	$R\bar{3}m$
AFLOW prototype command	:	`aflow --proto=A2BC4_hR7_166_c_a_2c [--hex] --params=`$a, c/a, x_2, x_3, x_4$

- We use the data taken at 10 K. Except for the magnetic ordering there is no substantial change in the structure up to room temperature.

Rhombohedral primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = -\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(1a)$	Mn
$\mathbf{B}_2$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$x_2 c\,\hat{\mathbf{z}}$	$(2c)$	Bi
$\mathbf{B}_3$	=	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 - x_2\,\mathbf{a}_3$	=	$-x_2 c\,\hat{\mathbf{z}}$	$(2c)$	Bi
$\mathbf{B}_4$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3 c\,\hat{\mathbf{z}}$	$(2c)$	Te I
$\mathbf{B}_5$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$-x_3 c\,\hat{\mathbf{z}}$	$(2c)$	Te I
$\mathbf{B}_6$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4 c\,\hat{\mathbf{z}}$	$(2c)$	Te II
$\mathbf{B}_7$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$-x_4 c\,\hat{\mathbf{z}}$	$(2c)$	Te II

References:

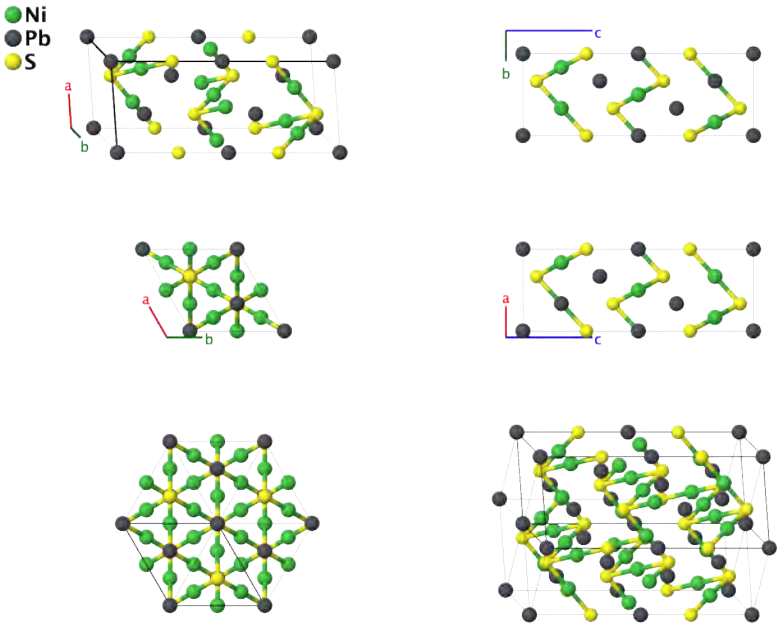
- J.-Q. Yan, Q. Zhang, T. Heitmann, Z. Huang, K. Y. Chen, J.-G. Cheng, W. Wu, D. Vaknin, B. C. Sales, and R. J. McQueeney, *Crystal growth and magnetic structure of* $MnBi_2Te_4$, Phys. Rev. Mater. **3**, 064202 (2019), doi:10.1103/PhysRevMaterials.3.064202.

Geometry files:

- CIF: pp. 1736
- POSCAR: pp. 1737

Shandite ($Ni_3Pb_2S_2$) Structure: A3B2C2_hR7_166_d_ab_c

http://aflow.org/prototype-encyclopedia/A3B2C2_hR7_166_d_ab_c



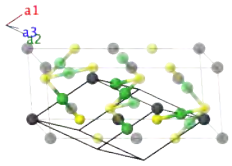
Prototype	:	$Ni_3Pb_2S_2$
AFLOW prototype label	:	A3B2C2_hR7_166_d_ab_c
***Strukturbericht* designation**	:	None
Pearson symbol	:	hR7
Space group number	:	166
Space group symbol	:	$R\bar{3}m$
AFLOW prototype command	:	`aflow --proto=A3B2C2_hR7_166_d_ab_c [--hex] --params=`$a, c/a, x_3$

Other compounds with this structure

- $Co_3In_2S_2$, $Co_3Sn_2S_2$, $Ni_3In_2S_2$, $Ni_3Sn_2S_2$, and $O_3K_2Sn_2$

Rhombohedral primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = -\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$ =	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($1a$)	Pb I

$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	(1*b*)	Pb II
$\mathbf{B}_3$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3 c\,\hat{\mathbf{z}}$	(2*c*)	S
$\mathbf{B}_4$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$-x_3 c\,\hat{\mathbf{z}}$	(2*c*)	S
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} - \frac{1}{4\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{6}c\,\hat{\mathbf{z}}$	(3*d*)	Ni
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{6}c\,\hat{\mathbf{z}}$	(3*d*)	Ni
$\mathbf{B}_7$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$-\frac{1}{4}a\,\hat{\mathbf{x}} - \frac{1}{4\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{6}c\,\hat{\mathbf{z}}$	(3*d*)	Ni

References:

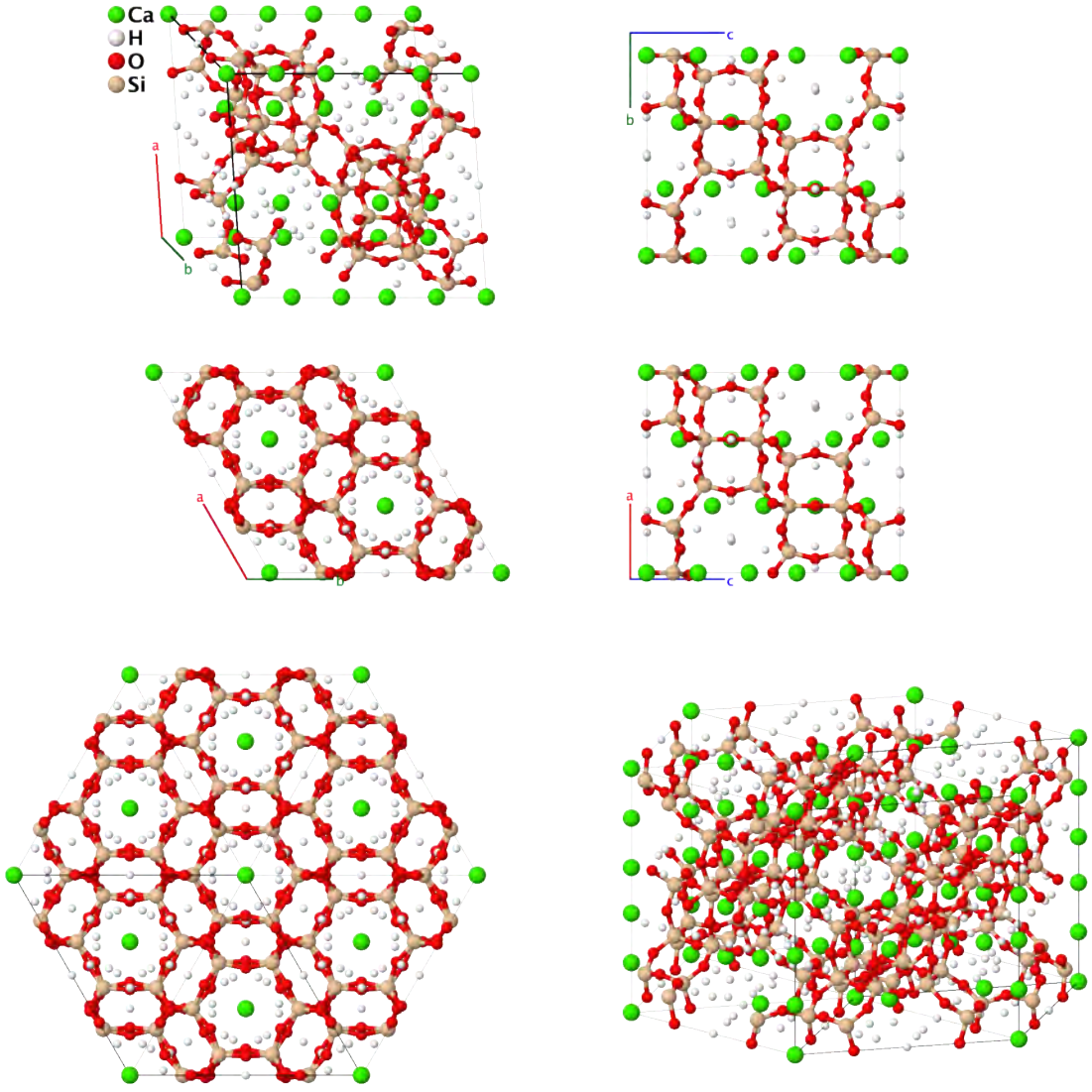
- R. Weihrich, S. F. Matar, V. Eyert, F. Rau, M. Zabel, M. Andratschke, I. Anusca, and T. Bernert, *Structure, ordering, and bonding of half antiperovskites: $PbNi_{3/2}S$ and $BiPd_{3/2}S$*, Prog. Solid State Chem. **35**, 309–327 (2007), doi:10.1016/j.progsolidstchem.2007.01.011.

Geometry files:

- CIF: pp. 1737
- POSCAR: pp. 1737

Chabazite ($Ca_{1.4}Sr_{0.3}Al_{3.8}Si_{8.3}O_{24}\cdot 13H_2O$, $S3_4$ (I)) Structure: A5B21C24D12_hR62_166_a2c_ehi_fg2h_i

http://aflow.org/prototype-encyclopedia/A5B21C24D12_hR62_166_a2c_ehi_fg2h_i



Prototype	:	$(Ca_{1.4},Sr_{0.3})(H_2O)_{13}O_{24}(Si_{8.3},Al_{3.8})$
AFLOW prototype label	:	A5B21C24D12_hR62_166_a2c_ehi_fg2h_i
***Strukturbericht* designation**	:	$S3_4$ (I)
Pearson symbol	:	hR62
Space group number	:	166
Space group symbol	:	$R\bar{3}m$
AFLOW prototype command	:	`aflow --proto=A5B21C24D12_hR62_166_a2c_ehi_fg2h_i [--hex] --params=`$a, c/a, x_2, x_3, x_5, x_6, x_7, z_7, x_8, z_8, x_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}$

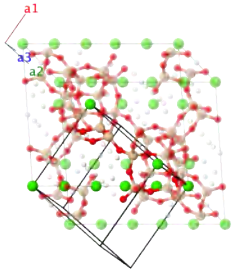
- The structure given here, from (Calligaris, 1982), has substantially the same Wyckoff positions as the structure desig-

nated $S\,3_4$ by (Gottfried, 1937). The largest change is the splitting of the calcium site from one stoichiometric ($2c$) site to three partially filled sites, one ($1a$) and two ($2c$).

- (Calligaris, 1982) did not give the exact positions of the water molecules, so we use the positions given by (Downs, 2003). However, this leaves us with only 7.7 water molecules per formula unit, rather than the 13 claimed by (Calligaris, 1982).
- Only the oxygen sites are fully occupied. For the remaining sites, according to (Downs, 2003),
  - The ($1a$) site is 9.06% Ca and 1.94% Sr, with the remaining sites vacant.
  - The first ($2c$) site is 43.65% Ca and 9.35% Sr.
  - The second ($2c$) site is 19.76% Ca and 4.24% Sr.
  - The ($3e$) is only occupied by a water molecule 50% of the time.
  - The ($6h$) H_2O site is only occupied 57% of the time.
  - The ($12i$) H_2O site is only occupied 23% of the time.
  - The remaining ($12i$) is 69% Si and 31% Al.
- (Gottfried, 1937) gave chabazite the *Strukturbericht* designation $S\,3_4$. However, (Gottfried, 1940) ignored this and designated catapleiite, $Na_2ZrSi_3O_9{\cdot}2H_2O$ as *Strukturbericht* $S\,3_4$. We resolve this by using $S\,3_4$ (I) to designate chabazite and $S\,3_4$ (II) to designate catapleiite.

Rhombohedral primitive vectors:

$$\begin{aligned}
\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2\sqrt{3}}\,a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,c\,\hat{\mathbf{z}} \\
\mathbf{a}_2 &= \tfrac{1}{\sqrt{3}}\,a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,c\,\hat{\mathbf{z}} \\
\mathbf{a}_3 &= -\tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2\sqrt{3}}\,a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,c\,\hat{\mathbf{z}}
\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($1a$)	Ca I
$\mathbf{B}_2$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$x_2 c\,\hat{\mathbf{z}}$	($2c$)	Ca II
$\mathbf{B}_3$	=	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 - x_2\,\mathbf{a}_3$	=	$-x_2 c\,\hat{\mathbf{z}}$	($2c$)	Ca II
$\mathbf{B}_4$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3 c\,\hat{\mathbf{z}}$	($2c$)	Ca III
$\mathbf{B}_5$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$-x_3 c\,\hat{\mathbf{z}}$	($2c$)	Ca III
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$	($3e$)	H_2O I
$\mathbf{B}_7$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$-\frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$	($3e$)	H_2O I
$\mathbf{B}_8$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$	($3e$)	H_2O I
$\mathbf{B}_9$	=	$x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2$	=	$\frac{1}{2}x_5 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}}$	($6f$)	O I
$\mathbf{B}_{10}$	=	$x_5\,\mathbf{a}_2 - x_5\,\mathbf{a}_3$	=	$\frac{1}{2}x_5 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}}$	($6f$)	O I
$\mathbf{B}_{11}$	=	$-x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}}$	($6f$)	O I
$\mathbf{B}_{12}$	=	$-x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2$	=	$-\frac{1}{2}x_5 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}}$	($6f$)	O I
$\mathbf{B}_{13}$	=	$-x_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$-\frac{1}{2}x_5 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}}$	($6f$)	O I
$\mathbf{B}_{14}$	=	$x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}}$	($6f$)	O I

$$\mathbf{B}_{15} = x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 + \tfrac{1}{2}\,\mathbf{a}_3 = \left(-\tfrac{1}{4} + \tfrac{1}{2}x_6\right) a\,\hat{\mathbf{x}} - \left(\tfrac{\sqrt{3}}{2}x_6 + \tfrac{1}{4\sqrt{3}}\right) a\,\hat{\mathbf{y}} + \tfrac{1}{6} c\,\hat{\mathbf{z}} \quad (6g) \quad \text{O II}$$

$$\mathbf{B}_{16} = \tfrac{1}{2}\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 - x_6\,\mathbf{a}_3 = \left(\tfrac{1}{4} + \tfrac{1}{2}x_6\right) a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{4\sqrt{3}} + \tfrac{\sqrt{3}}{2}x_6\right) a\,\hat{\mathbf{y}} + \tfrac{1}{6} c\,\hat{\mathbf{z}} \quad (6g) \quad \text{O II}$$

$$\mathbf{B}_{17} = -x_6\,\mathbf{a}_1 + \tfrac{1}{2}\,\mathbf{a}_2 + x_6\,\mathbf{a}_3 = -x_6 a\,\hat{\mathbf{x}} + \tfrac{1}{2\sqrt{3}} a\,\hat{\mathbf{y}} + \tfrac{1}{6} c\,\hat{\mathbf{z}} \quad (6g) \quad \text{O II}$$

$$\mathbf{B}_{18} = -x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + \tfrac{1}{2}\,\mathbf{a}_3 = -\left(\tfrac{1}{2}x_6 + \tfrac{1}{4}\right) a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{4\sqrt{3}} + \tfrac{\sqrt{3}}{2}x_6\right) a\,\hat{\mathbf{y}} + \tfrac{1}{6} c\,\hat{\mathbf{z}} \quad (6g) \quad \text{O II}$$

$$\mathbf{B}_{19} = \tfrac{1}{2}\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 + x_6\,\mathbf{a}_3 = \left(\tfrac{1}{4} - \tfrac{1}{2}x_6\right) a\,\hat{\mathbf{x}} - \left(\tfrac{\sqrt{3}}{2}x_6 + \tfrac{1}{4\sqrt{3}}\right) a\,\hat{\mathbf{y}} + \tfrac{1}{6} c\,\hat{\mathbf{z}} \quad (6g) \quad \text{O II}$$

$$\mathbf{B}_{20} = x_6\,\mathbf{a}_1 + \tfrac{1}{2}\,\mathbf{a}_2 - x_6\,\mathbf{a}_3 = x_6 a\,\hat{\mathbf{x}} + \tfrac{1}{2\sqrt{3}} a\,\hat{\mathbf{y}} + \tfrac{1}{6} c\,\hat{\mathbf{z}} \quad (6g) \quad \text{O II}$$

$$\mathbf{B}_{21} = x_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 = \tfrac{1}{2}(x_7 - z_7)\, a\,\hat{\mathbf{x}} + \tfrac{1}{2\sqrt{3}}(x_7 - z_7)\, a\,\hat{\mathbf{y}} + \left(\tfrac{2}{3}x_7 + \tfrac{1}{3}z_7\right) c\,\hat{\mathbf{z}} \quad (6h) \quad \text{H}_2\text{O II}$$

$$\mathbf{B}_{22} = z_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + x_7\,\mathbf{a}_3 = \tfrac{1}{2}(-x_7 + z_7)\, a\,\hat{\mathbf{x}} + \tfrac{1}{2\sqrt{3}}(x_7 - z_7)\, a\,\hat{\mathbf{y}} + \left(\tfrac{2}{3}x_7 + \tfrac{1}{3}z_7\right) c\,\hat{\mathbf{z}} \quad (6h) \quad \text{H}_2\text{O II}$$

$$\mathbf{B}_{23} = x_7\,\mathbf{a}_1 + z_7\,\mathbf{a}_2 + x_7\,\mathbf{a}_3 = \tfrac{1}{\sqrt{3}}(-x_7 + z_7)\, a\,\hat{\mathbf{y}} + \left(\tfrac{2}{3}x_7 + \tfrac{1}{3}z_7\right) c\,\hat{\mathbf{z}} \quad (6h) \quad \text{H}_2\text{O II}$$

$$\mathbf{B}_{24} = -z_7\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 - x_7\,\mathbf{a}_3 = \tfrac{1}{2}(x_7 - z_7)\, a\,\hat{\mathbf{x}} + \tfrac{1}{2\sqrt{3}}(-x_7 + z_7)\, a\,\hat{\mathbf{y}} - \left(\tfrac{2}{3}x_7 + \tfrac{1}{3}z_7\right) c\,\hat{\mathbf{z}} \quad (6h) \quad \text{H}_2\text{O II}$$

$$\mathbf{B}_{25} = -x_7\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3 = \tfrac{1}{2}(-x_7 + z_7)\, a\,\hat{\mathbf{x}} + \tfrac{1}{2\sqrt{3}}(-x_7 + z_7)\, a\,\hat{\mathbf{y}} - \left(\tfrac{2}{3}x_7 + \tfrac{1}{3}z_7\right) c\,\hat{\mathbf{z}} \quad (6h) \quad \text{H}_2\text{O II}$$

$$\mathbf{B}_{26} = -x_7\,\mathbf{a}_1 - z_7\,\mathbf{a}_2 - x_7\,\mathbf{a}_3 = \tfrac{1}{\sqrt{3}}(x_7 - z_7)\, a\,\hat{\mathbf{y}} - \left(\tfrac{2}{3}x_7 + \tfrac{1}{3}z_7\right) c\,\hat{\mathbf{z}} \quad (6h) \quad \text{H}_2\text{O II}$$

$$\mathbf{B}_{27} = x_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 = \tfrac{1}{2}(x_8 - z_8)\, a\,\hat{\mathbf{x}} + \tfrac{1}{2\sqrt{3}}(x_8 - z_8)\, a\,\hat{\mathbf{y}} + \left(\tfrac{2}{3}x_8 + \tfrac{1}{3}z_8\right) c\,\hat{\mathbf{z}} \quad (6h) \quad \text{O III}$$

$$\mathbf{B}_{28} = z_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 + x_8\,\mathbf{a}_3 = \tfrac{1}{2}(-x_8 + z_8)\, a\,\hat{\mathbf{x}} + \tfrac{1}{2\sqrt{3}}(x_8 - z_8)\, a\,\hat{\mathbf{y}} + \left(\tfrac{2}{3}x_8 + \tfrac{1}{3}z_8\right) c\,\hat{\mathbf{z}} \quad (6h) \quad \text{O III}$$

$$\mathbf{B}_{29} = x_8\,\mathbf{a}_1 + z_8\,\mathbf{a}_2 + x_8\,\mathbf{a}_3 = \tfrac{1}{\sqrt{3}}(-x_8 + z_8)\, a\,\hat{\mathbf{y}} + \left(\tfrac{2}{3}x_8 + \tfrac{1}{3}z_8\right) c\,\hat{\mathbf{z}} \quad (6h) \quad \text{O III}$$

$$\mathbf{B}_{30} = -z_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 - x_8\,\mathbf{a}_3 = \tfrac{1}{2}(x_8 - z_8)\, a\,\hat{\mathbf{x}} + \tfrac{1}{2\sqrt{3}}(-x_8 + z_8)\, a\,\hat{\mathbf{y}} - \left(\tfrac{2}{3}x_8 + \tfrac{1}{3}z_8\right) c\,\hat{\mathbf{z}} \quad (6h) \quad \text{O III}$$

$$\mathbf{B}_{31} = -x_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3 = \tfrac{1}{2}(-x_8 + z_8)\, a\,\hat{\mathbf{x}} + \tfrac{1}{2\sqrt{3}}(-x_8 + z_8)\, a\,\hat{\mathbf{y}} - \left(\tfrac{2}{3}x_8 + \tfrac{1}{3}z_8\right) c\,\hat{\mathbf{z}} \quad (6h) \quad \text{O III}$$

$$\mathbf{B}_{32} = -x_8\,\mathbf{a}_1 - z_8\,\mathbf{a}_2 - x_8\,\mathbf{a}_3 = \tfrac{1}{\sqrt{3}}(x_8 - z_8)\, a\,\hat{\mathbf{y}} - \left(\tfrac{2}{3}x_8 + \tfrac{1}{3}z_8\right) c\,\hat{\mathbf{z}} \quad (6h) \quad \text{O III}$$

$$\mathbf{B}_{33} = x_9\,\mathbf{a}_1 + x_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 = \tfrac{1}{2}(x_9 - z_9)\, a\,\hat{\mathbf{x}} + \tfrac{1}{2\sqrt{3}}(x_9 - z_9)\, a\,\hat{\mathbf{y}} + \left(\tfrac{2}{3}x_9 + \tfrac{1}{3}z_9\right) c\,\hat{\mathbf{z}} \quad (6h) \quad \text{O IV}$$

$$\mathbf{B}_{34} = z_9\,\mathbf{a}_1 + x_9\,\mathbf{a}_2 + x_9\,\mathbf{a}_3 = \tfrac{1}{2}(-x_9 + z_9)\, a\,\hat{\mathbf{x}} + \tfrac{1}{2\sqrt{3}}(x_9 - z_9)\, a\,\hat{\mathbf{y}} + \left(\tfrac{2}{3}x_9 + \tfrac{1}{3}z_9\right) c\,\hat{\mathbf{z}} \quad (6h) \quad \text{O IV}$$

$$\mathbf{B}_{35} = x_9\,\mathbf{a}_1 + z_9\,\mathbf{a}_2 + x_9\,\mathbf{a}_3 = \tfrac{1}{\sqrt{3}}(-x_9 + z_9)\, a\,\hat{\mathbf{y}} + \left(\tfrac{2}{3}x_9 + \tfrac{1}{3}z_9\right) c\,\hat{\mathbf{z}} \quad (6h) \quad \text{O IV}$$

$$\mathbf{B}_{36} = -z_9\,\mathbf{a}_1 - x_9\,\mathbf{a}_2 - x_9\,\mathbf{a}_3 = \tfrac{1}{2}(x_9 - z_9)\, a\,\hat{\mathbf{x}} + \tfrac{1}{2\sqrt{3}}(-x_9 + z_9)\, a\,\hat{\mathbf{y}} - \left(\tfrac{2}{3}x_9 + \tfrac{1}{3}z_9\right) c\,\hat{\mathbf{z}} \quad (6h) \quad \text{O IV}$$

$$\mathbf{B}_{37} = -x_9\,\mathbf{a}_1 - x_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3 = \tfrac{1}{2}(-x_9 + z_9)\, a\,\hat{\mathbf{x}} + \tfrac{1}{2\sqrt{3}}(-x_9 + z_9)\, a\,\hat{\mathbf{y}} - \left(\tfrac{2}{3}x_9 + \tfrac{1}{3}z_9\right) c\,\hat{\mathbf{z}} \quad (6h) \quad \text{O IV}$$

$$\mathbf{B}_{38} = -x_9\,\mathbf{a}_1 - z_9\,\mathbf{a}_2 - x_9\,\mathbf{a}_3 = \tfrac{1}{\sqrt{3}}(x_9 - z_9)\, a\,\hat{\mathbf{y}} - \left(\tfrac{2}{3}x_9 + \tfrac{1}{3}z_9\right) c\,\hat{\mathbf{z}} \quad (6h) \quad \text{O IV}$$

$$\mathbf{B}_{39} = x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 = \frac{1}{2}\,(x_{10} - z_{10})\,a\,\hat{\mathbf{x}} + \left(-\frac{1}{2\sqrt{3}}x_{10} + \frac{1}{\sqrt{3}}y_{10} - \frac{1}{2\sqrt{3}}z_{10}\right)a\,\hat{\mathbf{y}} + \frac{1}{3}\,(x_{10} + y_{10} + z_{10})\,c\,\hat{\mathbf{z}} \quad (12i) \quad \text{H}_2\text{O III}$$

$$\mathbf{B}_{40} = z_{10}\,\mathbf{a}_1 + x_{10}\,\mathbf{a}_2 + y_{10}\,\mathbf{a}_3 = \frac{1}{2}\,(-y_{10} + z_{10})\,a\,\hat{\mathbf{x}} + \left(\frac{1}{\sqrt{3}}x_{10} - \frac{1}{2\sqrt{3}}y_{10} - \frac{1}{2\sqrt{3}}z_{10}\right)a\,\hat{\mathbf{y}} + \frac{1}{3}\,(x_{10} + y_{10} + z_{10})\,c\,\hat{\mathbf{z}} \quad (12i) \quad \text{H}_2\text{O III}$$

$$\mathbf{B}_{41} = y_{10}\,\mathbf{a}_1 + z_{10}\,\mathbf{a}_2 + x_{10}\,\mathbf{a}_3 = \frac{1}{2}\,(-x_{10} + y_{10})\,a\,\hat{\mathbf{x}} + \left(-\frac{1}{2\sqrt{3}}x_{10} - \frac{1}{2\sqrt{3}}y_{10} + \frac{1}{\sqrt{3}}z_{10}\right)a\,\hat{\mathbf{y}} + \frac{1}{3}\,(x_{10} + y_{10} + z_{10})\,c\,\hat{\mathbf{z}} \quad (12i) \quad \text{H}_2\text{O III}$$

$$\mathbf{B}_{42} = -z_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 - x_{10}\,\mathbf{a}_3 = \frac{1}{2}\,(x_{10} - z_{10})\,a\,\hat{\mathbf{x}} + \left(\frac{1}{2\sqrt{3}}x_{10} - \frac{1}{\sqrt{3}}y_{10} + \frac{1}{2\sqrt{3}}z_{10}\right)a\,\hat{\mathbf{y}} - \frac{1}{3}\,(x_{10} + y_{10} + z_{10})\,c\,\hat{\mathbf{z}} \quad (12i) \quad \text{H}_2\text{O III}$$

$$\mathbf{B}_{43} = -y_{10}\,\mathbf{a}_1 - x_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 = \frac{1}{2}\,(-y_{10} + z_{10})\,a\,\hat{\mathbf{x}} + \left(-\frac{1}{\sqrt{3}}x_{10} + \frac{1}{2\sqrt{3}}y_{10} + \frac{1}{2\sqrt{3}}z_{10}\right)a\,\hat{\mathbf{y}} - \frac{1}{3}\,(x_{10} + y_{10} + z_{10})\,c\,\hat{\mathbf{z}} \quad (12i) \quad \text{H}_2\text{O III}$$

$$\mathbf{B}_{44} = -x_{10}\,\mathbf{a}_1 - z_{10}\,\mathbf{a}_2 - y_{10}\,\mathbf{a}_3 = \frac{1}{2}\,(-x_{10} + y_{10})\,a\,\hat{\mathbf{x}} + \left(\frac{1}{2\sqrt{3}}x_{10} + \frac{1}{2\sqrt{3}}y_{10} - \frac{1}{\sqrt{3}}z_{10}\right)a\,\hat{\mathbf{y}} - \frac{1}{3}\,(x_{10} + y_{10} + z_{10})\,c\,\hat{\mathbf{z}} \quad (12i) \quad \text{H}_2\text{O III}$$

$$\mathbf{B}_{45} = -x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 = \frac{1}{2}\,(-x_{10} + z_{10})\,a\,\hat{\mathbf{x}} + \left(\frac{1}{2\sqrt{3}}x_{10} - \frac{1}{\sqrt{3}}y_{10} + \frac{1}{2\sqrt{3}}z_{10}\right)a\,\hat{\mathbf{y}} - \frac{1}{3}\,(x_{10} + y_{10} + z_{10})\,c\,\hat{\mathbf{z}} \quad (12i) \quad \text{H}_2\text{O III}$$

$$\mathbf{B}_{46} = -z_{10}\,\mathbf{a}_1 - x_{10}\,\mathbf{a}_2 - y_{10}\,\mathbf{a}_3 = \frac{1}{2}\,(y_{10} - z_{10})\,a\,\hat{\mathbf{x}} + \left(-\frac{1}{\sqrt{3}}x_{10} + \frac{1}{2\sqrt{3}}y_{10} + \frac{1}{2\sqrt{3}}z_{10}\right)a\,\hat{\mathbf{y}} - \frac{1}{3}\,(x_{10} + y_{10} + z_{10})\,c\,\hat{\mathbf{z}} \quad (12i) \quad \text{H}_2\text{O III}$$

$$\mathbf{B}_{47} = -y_{10}\,\mathbf{a}_1 - z_{10}\,\mathbf{a}_2 - x_{10}\,\mathbf{a}_3 = \frac{1}{2}\,(x_{10} - y_{10})\,a\,\hat{\mathbf{x}} + \left(\frac{1}{2\sqrt{3}}x_{10} + \frac{1}{2\sqrt{3}}y_{10} - \frac{1}{\sqrt{3}}z_{10}\right)a\,\hat{\mathbf{y}} - \frac{1}{3}\,(x_{10} + y_{10} + z_{10})\,c\,\hat{\mathbf{z}} \quad (12i) \quad \text{H}_2\text{O III}$$

$$\mathbf{B}_{48} = z_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + x_{10}\,\mathbf{a}_3 = \frac{1}{2}\,(-x_{10} + z_{10})\,a\,\hat{\mathbf{x}} + \left(-\frac{1}{2\sqrt{3}}x_{10} + \frac{1}{\sqrt{3}}y_{10} - \frac{1}{2\sqrt{3}}z_{10}\right)a\,\hat{\mathbf{y}} + \frac{1}{3}\,(x_{10} + y_{10} + z_{10})\,c\,\hat{\mathbf{z}} \quad (12i) \quad \text{H}_2\text{O III}$$

$$\mathbf{B}_{49} = y_{10}\,\mathbf{a}_1 + x_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 = \frac{1}{2}\,(y_{10} - z_{10})\,a\,\hat{\mathbf{x}} + \left(\frac{1}{\sqrt{3}}x_{10} - \frac{1}{2\sqrt{3}}y_{10} - \frac{1}{2\sqrt{3}}z_{10}\right)a\,\hat{\mathbf{y}} + \frac{1}{3}\,(x_{10} + y_{10} + z_{10})\,c\,\hat{\mathbf{z}} \quad (12i) \quad \text{H}_2\text{O III}$$

$$\mathbf{B}_{50} = x_{10}\,\mathbf{a}_1 + z_{10}\,\mathbf{a}_2 + y_{10}\,\mathbf{a}_3 = \frac{1}{2}\,(x_{10} - y_{10})\,a\,\hat{\mathbf{x}} + \left(-\frac{1}{2\sqrt{3}}x_{10} - \frac{1}{2\sqrt{3}}y_{10} + \frac{1}{\sqrt{3}}z_{10}\right)a\,\hat{\mathbf{y}} + \frac{1}{3}\,(x_{10} + y_{10} + z_{10})\,c\,\hat{\mathbf{z}} \quad (12i) \quad \text{H}_2\text{O III}$$

$$\mathbf{B}_{51} = x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 = \frac{1}{2}\,(x_{11} - z_{11})\,a\,\hat{\mathbf{x}} + \left(-\frac{1}{2\sqrt{3}}x_{11} + \frac{1}{\sqrt{3}}y_{11} - \frac{1}{2\sqrt{3}}z_{11}\right)a\,\hat{\mathbf{y}} + \frac{1}{3}\,(x_{11} + y_{11} + z_{11})\,c\,\hat{\mathbf{z}} \quad (12i) \quad \text{Si}$$

$$\mathbf{B}_{52} = z_{11}\,\mathbf{a}_1 + x_{11}\,\mathbf{a}_2 + y_{11}\,\mathbf{a}_3 = \tfrac{1}{2}\,(-y_{11} + z_{11})\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{\sqrt{3}}x_{11} - \tfrac{1}{2\sqrt{3}}y_{11} - \tfrac{1}{2\sqrt{3}}z_{11}\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,(x_{11} + y_{11} + z_{11})\,c\,\hat{\mathbf{z}} \quad (12i) \quad \text{Si}$$

$$\mathbf{B}_{53} = y_{11}\,\mathbf{a}_1 + z_{11}\,\mathbf{a}_2 + x_{11}\,\mathbf{a}_3 = \tfrac{1}{2}\,(-x_{11} + y_{11})\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_{11} - \tfrac{1}{2\sqrt{3}}y_{11} + \tfrac{1}{\sqrt{3}}z_{11}\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,(x_{11} + y_{11} + z_{11})\,c\,\hat{\mathbf{z}} \quad (12i) \quad \text{Si}$$

$$\mathbf{B}_{54} = -z_{11}\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 - x_{11}\,\mathbf{a}_3 = \tfrac{1}{2}\,(x_{11} - z_{11})\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2\sqrt{3}}x_{11} - \tfrac{1}{\sqrt{3}}y_{11} + \tfrac{1}{2\sqrt{3}}z_{11}\right)a\,\hat{\mathbf{y}} - \tfrac{1}{3}\,(x_{11} + y_{11} + z_{11})\,c\,\hat{\mathbf{z}} \quad (12i) \quad \text{Si}$$

$$\mathbf{B}_{55} = -y_{11}\,\mathbf{a}_1 - x_{11}\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3 = \tfrac{1}{2}\,(-y_{11} + z_{11})\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{\sqrt{3}}x_{11} + \tfrac{1}{2\sqrt{3}}y_{11} + \tfrac{1}{2\sqrt{3}}z_{11}\right)a\,\hat{\mathbf{y}} - \tfrac{1}{3}\,(x_{11} + y_{11} + z_{11})\,c\,\hat{\mathbf{z}} \quad (12i) \quad \text{Si}$$

$$\mathbf{B}_{56} = -x_{11}\,\mathbf{a}_1 - z_{11}\,\mathbf{a}_2 - y_{11}\,\mathbf{a}_3 = \tfrac{1}{2}\,(-x_{11} + y_{11})\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2\sqrt{3}}x_{11} + \tfrac{1}{2\sqrt{3}}y_{11} - \tfrac{1}{\sqrt{3}}z_{11}\right)a\,\hat{\mathbf{y}} - \tfrac{1}{3}\,(x_{11} + y_{11} + z_{11})\,c\,\hat{\mathbf{z}} \quad (12i) \quad \text{Si}$$

$$\mathbf{B}_{57} = -x_{11}\,\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3 = \tfrac{1}{2}\,(-x_{11} + z_{11})\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2\sqrt{3}}x_{11} - \tfrac{1}{\sqrt{3}}y_{11} + \tfrac{1}{2\sqrt{3}}z_{11}\right)a\,\hat{\mathbf{y}} - \tfrac{1}{3}\,(x_{11} + y_{11} + z_{11})\,c\,\hat{\mathbf{z}} \quad (12i) \quad \text{Si}$$

$$\mathbf{B}_{58} = -z_{11}\,\mathbf{a}_1 - x_{11}\,\mathbf{a}_2 - y_{11}\,\mathbf{a}_3 = \tfrac{1}{2}\,(y_{11} - z_{11})\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{\sqrt{3}}x_{11} + \tfrac{1}{2\sqrt{3}}y_{11} + \tfrac{1}{2\sqrt{3}}z_{11}\right)a\,\hat{\mathbf{y}} - \tfrac{1}{3}\,(x_{11} + y_{11} + z_{11})\,c\,\hat{\mathbf{z}} \quad (12i) \quad \text{Si}$$

$$\mathbf{B}_{59} = -y_{11}\,\mathbf{a}_1 - z_{11}\,\mathbf{a}_2 - x_{11}\,\mathbf{a}_3 = \tfrac{1}{2}\,(x_{11} - y_{11})\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2\sqrt{3}}x_{11} + \tfrac{1}{2\sqrt{3}}y_{11} - \tfrac{1}{\sqrt{3}}z_{11}\right)a\,\hat{\mathbf{y}} - \tfrac{1}{3}\,(x_{11} + y_{11} + z_{11})\,c\,\hat{\mathbf{z}} \quad (12i) \quad \text{Si}$$

$$\mathbf{B}_{60} = z_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + x_{11}\,\mathbf{a}_3 = \tfrac{1}{2}\,(-x_{11} + z_{11})\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_{11} + \tfrac{1}{\sqrt{3}}y_{11} - \tfrac{1}{2\sqrt{3}}z_{11}\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,(x_{11} + y_{11} + z_{11})\,c\,\hat{\mathbf{z}} \quad (12i) \quad \text{Si}$$

$$\mathbf{B}_{61} = y_{11}\,\mathbf{a}_1 + x_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 = \tfrac{1}{2}\,(y_{11} - z_{11})\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{\sqrt{3}}x_{11} - \tfrac{1}{2\sqrt{3}}y_{11} - \tfrac{1}{2\sqrt{3}}z_{11}\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,(x_{11} + y_{11} + z_{11})\,c\,\hat{\mathbf{z}} \quad (12i) \quad \text{Si}$$

$$\mathbf{B}_{62} = x_{11}\,\mathbf{a}_1 + z_{11}\,\mathbf{a}_2 + y_{11}\,\mathbf{a}_3 = \tfrac{1}{2}\,(x_{11} - y_{11})\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_{11} - \tfrac{1}{2\sqrt{3}}y_{11} + \tfrac{1}{\sqrt{3}}z_{11}\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,(x_{11} + y_{11} + z_{11})\,c\,\hat{\mathbf{z}} \quad (12i) \quad \text{Si}$$

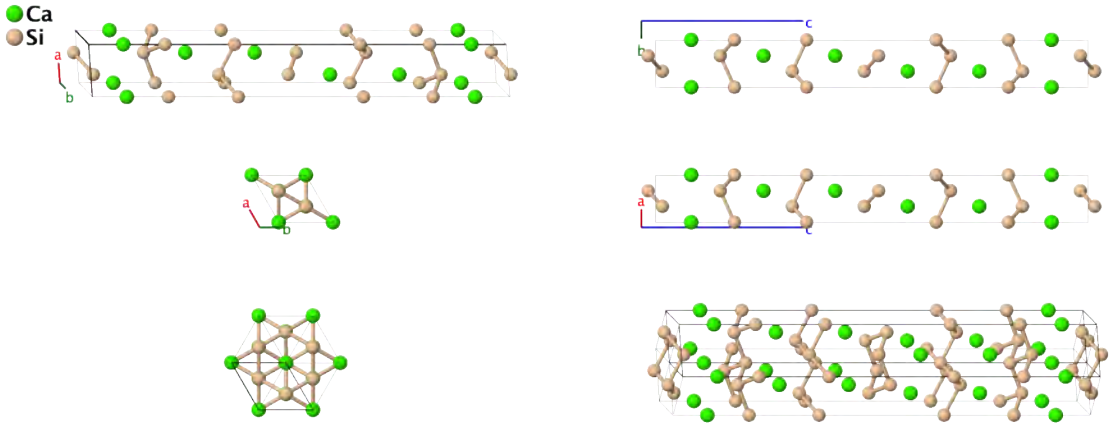
References:

- M. Calligaris, G. Nardin, L. Randaccio, and P. C. Chiaramonti, *Cation-site location in a natural chabazite*, Acta Crystallogr. Sect. B Struct. Sci. **38**, 602–605 (1982), doi:10.1107/S0567740882003483.
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).
- C. Gottfried and F. Schossberger, eds., *Strukturbericht Band III 1933-1935* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).
- C. Gottfried, ed., *Strukturbericht Band V 1937* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1940).

Geometry files:
- CIF: pp. 1737
- POSCAR: pp. 1738

$CaSi_2$ ($C12$) Structure: AB2_hR6_166_c_2c

http://aflow.org/prototype-encyclopedia/AB2_hR6_166_c_2c



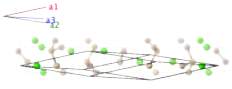
Prototype	:	$CaSi_2$
AFLOW prototype label	:	AB2_hR6_166_c_2c
***Strukturbericht* designation**	:	$C12$
Pearson symbol	:	hR6
Space group number	:	166
Space group symbol	:	$R\bar{3}m$
AFLOW prototype command	:	`aflow --proto=AB2_hR6_166_c_2c [--hex] --params=`$a, c/a, x_1, x_2, x_3$

Other compounds with this structure

- $CaGe_2$

Rhombohedral primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = -\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + x_1\,\mathbf{a}_2 + x_1\,\mathbf{a}_3$	=	$x_1 c\,\hat{\mathbf{z}}$	($2c$)	Ca
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_1 - x_1\,\mathbf{a}_2 - x_1\,\mathbf{a}_3$	=	$-x_1 c\,\hat{\mathbf{z}}$	($2c$)	Ca
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$x_2 c\,\hat{\mathbf{z}}$	($2c$)	Si I
$\mathbf{B}_4$	=	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 - x_2\,\mathbf{a}_3$	=	$-x_2 c\,\hat{\mathbf{z}}$	($2c$)	Si I
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3 c\,\hat{\mathbf{z}}$	($2c$)	Si II
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$-x_3 c\,\hat{\mathbf{z}}$	($2c$)	Si II

References:

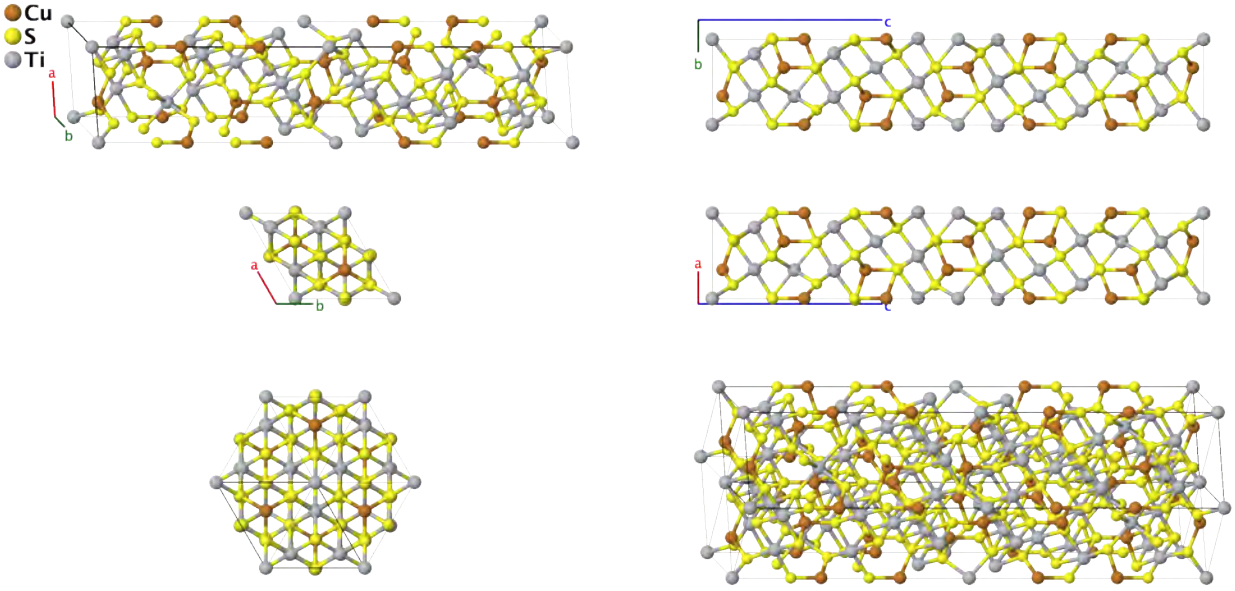
- S. M. Castillo, Z. Tang, A. P. Litvinchuk, and A. M. Guloy, *Lattice Dynamics of the Rhombohedral Polymorphs of $CaSi_2$*, Inorg. Chem. **55**, 10203–10207 (2016), doi:10.1021/acs.inorgchem.6b01399.

Geometry files:

- CIF: pp. 1738
- POSCAR: pp. 1739

Rhombohedral $CuTi_2S_4$ Structure: AB4C2_hR28_166_2c_2c2h_abh

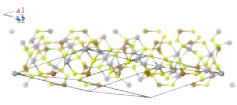
http://aflow.org/prototype-encyclopedia/AB4C2_hR28_166_2c_2c2h_abh



Prototype	:	CuS_4Ti_2
AFLOW prototype label	:	AB4C2_hR28_166_2c_2c2h_abh
***Strukturbericht* designation**	:	None
Pearson symbol	:	hR28
Space group number	:	166
Space group symbol	:	$R\bar{3}m$
AFLOW prototype command	:	`aflow --proto=AB4C2_hR28_166_2c_2c2h_abh [--hex] --params=`$a, c/a, x_3, x_4, x_5, x_6, x_7, z_7, x_8, z_8, x_9, z_9$

Rhombohedral primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = -\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(1a)$	Ti I
$\mathbf{B}_2$	$\frac{1}{2}\mathbf{a}_1 + \frac{1}{2}\mathbf{a}_2 + \frac{1}{2}\mathbf{a}_3$	$\frac{1}{2}c\,\hat{\mathbf{z}}$	$(1b)$	Ti II
$\mathbf{B}_3$	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	$x_3 c\,\hat{\mathbf{z}}$	$(2c)$	Cu I
$\mathbf{B}_4$	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	$-x_3 c\,\hat{\mathbf{z}}$	$(2c)$	Cu I
$\mathbf{B}_5$	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	$x_4 c\,\hat{\mathbf{z}}$	$(2c)$	Cu II
$\mathbf{B}_6$	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	$-x_4 c\,\hat{\mathbf{z}}$	$(2c)$	Cu II

$\mathbf{B}_7$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$x_5 c\,\hat{\mathbf{z}}$	($2c$)	S I
$\mathbf{B}_8$	=	$-x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - x_5\,\mathbf{a}_3$	=	$-x_5 c\,\hat{\mathbf{z}}$	($2c$)	S I
$\mathbf{B}_9$	=	$x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + x_6\,\mathbf{a}_3$	=	$x_6 c\,\hat{\mathbf{z}}$	($2c$)	S II
$\mathbf{B}_{10}$	=	$-x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 - x_6\,\mathbf{a}_3$	=	$-x_6 c\,\hat{\mathbf{z}}$	($2c$)	S II
$\mathbf{B}_{11}$	=	$x_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$\frac{1}{2}(x_7 - z_7)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}(x_7 - z_7)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_7 + \frac{1}{3}z_7\right)c\,\hat{\mathbf{z}}$	($6h$)	S III
$\mathbf{B}_{12}$	=	$z_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + x_7\,\mathbf{a}_3$	=	$\frac{1}{2}(-x_7 + z_7)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}(x_7 - z_7)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_7 + \frac{1}{3}z_7\right)c\,\hat{\mathbf{z}}$	($6h$)	S III
$\mathbf{B}_{13}$	=	$x_7\,\mathbf{a}_1 + z_7\,\mathbf{a}_2 + x_7\,\mathbf{a}_3$	=	$\frac{1}{\sqrt{3}}(-x_7 + z_7)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_7 + \frac{1}{3}z_7\right)c\,\hat{\mathbf{z}}$	($6h$)	S III
$\mathbf{B}_{14}$	=	$-z_7\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 - x_7\,\mathbf{a}_3$	=	$\frac{1}{2}(x_7 - z_7)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}(-x_7 + z_7)\,a\,\hat{\mathbf{y}} - \left(\frac{2}{3}x_7 + \frac{1}{3}z_7\right)c\,\hat{\mathbf{z}}$	($6h$)	S III
$\mathbf{B}_{15}$	=	$-x_7\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$\frac{1}{2}(-x_7 + z_7)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}(-x_7 + z_7)\,a\,\hat{\mathbf{y}} - \left(\frac{2}{3}x_7 + \frac{1}{3}z_7\right)c\,\hat{\mathbf{z}}$	($6h$)	S III
$\mathbf{B}_{16}$	=	$-x_7\,\mathbf{a}_1 - z_7\,\mathbf{a}_2 - x_7\,\mathbf{a}_3$	=	$\frac{1}{\sqrt{3}}(x_7 - z_7)\,a\,\hat{\mathbf{y}} - \left(\frac{2}{3}x_7 + \frac{1}{3}z_7\right)c\,\hat{\mathbf{z}}$	($6h$)	S III
$\mathbf{B}_{17}$	=	$x_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$\frac{1}{2}(x_8 - z_8)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}(x_8 - z_8)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_8 + \frac{1}{3}z_8\right)c\,\hat{\mathbf{z}}$	($6h$)	S IV
$\mathbf{B}_{18}$	=	$z_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 + x_8\,\mathbf{a}_3$	=	$\frac{1}{2}(-x_8 + z_8)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}(x_8 - z_8)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_8 + \frac{1}{3}z_8\right)c\,\hat{\mathbf{z}}$	($6h$)	S IV
$\mathbf{B}_{19}$	=	$x_8\,\mathbf{a}_1 + z_8\,\mathbf{a}_2 + x_8\,\mathbf{a}_3$	=	$\frac{1}{\sqrt{3}}(-x_8 + z_8)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_8 + \frac{1}{3}z_8\right)c\,\hat{\mathbf{z}}$	($6h$)	S IV
$\mathbf{B}_{20}$	=	$-z_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 - x_8\,\mathbf{a}_3$	=	$\frac{1}{2}(x_8 - z_8)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}(-x_8 + z_8)\,a\,\hat{\mathbf{y}} - \left(\frac{2}{3}x_8 + \frac{1}{3}z_8\right)c\,\hat{\mathbf{z}}$	($6h$)	S IV
$\mathbf{B}_{21}$	=	$-x_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$\frac{1}{2}(-x_8 + z_8)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}(-x_8 + z_8)\,a\,\hat{\mathbf{y}} - \left(\frac{2}{3}x_8 + \frac{1}{3}z_8\right)c\,\hat{\mathbf{z}}$	($6h$)	S IV
$\mathbf{B}_{22}$	=	$-x_8\,\mathbf{a}_1 - z_8\,\mathbf{a}_2 - x_8\,\mathbf{a}_3$	=	$\frac{1}{\sqrt{3}}(x_8 - z_8)\,a\,\hat{\mathbf{y}} - \left(\frac{2}{3}x_8 + \frac{1}{3}z_8\right)c\,\hat{\mathbf{z}}$	($6h$)	S IV
$\mathbf{B}_{23}$	=	$x_9\,\mathbf{a}_1 + x_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$\frac{1}{2}(x_9 - z_9)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}(x_9 - z_9)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_9 + \frac{1}{3}z_9\right)c\,\hat{\mathbf{z}}$	($6h$)	Ti III
$\mathbf{B}_{24}$	=	$z_9\,\mathbf{a}_1 + x_9\,\mathbf{a}_2 + x_9\,\mathbf{a}_3$	=	$\frac{1}{2}(-x_9 + z_9)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}(x_9 - z_9)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_9 + \frac{1}{3}z_9\right)c\,\hat{\mathbf{z}}$	($6h$)	Ti III
$\mathbf{B}_{25}$	=	$x_9\,\mathbf{a}_1 + z_9\,\mathbf{a}_2 + x_9\,\mathbf{a}_3$	=	$\frac{1}{\sqrt{3}}(-x_9 + z_9)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_9 + \frac{1}{3}z_9\right)c\,\hat{\mathbf{z}}$	($6h$)	Ti III
$\mathbf{B}_{26}$	=	$-z_9\,\mathbf{a}_1 - x_9\,\mathbf{a}_2 - x_9\,\mathbf{a}_3$	=	$\frac{1}{2}(x_9 - z_9)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}(-x_9 + z_9)\,a\,\hat{\mathbf{y}} - \left(\frac{2}{3}x_9 + \frac{1}{3}z_9\right)c\,\hat{\mathbf{z}}$	($6h$)	Ti III
$\mathbf{B}_{27}$	=	$-x_9\,\mathbf{a}_1 - x_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$\frac{1}{2}(-x_9 + z_9)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}(-x_9 + z_9)\,a\,\hat{\mathbf{y}} - \left(\frac{2}{3}x_9 + \frac{1}{3}z_9\right)c\,\hat{\mathbf{z}}$	($6h$)	Ti III
$\mathbf{B}_{28}$	=	$-x_9\,\mathbf{a}_1 - z_9\,\mathbf{a}_2 - x_9\,\mathbf{a}_3$	=	$\frac{1}{\sqrt{3}}(x_9 - z_9)\,a\,\hat{\mathbf{y}} - \left(\frac{2}{3}x_9 + \frac{1}{3}z_9\right)c\,\hat{\mathbf{z}}$	($6h$)	Ti III

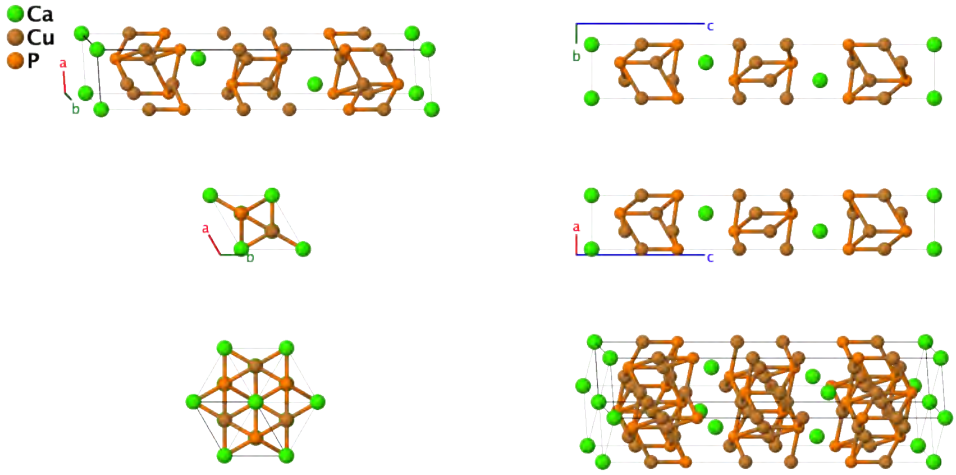
References:

- N. Soheilnia, K. M. Kleinke, E. Dashjav, H. L. Cuthbert, J. E. Greedan, and H. Kleinke, *Crystal Structure and Physical Properties of a New $CuTi_2S_4$ Modification in Comparison to the Thiospinel*, Inorg. Chem. **43**, 6473–6478 (2004), doi:10.1021/ic0495113.

Geometry files:
- CIF: pp. 1739
- POSCAR: pp. 1739

$CaCu_4P_2$ Structure: AB4C2_hR7_166_a_2c_c

http://aflow.org/prototype-encyclopedia/AB4C2_hR7_166_a_2c_c



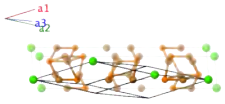
Prototype	:	$CaCu_4P_2$
AFLOW prototype label	:	AB4C2_hR7_166_a_2c_c
***Strukturbericht* designation**	:	None
Pearson symbol	:	hR7
Space group number	:	166
Space group symbol	:	$R\bar{3}m$
AFLOW prototype command	:	`aflow --proto=AB4C2_hR7_166_a_2c_c [--hex] --params=`$a, c/a, x_2, x_3, x_4$

Other compounds with this structure

- $BaAg_4As_2$, $CaAg_4As_2$, $EuAg_4As_2$, $EuAg_4Sb_2$, $SrAg_4As_2$, $SrAg_4Sb_2$, and $ZrLi_4Ge_2$

Rhombohedral primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = -\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(1a)$	Ca
$\mathbf{B}_2$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$x_2 c\,\hat{\mathbf{z}}$	$(2c)$	Cu I
$\mathbf{B}_3$	=	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 - x_2\,\mathbf{a}_3$	=	$-x_2 c\,\hat{\mathbf{z}}$	$(2c)$	Cu I
$\mathbf{B}_4$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3 c\,\hat{\mathbf{z}}$	$(2c)$	Cu II
$\mathbf{B}_5$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$-x_3 c\,\hat{\mathbf{z}}$	$(2c)$	Cu II
$\mathbf{B}_6$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4 c\,\hat{\mathbf{z}}$	$(2c)$	P
$\mathbf{B}_7$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$-x_4 c\,\hat{\mathbf{z}}$	$(2c)$	P

References:

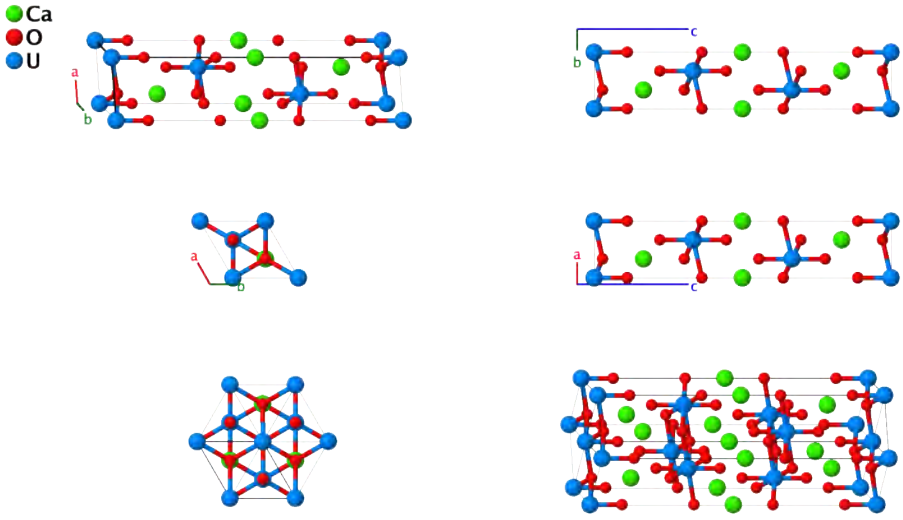
- A. Mewis, *Darstellung und Struktur der Verbindung* $CaCu_4P_2$, Z. Naturforsch. B **35**, 942–945 (1980), doi:10.1515/znb-1980-0802.

Geometry files:

- CIF: pp. 1739
- POSCAR: pp. 1740

$CaUO_4$ Structure: AB4C_hR6_166_b_2c_a

http://aflow.org/prototype-encyclopedia/AB4C_hR6_166_b_2c_a



Prototype	:	CaO_4U
AFLOW prototype label	:	AB4C_hR6_166_b_2c_a
***Strukturbericht* designation**	:	None
Pearson symbol	:	hR6
Space group number	:	166
Space group symbol	:	$R\bar{3}m$
AFLOW prototype command	:	`aflow --proto=AB4C_hR6_166_b_2c_a [--hex] --params=`$a, c/a, x_3, x_4$

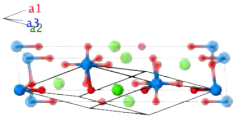
Other compounds with this structure

- $NaUO_4$ (clarkeite)

- Hexagonal settings of this structure can be obtained with the option `--hex`.

Rhombohedral primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = -\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(1a)$	U
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	$(1b)$	Ca
$\mathbf{B}_3$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3 c\,\hat{\mathbf{z}}$	$(2c)$	O I
$\mathbf{B}_4$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$-x_3 c\,\hat{\mathbf{z}}$	$(2c)$	O I

$\mathbf{B}_5$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4 c\,\hat{\mathbf{z}}$	$(2c)$	O II
$\mathbf{B}_6$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$-x_4 c\,\hat{\mathbf{z}}$	$(2c)$	O II

References:

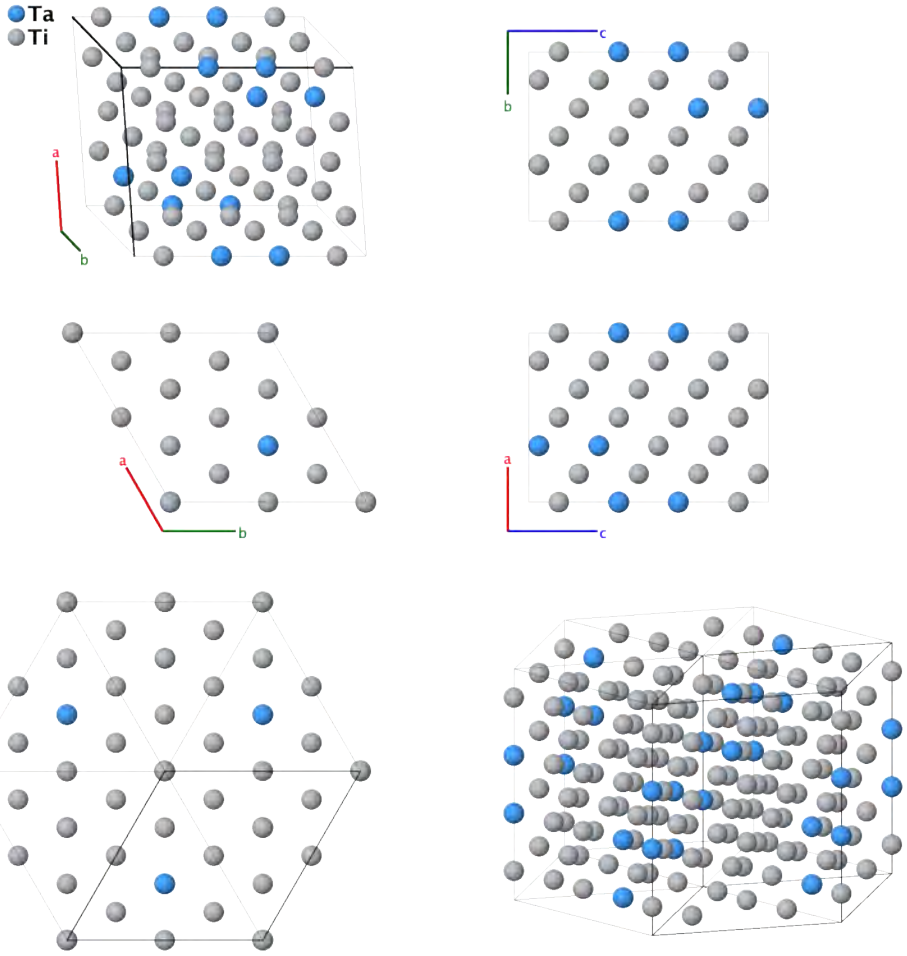
- B. O. Loopstra and H. M. Rietveld, *The structure of some alkaline-earth metal uranates*, Acta Crystallogr. Sect. B Struct. Sci. **25**, 787–791 (1969), doi:10.1107/S0567740869002974.

Geometry files:

- CIF: pp. 1740
- POSCAR: pp. 1741

$TaTi_7$ (BCC SQS-16) Structure: AB7_hR16_166_c_c2h

http://aflow.org/prototype-encyclopedia/AB7_hR16_166_c_c2h

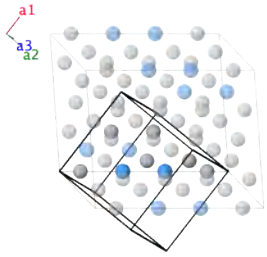


Prototype	:	$TaTi_7$
AFLOW prototype label	:	AB7_hR16_166_c_c2h
***Strukturbericht* designation**	:	None
Pearson symbol	:	hR16
Space group number	:	166
Space group symbol	:	$R\bar{3}m$
AFLOW prototype command	:	`aflow --proto=AB7_hR16_166_c_c2h [--hex] --params=`$a, c/a, x_1, x_2, x_3, z_3, x_4, z_4$

- This is a special quasirandom structure with 16 atoms per unit cell (SQS-16) for the β-phase (high-temperature austenite) bcc substitutional Ti-Ta alloy (Chakraborty, 2016). This protoype contains 12.5% Ta. Prototypes are listed for other Ta-Ti concentrations: 18.75% Ta (A3B13_oC32_38_ac_a2bcdef), 25% Ta (AB3_mC32_8_4a_4a4b), 31.25% Ta (A5B11_mP16_6_2abc_2a3b3c), and 37.5% Ta (A3B5_oC32_38_abce_abcdf).

Rhombohedral primitive vectors:

$$\begin{aligned}
\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2\sqrt{3}}\,a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,c\,\hat{\mathbf{z}} \\
\mathbf{a}_2 &= \tfrac{1}{\sqrt{3}}\,a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,c\,\hat{\mathbf{z}} \\
\mathbf{a}_3 &= -\tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2\sqrt{3}}\,a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,c\,\hat{\mathbf{z}}
\end{aligned}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$x_1\,\mathbf{a}_1 + x_1\,\mathbf{a}_2 + x_1\,\mathbf{a}_3$	$x_1 c\,\hat{\mathbf{z}}$	$(2c)$	Ta
$\mathbf{B}_2$	$-x_1\,\mathbf{a}_1 - x_1\,\mathbf{a}_2 - x_1\,\mathbf{a}_3$	$-x_1 c\,\hat{\mathbf{z}}$	$(2c)$	Ta
$\mathbf{B}_3$	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	$x_2 c\,\hat{\mathbf{z}}$	$(2c)$	Ti I
$\mathbf{B}_4$	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 - x_2\,\mathbf{a}_3$	$-x_2 c\,\hat{\mathbf{z}}$	$(2c)$	Ti I
$\mathbf{B}_5$	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$\frac{1}{2}(x_3 - z_3)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}(x_3 - z_3)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_3 + \frac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	$(6h)$	Ti II
$\mathbf{B}_6$	$z_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	$\frac{1}{2}(-x_3 + z_3)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}(x_3 - z_3)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_3 + \frac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	$(6h)$	Ti II
$\mathbf{B}_7$	$x_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	$\frac{1}{\sqrt{3}}(-x_3 + z_3)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_3 + \frac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	$(6h)$	Ti II
$\mathbf{B}_8$	$-z_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	$\frac{1}{2}(x_3 - z_3)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}(-x_3 + z_3)\,a\,\hat{\mathbf{y}} - \left(\frac{2}{3}x_3 + \frac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	$(6h)$	Ti II
$\mathbf{B}_9$	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	$\frac{1}{2}(-x_3 + z_3)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}(-x_3 + z_3)\,a\,\hat{\mathbf{y}} - \left(\frac{2}{3}x_3 + \frac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	$(6h)$	Ti II
$\mathbf{B}_{10}$	$-x_3\,\mathbf{a}_1 - z_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	$\frac{1}{\sqrt{3}}(x_3 - z_3)\,a\,\hat{\mathbf{y}} - \left(\frac{2}{3}x_3 + \frac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	$(6h)$	Ti II
$\mathbf{B}_{11}$	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	$\frac{1}{2}(x_4 - z_4)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}(x_4 - z_4)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_4 + \frac{1}{3}z_4\right)c\,\hat{\mathbf{z}}$	$(6h)$	Ti III
$\mathbf{B}_{12}$	$z_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	$\frac{1}{2}(-x_4 + z_4)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}(x_4 - z_4)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_4 + \frac{1}{3}z_4\right)c\,\hat{\mathbf{z}}$	$(6h)$	Ti III
$\mathbf{B}_{13}$	$x_4\,\mathbf{a}_1 + z_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	$\frac{1}{\sqrt{3}}(-x_4 + z_4)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_4 + \frac{1}{3}z_4\right)c\,\hat{\mathbf{z}}$	$(6h)$	Ti III
$\mathbf{B}_{14}$	$-z_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	$\frac{1}{2}(x_4 - z_4)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}(-x_4 + z_4)\,a\,\hat{\mathbf{y}} - \left(\frac{2}{3}x_4 + \frac{1}{3}z_4\right)c\,\hat{\mathbf{z}}$	$(6h)$	Ti III
$\mathbf{B}_{15}$	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	$\frac{1}{2}(-x_4 + z_4)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}(-x_4 + z_4)\,a\,\hat{\mathbf{y}} - \left(\frac{2}{3}x_4 + \frac{1}{3}z_4\right)c\,\hat{\mathbf{z}}$	$(6h)$	Ti III
$\mathbf{B}_{16}$	$-x_4\,\mathbf{a}_1 - z_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	$\frac{1}{\sqrt{3}}(x_4 - z_4)\,a\,\hat{\mathbf{y}} - \left(\frac{2}{3}x_4 + \frac{1}{3}z_4\right)c\,\hat{\mathbf{z}}$	$(6h)$	Ti III

References:

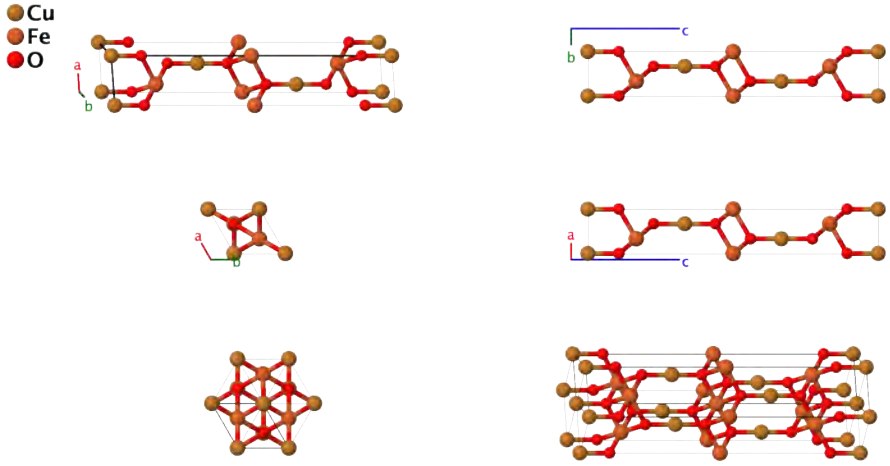
- T. Chakraborty, J. Rogal, and R. Drautz, *Unraveling the composition dependence of the martensitic transformation temperature: A first-principles study of Ti-Ta alloys*, Phys. Rev. B **94**, 224104 (2016), doi:10.1103/PhysRevB.94.224104.

Geometry files:

- CIF: pp. 1741
- POSCAR: pp. 1741

Rhombohedral Delafossite ($CuFeO_2$) Structure: ABC2_hR4_166_a_b_c

http://aflow.org/prototype-encyclopedia/ABC2_hR4_166_a_b_c.CuFeO2



Prototype	:	$CuFeO_2$
AFLOW prototype label	:	ABC2_hR4_166_a_b_c
***Strukturbericht* designation**	:	None
Pearson symbol	:	hR4
Space group number	:	166
Space group symbol	:	$R\bar{3}m$
AFLOW prototype command	:	`aflow --proto=ABC2_hR4_166_a_b_c [--hex] --params=`$a, c/a, x_3$

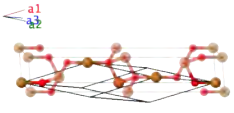
Other compounds with this structure

- $AgAlO_2$, $AgCoO_2$, $AgCrO_2$, $AgFeO_2$, $AgGaO_2$, $AgInO_2$, $AgNiO_2$, $AgRhO_2$, $AgScO_2$, $AgTlO_2$, $CuAlO_2$, $CuCoO_2$, $CuCrO_2$, $CuEuO_2$, $CuGaO_2$, $CuInO_2$, $CuLaO_2$, $CuRhO_2$, $CuScO_2$, $CuYO_2$, $PdCoO_2$, $PdCrO_2$, $PdRhO_2$, and $PtCoO_2$

- Delafossite appears in two forms which differ in the stacking of the layers: rhombohedral, shown here, and hexagonal, prototype $CuAlO_2$. Most of the compounds found in the hexagonal phase can also be found in the rhombohedral structure (Marquardt, 2006).
- Rhombohedral delafossite has the same AFLOW label, ABC2_hR4_166_a_b_c, as caswellsilverite $F5_1$. The difference in the internal parameter z_3 causes a large change in the bonding of the two crystals, so we present them as different structures.

Rhombohedral primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = -\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($1a$)	Cu
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	($1b$)	Fe
$\mathbf{B}_3$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3 c\,\hat{\mathbf{z}}$	($2c$)	O
$\mathbf{B}_4$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$-x_3 c\,\hat{\mathbf{z}}$	($2c$)	O

References:

- C. T. Prewitt, R. D. Shannon, and D. B. Rogers, *Chemistry of noble metal oxides. II. Crystal structures of platinum cobalt dioxide, palladium cobalt dioxide, coppper iron dioxide, and silver iron dioxide*, Inorg. Chem. **10**, 719–723 (1971), doi:10.1021/ic50098a012.

Found in:

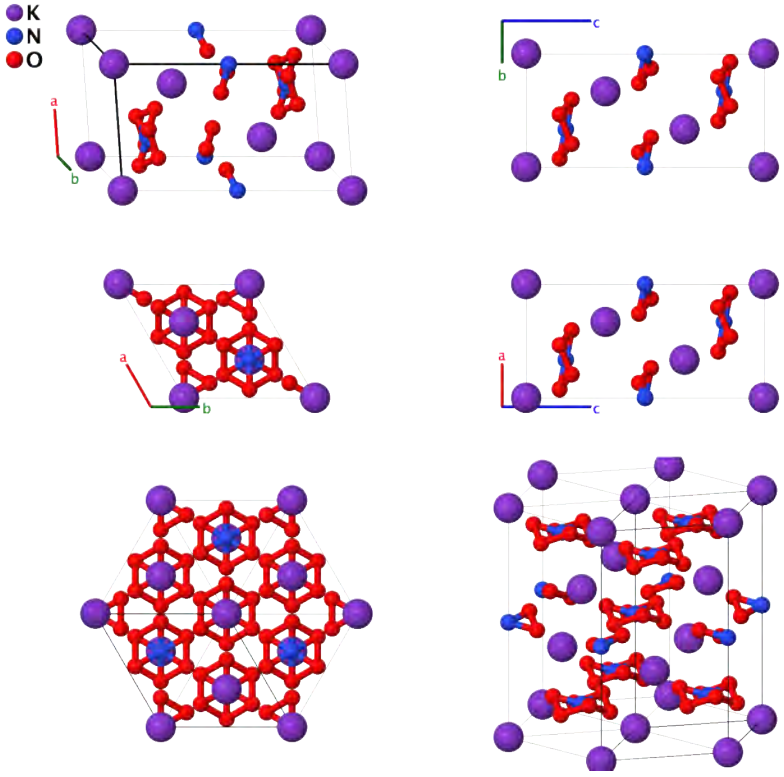
- M. A. Marquardt, N. A. Ashmore, and D. P. Cann, *Crystal chemistry and electrical properties of the delafossite structure*, Thin Solid Films **496**, 146–156 (2006), doi:10.1016/j.tsf.2005.08.316.

Geometry files:

- CIF: pp. 1741
- POSCAR: pp. 1742

β-Potassium Nitrate (KNO_3) Structure: ABC6_hR8_166_a_b_h

http://aflow.org/prototype-encyclopedia/ABC6_hR8_166_a_b_h

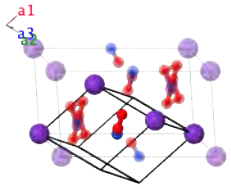


Prototype	:	KNO_3
AFLOW prototype label	:	ABC6_hR8_166_a_b_h
***Strukturbericht* designation**	:	None
Pearson symbol	:	hR8
Space group number	:	166
Space group symbol	:	$R\bar{3}m$
AFLOW prototype command	:	`aflow --proto=ABC6_hR8_166_a_b_h [--hex] --params=`$a, c/a, x_3, z_3$

- On heating, α-KNO_3 (either Structure I or Structure II) transforms into β-KNO_3 at 128 °C. When heated above 200 °C and then cooled, the β phase transforms into the metastable ferroelectric γ-KNO_3 phase, which can remain down to room temperature.
- In the β-phase the oxygen ($6h$) sites are only 50% filled. Presumably only the $+z$ or $-z$ sites will be occupied around a given nitrogen atom.
- (Nimmo, 1976) give the data for β-KNO_3 taken at 151 °C.

Rhombohedral primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}\,a\,\hat{\mathbf{y}} + \frac{1}{3}\,c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{\sqrt{3}}\,a\,\hat{\mathbf{y}} + \frac{1}{3}\,c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = -\frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}\,a\,\hat{\mathbf{y}} + \frac{1}{3}\,c\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(1*a*)	K
$\mathbf{B}_2$	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	$\frac{1}{2}c\,\hat{\mathbf{z}}$	(1*b*)	N
$\mathbf{B}_3$	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$\frac{1}{2}\,(x_3 - z_3)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}\,(x_3 - z_3)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_3 + \frac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	(6*h*)	O
$\mathbf{B}_4$	$z_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	$\frac{1}{2}\,(-x_3 + z_3)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}\,(x_3 - z_3)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_3 + \frac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	(6*h*)	O
$\mathbf{B}_5$	$x_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	$\frac{1}{\sqrt{3}}\,(-x_3 + z_3)\,a\,\hat{\mathbf{y}} + \left(\frac{2}{3}x_3 + \frac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	(6*h*)	O
$\mathbf{B}_6$	$-z_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	$\frac{1}{2}\,(x_3 - z_3)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}\,(-x_3 + z_3)\,a\,\hat{\mathbf{y}} - \left(\frac{2}{3}x_3 + \frac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	(6*h*)	O
$\mathbf{B}_7$	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	$\frac{1}{2}\,(-x_3 + z_3)\,a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}\,(-x_3 + z_3)\,a\,\hat{\mathbf{y}} - \left(\frac{2}{3}x_3 + \frac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	(6*h*)	O
$\mathbf{B}_8$	$-x_3\,\mathbf{a}_1 - z_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	$\frac{1}{\sqrt{3}}\,(x_3 - z_3)\,a\,\hat{\mathbf{y}} - \left(\frac{2}{3}x_3 + \frac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	(6*h*)	O

References:

- J. K. Nimmo and B. W. Lucas, *The crystal structures of γ-and β-KNO_3 and the α-β-γ phase transformations*, Acta Crystallogr. Sect. B Struct. Sci. **32**, 1968–1971 (1976), doi:10.1107/S0567740876006894.

Found in:

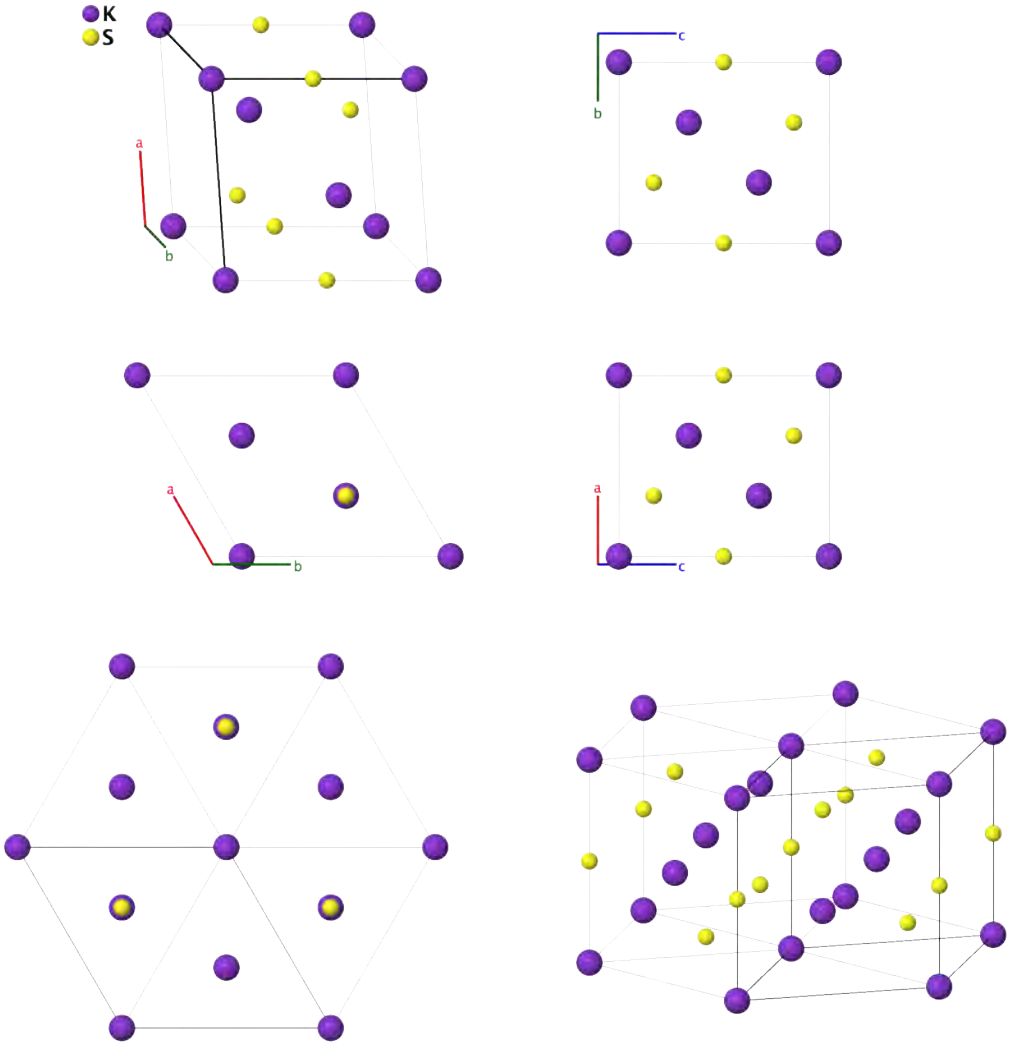
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1742
- POSCAR: pp. 1742

K(SH) (*B*22) Structure: AB_hR2_166_a_b

http://aflow.org/prototype-encyclopedia/AB_hR2_166_a_b.KSH



Prototype	:	KS
AFLOW prototype label	:	AB_hR2_166_a_b
***Strukturbericht* designation**	:	*B*22
Pearson symbol	:	hR2
Space group number	:	166
Space group symbol	:	$R\bar{3}m$
AFLOW prototype command	:	`aflow --proto=AB_hR2_166_a_b [--hex] --params=`$a, c/a$

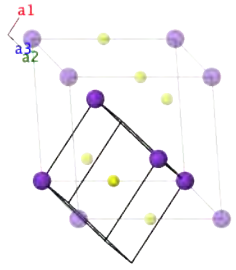
Other compounds with this structure

- Na(SH), Rb(SH), and Cs(SH)

- This is the room-temperature structure of KSH. The S and H atoms form an SH^- ion, and so are listed together at the (1*b*) Wyckoff postition, leading to the listing of this compound in the *B Strukturbericht* category.

Rhombohedral primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}\,a\,\hat{\mathbf{y}} + \frac{1}{3}\,c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{\sqrt{3}}\,a\,\hat{\mathbf{y}} + \frac{1}{3}\,c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = -\frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}\,a\,\hat{\mathbf{y}} + \frac{1}{3}\,c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(1*a*)	K
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	(1*b*)	S

References:

- C. D. West, *The Crystal Structures of Some Alkali Hydrosulfides and Monosulfides*, Zeitschrift für Kristallographie - Crystalline Materials **88**, 97–115 (1934), doi:10.1524/zkri.1934.88.1.97.

Found in:

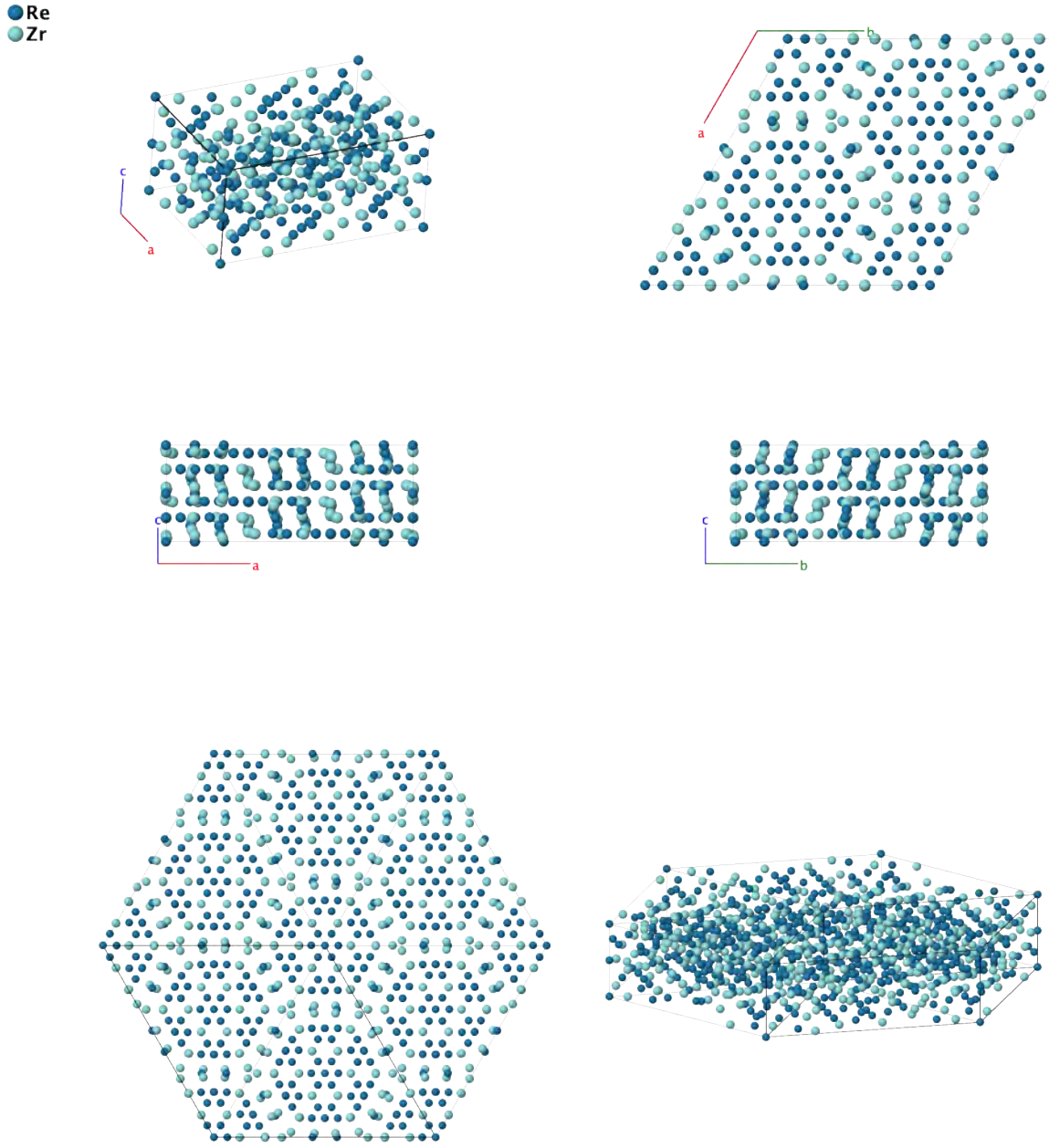
- C. Gottfried and F. Schossberger, eds., *Strukturbericht Band III 1933-1935* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1742
- POSCAR: pp. 1743

$Zr_{21}Re_{25}$ Structure: A25B21_hR92_167_b2e3f_e3f

http://aflow.org/prototype-encyclopedia/A25B21_hR92_167_b2e3f_e3f



Prototype	:	$Re_{25}Zr_{21}$
AFLOW prototype label	:	A25B21_hR92_167_b2e3f_e3f
***Strukturbericht* designation**	:	None
Pearson symbol	:	hR92
Space group number	:	167
Space group symbol	:	$R\bar{3}c$

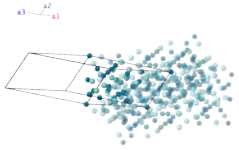
AFLOW prototype command : `aflow --proto=A25B21_hR92_167_b2e3f_e3f [--hex] --params=`$a, c/a, x_2, x_3, x_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}$

Other compounds with this structure

- $Mg_{21}Zn_{25}$

Rhombohedral primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2\sqrt{3}}\,a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,c\,\hat{\mathbf{z}}\\ \mathbf{a}_2 &= \tfrac{1}{\sqrt{3}}\,a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,c\,\hat{\mathbf{z}}\\ \mathbf{a}_3 &= -\tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2\sqrt{3}}\,a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($2b$)	Re I
$\mathbf{B}_2$	=	$\tfrac{1}{2}\,\mathbf{a}_1 + \tfrac{1}{2}\,\mathbf{a}_2 + \tfrac{1}{2}\,\mathbf{a}_3$	=	$\tfrac{1}{2}c\,\hat{\mathbf{z}}$	($2b$)	Re I
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_2\right)\mathbf{a}_2 + \tfrac{1}{4}\,\mathbf{a}_3$	=	$\left(-\tfrac{1}{8} + \tfrac{1}{2}x_2\right)a\,\hat{\mathbf{x}} + \left(\tfrac{\sqrt{3}}{8} - \tfrac{\sqrt{3}}{2}x_2\right)a\,\hat{\mathbf{y}} + \tfrac{1}{4}c\,\hat{\mathbf{z}}$	($6e$)	Re II
$\mathbf{B}_4$	=	$\tfrac{1}{4}\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_2\right)\mathbf{a}_3$	=	$\left(-\tfrac{1}{8} + \tfrac{1}{2}x_2\right)a\,\hat{\mathbf{x}} + \left(-\tfrac{\sqrt{3}}{8} + \tfrac{\sqrt{3}}{2}x_2\right)a\,\hat{\mathbf{y}} + \tfrac{1}{4}c\,\hat{\mathbf{z}}$	($6e$)	Re II
$\mathbf{B}_5$	=	$\left(\tfrac{1}{2} - x_2\right)\mathbf{a}_1 + \tfrac{1}{4}\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$\left(\tfrac{1}{4} - x_2\right)a\,\hat{\mathbf{x}} + \tfrac{1}{4}c\,\hat{\mathbf{z}}$	($6e$)	Re II
$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_2\right)\mathbf{a}_2 + \tfrac{3}{4}\,\mathbf{a}_3$	=	$-\left(\tfrac{1}{2}x_2 + \tfrac{3}{8}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{8\sqrt{3}} + \tfrac{\sqrt{3}}{2}x_2\right)a\,\hat{\mathbf{y}} + \tfrac{5}{12}c\,\hat{\mathbf{z}}$	($6e$)	Re II
$\mathbf{B}_7$	=	$\tfrac{3}{4}\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_2\right)\mathbf{a}_3$	=	$\left(\tfrac{1}{8} - \tfrac{1}{2}x_2\right)a\,\hat{\mathbf{x}} - \left(\tfrac{\sqrt{3}}{2}x_2 + \tfrac{5}{8\sqrt{3}}\right)a\,\hat{\mathbf{y}} + \tfrac{5}{12}c\,\hat{\mathbf{z}}$	($6e$)	Re II
$\mathbf{B}_8$	=	$\left(\tfrac{1}{2} + x_2\right)\mathbf{a}_1 + \tfrac{3}{4}\,\mathbf{a}_2 - x_2\,\mathbf{a}_3$	=	$\left(\tfrac{1}{4} + x_2\right)a\,\hat{\mathbf{x}} + \tfrac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \tfrac{5}{12}c\,\hat{\mathbf{z}}$	($6e$)	Re II
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_3\right)\mathbf{a}_2 + \tfrac{1}{4}\,\mathbf{a}_3$	=	$\left(-\tfrac{1}{8} + \tfrac{1}{2}x_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{\sqrt{3}}{8} - \tfrac{\sqrt{3}}{2}x_3\right)a\,\hat{\mathbf{y}} + \tfrac{1}{4}c\,\hat{\mathbf{z}}$	($6e$)	Re III
$\mathbf{B}_{10}$	=	$\tfrac{1}{4}\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_3\right)\mathbf{a}_3$	=	$\left(-\tfrac{1}{8} + \tfrac{1}{2}x_3\right)a\,\hat{\mathbf{x}} + \left(-\tfrac{\sqrt{3}}{8} + \tfrac{\sqrt{3}}{2}x_3\right)a\,\hat{\mathbf{y}} + \tfrac{1}{4}c\,\hat{\mathbf{z}}$	($6e$)	Re III
$\mathbf{B}_{11}$	=	$\left(\tfrac{1}{2} - x_3\right)\mathbf{a}_1 + \tfrac{1}{4}\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$\left(\tfrac{1}{4} - x_3\right)a\,\hat{\mathbf{x}} + \tfrac{1}{4}c\,\hat{\mathbf{z}}$	($6e$)	Re III
$\mathbf{B}_{12}$	=	$-x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_3\right)\mathbf{a}_2 + \tfrac{3}{4}\,\mathbf{a}_3$	=	$-\left(\tfrac{1}{2}x_3 + \tfrac{3}{8}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{8\sqrt{3}} + \tfrac{\sqrt{3}}{2}x_3\right)a\,\hat{\mathbf{y}} + \tfrac{5}{12}c\,\hat{\mathbf{z}}$	($6e$)	Re III
$\mathbf{B}_{13}$	=	$\tfrac{3}{4}\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_3\right)\mathbf{a}_3$	=	$\left(\tfrac{1}{8} - \tfrac{1}{2}x_3\right)a\,\hat{\mathbf{x}} - \left(\tfrac{\sqrt{3}}{2}x_3 + \tfrac{5}{8\sqrt{3}}\right)a\,\hat{\mathbf{y}} + \tfrac{5}{12}c\,\hat{\mathbf{z}}$	($6e$)	Re III
$\mathbf{B}_{14}$	=	$\left(\tfrac{1}{2} + x_3\right)\mathbf{a}_1 + \tfrac{3}{4}\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$\left(\tfrac{1}{4} + x_3\right)a\,\hat{\mathbf{x}} + \tfrac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \tfrac{5}{12}c\,\hat{\mathbf{z}}$	($6e$)	Re III

$$\mathbf{B}_{15} = x_4\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_4\right)\mathbf{a}_2 + \tfrac{1}{4}\,\mathbf{a}_3 = \left(-\tfrac{1}{8} + \tfrac{1}{2}x_4\right)a\,\hat{\mathbf{x}} + \left(\tfrac{\sqrt{3}}{8} - \tfrac{\sqrt{3}}{2}x_4\right)a\,\hat{\mathbf{y}} + \tfrac{1}{4}c\,\hat{\mathbf{z}} \quad (6e) \quad \text{Zr I}$$

$$\mathbf{B}_{16} = \tfrac{1}{4}\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_4\right)\mathbf{a}_3 = \left(-\tfrac{1}{8} + \tfrac{1}{2}x_4\right)a\,\hat{\mathbf{x}} + \left(-\tfrac{\sqrt{3}}{8} + \tfrac{\sqrt{3}}{2}x_4\right)a\,\hat{\mathbf{y}} + \tfrac{1}{4}c\,\hat{\mathbf{z}} \quad (6e) \quad \text{Zr I}$$

$$\mathbf{B}_{17} = \left(\tfrac{1}{2} - x_4\right)\mathbf{a}_1 + \tfrac{1}{4}\,\mathbf{a}_2 + x_4\,\mathbf{a}_3 = \left(\tfrac{1}{4} - x_4\right)a\,\hat{\mathbf{x}} + \tfrac{1}{4}c\,\hat{\mathbf{z}} \quad (6e) \quad \text{Zr I}$$

$$\mathbf{B}_{18} = -x_4\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_4\right)\mathbf{a}_2 + \tfrac{3}{4}\,\mathbf{a}_3 = -\left(\tfrac{1}{2}x_4 + \tfrac{3}{8}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{8\sqrt{3}} + \tfrac{\sqrt{3}}{2}x_4\right)a\,\hat{\mathbf{y}} + \tfrac{5}{12}c\,\hat{\mathbf{z}} \quad (6e) \quad \text{Zr I}$$

$$\mathbf{B}_{19} = \tfrac{3}{4}\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_4\right)\mathbf{a}_3 = \left(\tfrac{1}{8} - \tfrac{1}{2}x_4\right)a\,\hat{\mathbf{x}} - \left(\tfrac{\sqrt{3}}{2}x_4 + \tfrac{5}{8\sqrt{3}}\right)a\,\hat{\mathbf{y}} + \tfrac{5}{12}c\,\hat{\mathbf{z}} \quad (6e) \quad \text{Zr I}$$

$$\mathbf{B}_{20} = \left(\tfrac{1}{2} + x_4\right)\mathbf{a}_1 + \tfrac{3}{4}\,\mathbf{a}_2 - x_4\,\mathbf{a}_3 = \left(\tfrac{1}{4} + x_4\right)a\,\hat{\mathbf{x}} + \tfrac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \tfrac{5}{12}c\,\hat{\mathbf{z}} \quad (6e) \quad \text{Zr I}$$

$$\mathbf{B}_{21} = x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3 = \tfrac{1}{2}\left(x_5 - z_5\right)a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_5 + \tfrac{1}{\sqrt{3}}y_5 - \tfrac{1}{2\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\left(x_5 + y_5 + z_5\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re IV}$$

$$\mathbf{B}_{22} = z_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + y_5\,\mathbf{a}_3 = \tfrac{1}{2}\left(-y_5 + z_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{\sqrt{3}}x_5 - \tfrac{1}{2\sqrt{3}}y_5 - \tfrac{1}{2\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\left(x_5 + y_5 + z_5\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re IV}$$

$$\mathbf{B}_{23} = y_5\,\mathbf{a}_1 + z_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3 = \tfrac{1}{2}\left(-x_5 + y_5\right)a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_5 - \tfrac{1}{2\sqrt{3}}y_5 + \tfrac{1}{\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\left(x_5 + y_5 + z_5\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re IV}$$

$$\mathbf{B}_{24} = \left(\tfrac{1}{2} - z_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_5\right)\mathbf{a}_3 = \tfrac{1}{2}\left(x_5 - z_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2\sqrt{3}}x_5 - \tfrac{1}{\sqrt{3}}y_5 + \tfrac{1}{2\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - \tfrac{1}{3}x_5 - \tfrac{1}{3}y_5 - \tfrac{1}{3}z_5\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re IV}$$

$$\mathbf{B}_{25} = \left(\tfrac{1}{2} - y_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_5\right)\mathbf{a}_3 = \tfrac{1}{2}\left(-y_5 + z_5\right)a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{\sqrt{3}}x_5 + \tfrac{1}{2\sqrt{3}}y_5 + \tfrac{1}{2\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - \tfrac{1}{3}x_5 - \tfrac{1}{3}y_5 - \tfrac{1}{3}z_5\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re IV}$$

$$\mathbf{B}_{26} = \left(\tfrac{1}{2} - x_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - z_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - y_5\right)\mathbf{a}_3 = \tfrac{1}{2}\left(-x_5 + y_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2\sqrt{3}}x_5 + \tfrac{1}{2\sqrt{3}}y_5 - \tfrac{1}{\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - \tfrac{1}{3}x_5 - \tfrac{1}{3}y_5 - \tfrac{1}{3}z_5\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re IV}$$

$$\mathbf{B}_{27} = -x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3 = \tfrac{1}{2}\left(-x_5 + z_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2\sqrt{3}}x_5 - \tfrac{1}{\sqrt{3}}y_5 + \tfrac{1}{2\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}} - \tfrac{1}{3}\left(x_5 + y_5 + z_5\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re IV}$$

$$\mathbf{B}_{28} = -z_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - y_5\,\mathbf{a}_3 = \tfrac{1}{2}\left(y_5 - z_5\right)a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{\sqrt{3}}x_5 + \tfrac{1}{2\sqrt{3}}y_5 + \tfrac{1}{2\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}} - \tfrac{1}{3}\left(x_5 + y_5 + z_5\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re IV}$$

$$\mathbf{B}_{29} = -y_5\,\mathbf{a}_1 - z_5\,\mathbf{a}_2 - x_5\,\mathbf{a}_3 = \tfrac{1}{2}\left(x_5 - y_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2\sqrt{3}}x_5 + \tfrac{1}{2\sqrt{3}}y_5 - \tfrac{1}{\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}} - \tfrac{1}{3}\left(x_5 + y_5 + z_5\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re IV}$$

$$\mathbf{B}_{30} = \left(\tfrac{1}{2}+z_5\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_5\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_5\right)\mathbf{a}_3 = \tfrac{1}{2}(-x_5+z_5)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_5+\tfrac{1}{\sqrt{3}}y_5-\tfrac{1}{2\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+\tfrac{1}{3}x_5+\tfrac{1}{3}y_5+\tfrac{1}{3}z_5\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re IV}$$

$$\mathbf{B}_{31} = \left(\tfrac{1}{2}+y_5\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_5\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_5\right)\mathbf{a}_3 = \tfrac{1}{2}(y_5-z_5)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{\sqrt{3}}x_5-\tfrac{1}{2\sqrt{3}}y_5-\tfrac{1}{2\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+\tfrac{1}{3}x_5+\tfrac{1}{3}y_5+\tfrac{1}{3}z_5\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re IV}$$

$$\mathbf{B}_{32} = \left(\tfrac{1}{2}+x_5\right)\mathbf{a}_1+\left(\tfrac{1}{2}+z_5\right)\mathbf{a}_2+\left(\tfrac{1}{2}+y_5\right)\mathbf{a}_3 = \tfrac{1}{2}(x_5-y_5)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_5-\tfrac{1}{2\sqrt{3}}y_5+\tfrac{1}{\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+\tfrac{1}{3}x_5+\tfrac{1}{3}y_5+\tfrac{1}{3}z_5\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re IV}$$

$$\mathbf{B}_{33} = x_6\,\mathbf{a}_1+y_6\,\mathbf{a}_2+z_6\,\mathbf{a}_3 = \tfrac{1}{2}(x_6-z_6)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_6+\tfrac{1}{\sqrt{3}}y_6-\tfrac{1}{2\sqrt{3}}z_6\right)a\,\hat{\mathbf{y}}+\tfrac{1}{3}(x_6+y_6+z_6)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re V}$$

$$\mathbf{B}_{34} = z_6\,\mathbf{a}_1+x_6\,\mathbf{a}_2+y_6\,\mathbf{a}_3 = \tfrac{1}{2}(-y_6+z_6)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{\sqrt{3}}x_6-\tfrac{1}{2\sqrt{3}}y_6-\tfrac{1}{2\sqrt{3}}z_6\right)a\,\hat{\mathbf{y}}+\tfrac{1}{3}(x_6+y_6+z_6)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re V}$$

$$\mathbf{B}_{35} = y_6\,\mathbf{a}_1+z_6\,\mathbf{a}_2+x_6\,\mathbf{a}_3 = \tfrac{1}{2}(-x_6+y_6)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_6-\tfrac{1}{2\sqrt{3}}y_6+\tfrac{1}{\sqrt{3}}z_6\right)a\,\hat{\mathbf{y}}+\tfrac{1}{3}(x_6+y_6+z_6)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re V}$$

$$\mathbf{B}_{36} = \left(\tfrac{1}{2}-z_6\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_6\right)\mathbf{a}_2+\left(\tfrac{1}{2}-x_6\right)\mathbf{a}_3 = \tfrac{1}{2}(x_6-z_6)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2\sqrt{3}}x_6-\tfrac{1}{\sqrt{3}}y_6+\tfrac{1}{2\sqrt{3}}z_6\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-\tfrac{1}{3}x_6-\tfrac{1}{3}y_6-\tfrac{1}{3}z_6\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re V}$$

$$\mathbf{B}_{37} = \left(\tfrac{1}{2}-y_6\right)\mathbf{a}_1+\left(\tfrac{1}{2}-x_6\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_6\right)\mathbf{a}_3 = \tfrac{1}{2}(-y_6+z_6)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{\sqrt{3}}x_6+\tfrac{1}{2\sqrt{3}}y_6+\tfrac{1}{2\sqrt{3}}z_6\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-\tfrac{1}{3}x_6-\tfrac{1}{3}y_6-\tfrac{1}{3}z_6\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re V}$$

$$\mathbf{B}_{38} = \left(\tfrac{1}{2}-x_6\right)\mathbf{a}_1+\left(\tfrac{1}{2}-z_6\right)\mathbf{a}_2+\left(\tfrac{1}{2}-y_6\right)\mathbf{a}_3 = \tfrac{1}{2}(-x_6+y_6)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2\sqrt{3}}x_6+\tfrac{1}{2\sqrt{3}}y_6-\tfrac{1}{\sqrt{3}}z_6\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-\tfrac{1}{3}x_6-\tfrac{1}{3}y_6-\tfrac{1}{3}z_6\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re V}$$

$$\mathbf{B}_{39} = -x_6\,\mathbf{a}_1-y_6\,\mathbf{a}_2-z_6\,\mathbf{a}_3 = \tfrac{1}{2}(-x_6+z_6)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2\sqrt{3}}x_6-\tfrac{1}{\sqrt{3}}y_6+\tfrac{1}{2\sqrt{3}}z_6\right)a\,\hat{\mathbf{y}}-\tfrac{1}{3}(x_6+y_6+z_6)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re V}$$

$$\mathbf{B}_{40} = -z_6\,\mathbf{a}_1-x_6\,\mathbf{a}_2-y_6\,\mathbf{a}_3 = \tfrac{1}{2}(y_6-z_6)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{\sqrt{3}}x_6+\tfrac{1}{2\sqrt{3}}y_6+\tfrac{1}{2\sqrt{3}}z_6\right)a\,\hat{\mathbf{y}}-\tfrac{1}{3}(x_6+y_6+z_6)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re V}$$

$$\mathbf{B}_{41} = -y_6\,\mathbf{a}_1-z_6\,\mathbf{a}_2-x_6\,\mathbf{a}_3 = \tfrac{1}{2}(x_6-y_6)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2\sqrt{3}}x_6+\tfrac{1}{2\sqrt{3}}y_6-\tfrac{1}{\sqrt{3}}z_6\right)a\,\hat{\mathbf{y}}-\tfrac{1}{3}(x_6+y_6+z_6)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re V}$$

$$\mathbf{B}_{42} = \left(\tfrac{1}{2}+z_6\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_6\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_6\right)\mathbf{a}_3 = \tfrac{1}{2}(-x_6+z_6)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_6+\tfrac{1}{\sqrt{3}}y_6-\tfrac{1}{2\sqrt{3}}z_6\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+\tfrac{1}{3}x_6+\tfrac{1}{3}y_6+\tfrac{1}{3}z_6\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re V}$$

$$\mathbf{B}_{43} = \left(\tfrac{1}{2}+y_6\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_6\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_6\right)\mathbf{a}_3 = \tfrac{1}{2}(y_6-z_6)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{\sqrt{3}}x_6-\tfrac{1}{2\sqrt{3}}y_6-\tfrac{1}{2\sqrt{3}}z_6\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+\tfrac{1}{3}x_6+\tfrac{1}{3}y_6+\tfrac{1}{3}z_6\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re V}$$

$$\mathbf{B}_{44} = \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_1+\left(\tfrac{1}{2}+z_6\right)\mathbf{a}_2+\left(\tfrac{1}{2}+y_6\right)\mathbf{a}_3 = \tfrac{1}{2}(x_6-y_6)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_6-\tfrac{1}{2\sqrt{3}}y_6+\tfrac{1}{\sqrt{3}}z_6\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+\tfrac{1}{3}x_6+\tfrac{1}{3}y_6+\tfrac{1}{3}z_6\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re V}$$

$$\mathbf{B}_{45} = x_7\,\mathbf{a}_1+y_7\,\mathbf{a}_2+z_7\,\mathbf{a}_3 = \tfrac{1}{2}(x_7-z_7)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_7+\tfrac{1}{\sqrt{3}}y_7-\tfrac{1}{2\sqrt{3}}z_7\right)a\,\hat{\mathbf{y}}+\tfrac{1}{3}(x_7+y_7+z_7)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re VI}$$

$$\mathbf{B}_{46} = z_7\,\mathbf{a}_1+x_7\,\mathbf{a}_2+y_7\,\mathbf{a}_3 = \tfrac{1}{2}(-y_7+z_7)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{\sqrt{3}}x_7-\tfrac{1}{2\sqrt{3}}y_7-\tfrac{1}{2\sqrt{3}}z_7\right)a\,\hat{\mathbf{y}}+\tfrac{1}{3}(x_7+y_7+z_7)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re VI}$$

$$\mathbf{B}_{47} = y_7\,\mathbf{a}_1+z_7\,\mathbf{a}_2+x_7\,\mathbf{a}_3 = \tfrac{1}{2}(-x_7+y_7)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_7-\tfrac{1}{2\sqrt{3}}y_7+\tfrac{1}{\sqrt{3}}z_7\right)a\,\hat{\mathbf{y}}+\tfrac{1}{3}(x_7+y_7+z_7)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re VI}$$

$$\mathbf{B}_{48} = \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}-x_7\right)\mathbf{a}_3 = \tfrac{1}{2}(x_7-z_7)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2\sqrt{3}}x_7-\tfrac{1}{\sqrt{3}}y_7+\tfrac{1}{2\sqrt{3}}z_7\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-\tfrac{1}{3}x_7-\tfrac{1}{3}y_7-\tfrac{1}{3}z_7\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re VI}$$

$$\mathbf{B}_{49} = \left(\tfrac{1}{2}-y_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}-x_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_7\right)\mathbf{a}_3 = \tfrac{1}{2}(-y_7+z_7)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{\sqrt{3}}x_7+\tfrac{1}{2\sqrt{3}}y_7+\tfrac{1}{2\sqrt{3}}z_7\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-\tfrac{1}{3}x_7-\tfrac{1}{3}y_7-\tfrac{1}{3}z_7\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re VI}$$

$$\mathbf{B}_{50} = \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}-z_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}-y_7\right)\mathbf{a}_3 = \tfrac{1}{2}(-x_7+y_7)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2\sqrt{3}}x_7+\tfrac{1}{2\sqrt{3}}y_7-\tfrac{1}{\sqrt{3}}z_7\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-\tfrac{1}{3}x_7-\tfrac{1}{3}y_7-\tfrac{1}{3}z_7\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re VI}$$

$$\mathbf{B}_{51} = -x_7\,\mathbf{a}_1-y_7\,\mathbf{a}_2-z_7\,\mathbf{a}_3 = \tfrac{1}{2}(-x_7+z_7)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2\sqrt{3}}x_7-\tfrac{1}{\sqrt{3}}y_7+\tfrac{1}{2\sqrt{3}}z_7\right)a\,\hat{\mathbf{y}}-\tfrac{1}{3}(x_7+y_7+z_7)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re VI}$$

$$\mathbf{B}_{52} = -z_7\,\mathbf{a}_1-x_7\,\mathbf{a}_2-y_7\,\mathbf{a}_3 = \tfrac{1}{2}(y_7-z_7)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{\sqrt{3}}x_7+\tfrac{1}{2\sqrt{3}}y_7+\tfrac{1}{2\sqrt{3}}z_7\right)a\,\hat{\mathbf{y}}-\tfrac{1}{3}(x_7+y_7+z_7)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re VI}$$

$$\mathbf{B}_{53} = -y_7\,\mathbf{a}_1-z_7\,\mathbf{a}_2-x_7\,\mathbf{a}_3 = \tfrac{1}{2}(x_7-y_7)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2\sqrt{3}}x_7+\tfrac{1}{2\sqrt{3}}y_7-\tfrac{1}{\sqrt{3}}z_7\right)a\,\hat{\mathbf{y}}-\tfrac{1}{3}(x_7+y_7+z_7)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re VI}$$

$$\mathbf{B}_{54} = \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_7\right)\mathbf{a}_3 = \tfrac{1}{2}(-x_7+z_7)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_7+\tfrac{1}{\sqrt{3}}y_7-\tfrac{1}{2\sqrt{3}}z_7\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+\tfrac{1}{3}x_7+\tfrac{1}{3}y_7+\tfrac{1}{3}z_7\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re VI}$$

$$\mathbf{B}_{55} = \left(\tfrac{1}{2}+y_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 = \tfrac{1}{2}(y_7-z_7)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{\sqrt{3}}x_7-\tfrac{1}{2\sqrt{3}}y_7-\tfrac{1}{2\sqrt{3}}z_7\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+\tfrac{1}{3}x_7+\tfrac{1}{3}y_7+\tfrac{1}{3}z_7\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re VI}$$

$$\mathbf{B}_{56} = \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}+z_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}+y_7\right)\mathbf{a}_3 = \tfrac{1}{2}(x_7-y_7)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_7-\tfrac{1}{2\sqrt{3}}y_7+\tfrac{1}{\sqrt{3}}z_7\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+\tfrac{1}{3}x_7+\tfrac{1}{3}y_7+\tfrac{1}{3}z_7\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Re VI}$$

$$\mathbf{B}_{57} = x_8\,\mathbf{a}_1+y_8\,\mathbf{a}_2+z_8\,\mathbf{a}_3 = \tfrac{1}{2}(x_8-z_8)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_8+\tfrac{1}{\sqrt{3}}y_8-\tfrac{1}{2\sqrt{3}}z_8\right)a\,\hat{\mathbf{y}}+\tfrac{1}{3}(x_8+y_8+z_8)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Zr II}$$

$$\mathbf{B}_{58} = z_8\,\mathbf{a}_1+x_8\,\mathbf{a}_2+y_8\,\mathbf{a}_3 = \tfrac{1}{2}(-y_8+z_8)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{\sqrt{3}}x_8-\tfrac{1}{2\sqrt{3}}y_8-\tfrac{1}{2\sqrt{3}}z_8\right)a\,\hat{\mathbf{y}}+\tfrac{1}{3}(x_8+y_8+z_8)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Zr II}$$

$$\mathbf{B}_{59} = y_8\,\mathbf{a}_1+z_8\,\mathbf{a}_2+x_8\,\mathbf{a}_3 = \tfrac{1}{2}(-x_8+y_8)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_8-\tfrac{1}{2\sqrt{3}}y_8+\tfrac{1}{\sqrt{3}}z_8\right)a\,\hat{\mathbf{y}}+\tfrac{1}{3}(x_8+y_8+z_8)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Zr II}$$

$$\mathbf{B}_{60} = \left(\tfrac{1}{2}-z_8\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}-x_8\right)\mathbf{a}_3 = \tfrac{1}{2}(x_8-z_8)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2\sqrt{3}}x_8-\tfrac{1}{\sqrt{3}}y_8+\tfrac{1}{2\sqrt{3}}z_8\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-\tfrac{1}{3}x_8-\tfrac{1}{3}y_8-\tfrac{1}{3}z_8\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Zr II}$$

$$\mathbf{B}_{61} = \left(\tfrac{1}{2}-y_8\right)\mathbf{a}_1+\left(\tfrac{1}{2}-x_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_8\right)\mathbf{a}_3 = \tfrac{1}{2}(-y_8+z_8)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{\sqrt{3}}x_8+\tfrac{1}{2\sqrt{3}}y_8+\tfrac{1}{2\sqrt{3}}z_8\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-\tfrac{1}{3}x_8-\tfrac{1}{3}y_8-\tfrac{1}{3}z_8\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Zr II}$$

$$\mathbf{B}_{62} = \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_1+\left(\tfrac{1}{2}-z_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}-y_8\right)\mathbf{a}_3 = \tfrac{1}{2}(-x_8+y_8)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2\sqrt{3}}x_8+\tfrac{1}{2\sqrt{3}}y_8-\tfrac{1}{\sqrt{3}}z_8\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-\tfrac{1}{3}x_8-\tfrac{1}{3}y_8-\tfrac{1}{3}z_8\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Zr II}$$

$$\mathbf{B}_{63} = -x_8\,\mathbf{a}_1-y_8\,\mathbf{a}_2-z_8\,\mathbf{a}_3 = \tfrac{1}{2}(-x_8+z_8)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2\sqrt{3}}x_8-\tfrac{1}{\sqrt{3}}y_8+\tfrac{1}{2\sqrt{3}}z_8\right)a\,\hat{\mathbf{y}}-\tfrac{1}{3}(x_8+y_8+z_8)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Zr II}$$

$$\mathbf{B}_{64} = -z_8\,\mathbf{a}_1-x_8\,\mathbf{a}_2-y_8\,\mathbf{a}_3 = \tfrac{1}{2}(y_8-z_8)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{\sqrt{3}}x_8+\tfrac{1}{2\sqrt{3}}y_8+\tfrac{1}{2\sqrt{3}}z_8\right)a\,\hat{\mathbf{y}}-\tfrac{1}{3}(x_8+y_8+z_8)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Zr II}$$

$$\mathbf{B}_{65} = -y_8\,\mathbf{a}_1-z_8\,\mathbf{a}_2-x_8\,\mathbf{a}_3 = \tfrac{1}{2}(x_8-y_8)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2\sqrt{3}}x_8+\tfrac{1}{2\sqrt{3}}y_8-\tfrac{1}{\sqrt{3}}z_8\right)a\,\hat{\mathbf{y}}-\tfrac{1}{3}(x_8+y_8+z_8)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Zr II}$$

$$\mathbf{B}_{66} = \left(\tfrac{1}{2}+z_8\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_8\right)\mathbf{a}_3 = \tfrac{1}{2}(-x_8+z_8)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_8+\tfrac{1}{\sqrt{3}}y_8-\tfrac{1}{2\sqrt{3}}z_8\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+\tfrac{1}{3}x_8+\tfrac{1}{3}y_8+\tfrac{1}{3}z_8\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Zr II}$$

$$\mathbf{B}_{67} = \left(\tfrac{1}{2}+y_8\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_8\right)\mathbf{a}_3 = \tfrac{1}{2}(y_8-z_8)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{\sqrt{3}}x_8-\tfrac{1}{2\sqrt{3}}y_8-\tfrac{1}{2\sqrt{3}}z_8\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+\tfrac{1}{3}x_8+\tfrac{1}{3}y_8+\tfrac{1}{3}z_8\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Zr II}$$

$$\mathbf{B}_{68} = \left(\tfrac{1}{2}+x_8\right)\mathbf{a}_1+\left(\tfrac{1}{2}+z_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}+y_8\right)\mathbf{a}_3 = \tfrac{1}{2}(x_8-y_8)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_8-\tfrac{1}{2\sqrt{3}}y_8+\tfrac{1}{\sqrt{3}}z_8\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+\tfrac{1}{3}x_8+\tfrac{1}{3}y_8+\tfrac{1}{3}z_8\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Zr II}$$

$$\mathbf{B}_{69} = x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 = \tfrac{1}{2}\,(x_9 - z_9)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_9 + \tfrac{1}{\sqrt{3}}y_9 - \tfrac{1}{2\sqrt{3}}z_9\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,(x_9 + y_9 + z_9)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Zr III}$$

$$\mathbf{B}_{70} = z_9\,\mathbf{a}_1 + x_9\,\mathbf{a}_2 + y_9\,\mathbf{a}_3 = \tfrac{1}{2}\,(-y_9 + z_9)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{\sqrt{3}}x_9 - \tfrac{1}{2\sqrt{3}}y_9 - \tfrac{1}{2\sqrt{3}}z_9\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,(x_9 + y_9 + z_9)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Zr III}$$

$$\mathbf{B}_{71} = y_9\,\mathbf{a}_1 + z_9\,\mathbf{a}_2 + x_9\,\mathbf{a}_3 = \tfrac{1}{2}\,(-x_9 + y_9)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_9 - \tfrac{1}{2\sqrt{3}}y_9 + \tfrac{1}{\sqrt{3}}z_9\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,(x_9 + y_9 + z_9)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Zr III}$$

$$\mathbf{B}_{72} = \left(\tfrac{1}{2} - z_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_9\right)\mathbf{a}_3 = \tfrac{1}{2}\,(x_9 - z_9)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2\sqrt{3}}x_9 - \tfrac{1}{\sqrt{3}}y_9 + \tfrac{1}{2\sqrt{3}}z_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - \tfrac{1}{3}x_9 - \tfrac{1}{3}y_9 - \tfrac{1}{3}z_9\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Zr III}$$

$$\mathbf{B}_{73} = \left(\tfrac{1}{2} - y_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_9\right)\mathbf{a}_3 = \tfrac{1}{2}\,(-y_9 + z_9)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{\sqrt{3}}x_9 + \tfrac{1}{2\sqrt{3}}y_9 + \tfrac{1}{2\sqrt{3}}z_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - \tfrac{1}{3}x_9 - \tfrac{1}{3}y_9 - \tfrac{1}{3}z_9\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Zr III}$$

$$\mathbf{B}_{74} = \left(\tfrac{1}{2} - x_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - z_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - y_9\right)\mathbf{a}_3 = \tfrac{1}{2}\,(-x_9 + y_9)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2\sqrt{3}}x_9 + \tfrac{1}{2\sqrt{3}}y_9 - \tfrac{1}{\sqrt{3}}z_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - \tfrac{1}{3}x_9 - \tfrac{1}{3}y_9 - \tfrac{1}{3}z_9\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Zr III}$$

$$\mathbf{B}_{75} = -x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3 = \tfrac{1}{2}\,(-x_9 + z_9)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2\sqrt{3}}x_9 - \tfrac{1}{\sqrt{3}}y_9 + \tfrac{1}{2\sqrt{3}}z_9\right)a\,\hat{\mathbf{y}} - \tfrac{1}{3}\,(x_9 + y_9 + z_9)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Zr III}$$

$$\mathbf{B}_{76} = -z_9\,\mathbf{a}_1 - x_9\,\mathbf{a}_2 - y_9\,\mathbf{a}_3 = \tfrac{1}{2}\,(y_9 - z_9)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{\sqrt{3}}x_9 + \tfrac{1}{2\sqrt{3}}y_9 + \tfrac{1}{2\sqrt{3}}z_9\right)a\,\hat{\mathbf{y}} - \tfrac{1}{3}\,(x_9 + y_9 + z_9)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Zr III}$$

$$\mathbf{B}_{77} = -y_9\,\mathbf{a}_1 - z_9\,\mathbf{a}_2 - x_9\,\mathbf{a}_3 = \tfrac{1}{2}\,(x_9 - y_9)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2\sqrt{3}}x_9 + \tfrac{1}{2\sqrt{3}}y_9 - \tfrac{1}{\sqrt{3}}z_9\right)a\,\hat{\mathbf{y}} - \tfrac{1}{3}\,(x_9 + y_9 + z_9)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Zr III}$$

$$\mathbf{B}_{78} = \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_9\right)\mathbf{a}_3 = \tfrac{1}{2}\,(-x_9 + z_9)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_9 + \tfrac{1}{\sqrt{3}}y_9 - \tfrac{1}{2\sqrt{3}}z_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + \tfrac{1}{3}x_9 + \tfrac{1}{3}y_9 + \tfrac{1}{3}z_9\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Zr III}$$

$$\mathbf{B}_{79} = \left(\tfrac{1}{2} + y_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_3 = \tfrac{1}{2}\,(y_9 - z_9)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{\sqrt{3}}x_9 - \tfrac{1}{2\sqrt{3}}y_9 - \tfrac{1}{2\sqrt{3}}z_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + \tfrac{1}{3}x_9 + \tfrac{1}{3}y_9 + \tfrac{1}{3}z_9\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Zr III}$$

$$\mathbf{B}_{80} = \left(\tfrac{1}{2} + x_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + y_9\right)\mathbf{a}_3 = \tfrac{1}{2}\,(x_9 - y_9)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_9 - \tfrac{1}{2\sqrt{3}}y_9 + \tfrac{1}{\sqrt{3}}z_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + \tfrac{1}{3}x_9 + \tfrac{1}{3}y_9 + \tfrac{1}{3}z_9\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Zr III}$$

$$\mathbf{B}_{81} = x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 = \tfrac{1}{2}\,(x_{10} - z_{10})\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_{10} + \tfrac{1}{\sqrt{3}}y_{10} - \tfrac{1}{2\sqrt{3}}z_{10}\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,(x_{10} + y_{10} + z_{10})\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Zr IV}$$

$\mathbf{B}_{82}$	=	$z_{10}\,\mathbf{a}_1 + x_{10}\,\mathbf{a}_2 + y_{10}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(-y_{10}+z_{10})\,a\,\hat{\mathbf{x}} + \left(\frac{1}{\sqrt{3}}x_{10} - \frac{1}{2\sqrt{3}}y_{10} - \frac{1}{2\sqrt{3}}z_{10}\right)a\,\hat{\mathbf{y}} + \frac{1}{3}\,(x_{10}+y_{10}+z_{10})\,c\,\hat{\mathbf{z}}$	$(12f)$	Zr IV
$\mathbf{B}_{83}$	=	$y_{10}\,\mathbf{a}_1 + z_{10}\,\mathbf{a}_2 + x_{10}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(-x_{10}+y_{10})\,a\,\hat{\mathbf{x}} + \left(-\frac{1}{2\sqrt{3}}x_{10} - \frac{1}{2\sqrt{3}}y_{10} + \frac{1}{\sqrt{3}}z_{10}\right)a\,\hat{\mathbf{y}} + \frac{1}{3}\,(x_{10}+y_{10}+z_{10})\,c\,\hat{\mathbf{z}}$	$(12f)$	Zr IV
$\mathbf{B}_{84}$	=	$\left(\frac{1}{2}-z_{10}\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_{10}\right)\mathbf{a}_2 + \left(\frac{1}{2}-x_{10}\right)\mathbf{a}_3$	=	$\frac{1}{2}\,(x_{10}-z_{10})\,a\,\hat{\mathbf{x}} + \left(\frac{1}{2\sqrt{3}}x_{10} - \frac{1}{\sqrt{3}}y_{10} + \frac{1}{2\sqrt{3}}z_{10}\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - \frac{1}{3}x_{10} - \frac{1}{3}y_{10} - \frac{1}{3}z_{10}\right)c\,\hat{\mathbf{z}}$	$(12f)$	Zr IV
$\mathbf{B}_{85}$	=	$\left(\frac{1}{2}-y_{10}\right)\mathbf{a}_1 + \left(\frac{1}{2}-x_{10}\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_{10}\right)\mathbf{a}_3$	=	$\frac{1}{2}\,(-y_{10}+z_{10})\,a\,\hat{\mathbf{x}} + \left(-\frac{1}{\sqrt{3}}x_{10} + \frac{1}{2\sqrt{3}}y_{10} + \frac{1}{2\sqrt{3}}z_{10}\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - \frac{1}{3}x_{10} - \frac{1}{3}y_{10} - \frac{1}{3}z_{10}\right)c\,\hat{\mathbf{z}}$	$(12f)$	Zr IV
$\mathbf{B}_{86}$	=	$\left(\frac{1}{2}-x_{10}\right)\mathbf{a}_1 + \left(\frac{1}{2}-z_{10}\right)\mathbf{a}_2 + \left(\frac{1}{2}-y_{10}\right)\mathbf{a}_3$	=	$\frac{1}{2}\,(-x_{10}+y_{10})\,a\,\hat{\mathbf{x}} + \left(\frac{1}{2\sqrt{3}}x_{10} + \frac{1}{2\sqrt{3}}y_{10} - \frac{1}{\sqrt{3}}z_{10}\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - \frac{1}{3}x_{10} - \frac{1}{3}y_{10} - \frac{1}{3}z_{10}\right)c\,\hat{\mathbf{z}}$	$(12f)$	Zr IV
$\mathbf{B}_{87}$	=	$-x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(-x_{10}+z_{10})\,a\,\hat{\mathbf{x}} + \left(\frac{1}{2\sqrt{3}}x_{10} - \frac{1}{\sqrt{3}}y_{10} + \frac{1}{2\sqrt{3}}z_{10}\right)a\,\hat{\mathbf{y}} - \frac{1}{3}\,(x_{10}+y_{10}+z_{10})\,c\,\hat{\mathbf{z}}$	$(12f)$	Zr IV
$\mathbf{B}_{88}$	=	$-z_{10}\,\mathbf{a}_1 - x_{10}\,\mathbf{a}_2 - y_{10}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(y_{10}-z_{10})\,a\,\hat{\mathbf{x}} + \left(-\frac{1}{\sqrt{3}}x_{10} + \frac{1}{2\sqrt{3}}y_{10} + \frac{1}{2\sqrt{3}}z_{10}\right)a\,\hat{\mathbf{y}} - \frac{1}{3}\,(x_{10}+y_{10}+z_{10})\,c\,\hat{\mathbf{z}}$	$(12f)$	Zr IV
$\mathbf{B}_{89}$	=	$-y_{10}\,\mathbf{a}_1 - z_{10}\,\mathbf{a}_2 - x_{10}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_{10}-y_{10})\,a\,\hat{\mathbf{x}} + \left(\frac{1}{2\sqrt{3}}x_{10} + \frac{1}{2\sqrt{3}}y_{10} - \frac{1}{\sqrt{3}}z_{10}\right)a\,\hat{\mathbf{y}} - \frac{1}{3}\,(x_{10}+y_{10}+z_{10})\,c\,\hat{\mathbf{z}}$	$(12f)$	Zr IV
$\mathbf{B}_{90}$	=	$\left(\frac{1}{2}+z_{10}\right)\mathbf{a}_1 + \left(\frac{1}{2}+y_{10}\right)\mathbf{a}_2 + \left(\frac{1}{2}+x_{10}\right)\mathbf{a}_3$	=	$\frac{1}{2}\,(-x_{10}+z_{10})\,a\,\hat{\mathbf{x}} + \left(-\frac{1}{2\sqrt{3}}x_{10} + \frac{1}{\sqrt{3}}y_{10} - \frac{1}{2\sqrt{3}}z_{10}\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + \frac{1}{3}x_{10} + \frac{1}{3}y_{10} + \frac{1}{3}z_{10}\right)c\,\hat{\mathbf{z}}$	$(12f)$	Zr IV
$\mathbf{B}_{91}$	=	$\left(\frac{1}{2}+y_{10}\right)\mathbf{a}_1 + \left(\frac{1}{2}+x_{10}\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_{10}\right)\mathbf{a}_3$	=	$\frac{1}{2}\,(y_{10}-z_{10})\,a\,\hat{\mathbf{x}} + \left(\frac{1}{\sqrt{3}}x_{10} - \frac{1}{2\sqrt{3}}y_{10} - \frac{1}{2\sqrt{3}}z_{10}\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + \frac{1}{3}x_{10} + \frac{1}{3}y_{10} + \frac{1}{3}z_{10}\right)c\,\hat{\mathbf{z}}$	$(12f)$	Zr IV
$\mathbf{B}_{92}$	=	$\left(\frac{1}{2}+x_{10}\right)\mathbf{a}_1 + \left(\frac{1}{2}+z_{10}\right)\mathbf{a}_2 + \left(\frac{1}{2}+y_{10}\right)\mathbf{a}_3$	=	$\frac{1}{2}\,(x_{10}-y_{10})\,a\,\hat{\mathbf{x}} + \left(-\frac{1}{2\sqrt{3}}x_{10} - \frac{1}{2\sqrt{3}}y_{10} + \frac{1}{\sqrt{3}}z_{10}\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + \frac{1}{3}x_{10} + \frac{1}{3}y_{10} + \frac{1}{3}z_{10}\right)c\,\hat{\mathbf{z}}$	$(12f)$	Zr IV

References:

- K. Cenzual, E. Parthé, and R. M. Waterstrat, *$Zr_{21}Re_{25}$, a new rhombohedral structure type containing 12 Å-thick infinite $MgZn_2$(Laves)-type columns*, Acta Crystallogr. C **42**, 261–266 (1986), doi:10.1107/S0108270186096555.

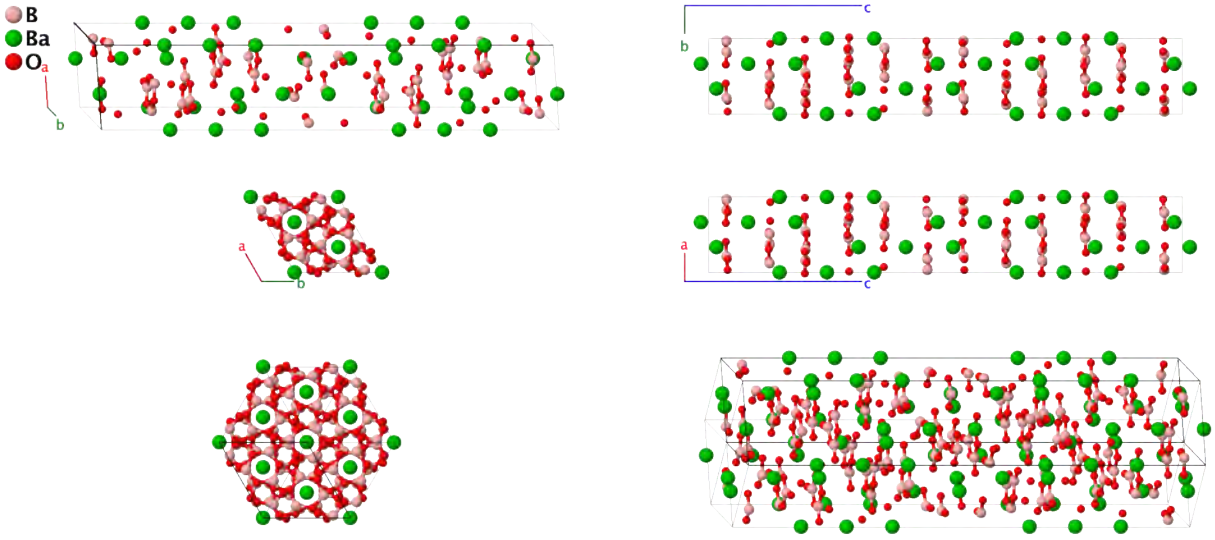
Found in:

- R. Cerný and G. Renaudin, *The intermetallic compound $Mg_{21}Zn_{25}$*, Acta Crystallogr. C **58**, i154–i155 (2002), doi:10.1107/S0108270102018103.

Geometry files:
- CIF: pp. 1743
- POSCAR: pp. 1743

β-BaB_2O_4 (High-Temperature) Structure: A2BC4_hR42_167_f_ac_2f

http://aflow.org/prototype-encyclopedia/A2BC4_hR42_167_f_ac_2f

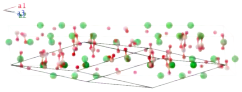


Prototype	:	B_2BaO_4
AFLOW prototype label	:	A2BC4_hR42_167_f_ac_2f
***Strukturbericht* designation**	:	None
Pearson symbol	:	hR42
Space group number	:	167
Space group symbol	:	$R\bar{3}c$
AFLOW prototype command	:	`aflow --proto=A2BC4_hR42_167_f_ac_2f [--hex] --params=`$a, c/a, x_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5$

- This is the high-temperature structure of BaB_2O_4. Cooling to temperatures between 100-400 °C, it transforms into α-BaB_2O_4. The principle difference between the two forms is the lack of inversion symmetry in the low-temperature structure.
- This structure was determined by cooling liquid BaB_2O_4 to room temperature.

Rhombohedral primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = -\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{4}\mathbf{a}_1 + \frac{1}{4}\mathbf{a}_2 + \frac{1}{4}\mathbf{a}_3$	=	$\frac{1}{4}c\,\hat{\mathbf{z}}$	(2*a*)	Ba I
$\mathbf{B}_2$	=	$\frac{3}{4}\mathbf{a}_1 + \frac{3}{4}\mathbf{a}_2 + \frac{3}{4}\mathbf{a}_3$	=	$\frac{3}{4}c\,\hat{\mathbf{z}}$	(2*a*)	Ba I

$$\mathbf{B}_3 = x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + x_2\,\mathbf{a}_3 = x_2 c\,\hat{\mathbf{z}} \quad (4c) \quad \text{Ba II}$$

$$\mathbf{B}_4 = \left(\tfrac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_2\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_2\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_2\right) c\,\hat{\mathbf{z}} \quad (4c) \quad \text{Ba II}$$

$$\mathbf{B}_5 = -x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 - x_2\,\mathbf{a}_3 = -x_2 c\,\hat{\mathbf{z}} \quad (4c) \quad \text{Ba II}$$

$$\mathbf{B}_6 = \left(\tfrac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_2\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_2\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_2\right) c\,\hat{\mathbf{z}} \quad (4c) \quad \text{Ba II}$$

$$\mathbf{B}_7 = x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3 = \tfrac{1}{2}\left(x_3 - z_3\right) a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}} x_3 + \tfrac{1}{\sqrt{3}} y_3 - \tfrac{1}{2\sqrt{3}} z_3\right) a\,\hat{\mathbf{y}} + \tfrac{1}{3}\left(x_3 + y_3 + z_3\right) c\,\hat{\mathbf{z}} \quad (12f) \quad \text{B}$$

$$\mathbf{B}_8 = z_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + y_3\,\mathbf{a}_3 = \tfrac{1}{2}\left(-y_3 + z_3\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{\sqrt{3}} x_3 - \tfrac{1}{2\sqrt{3}} y_3 - \tfrac{1}{2\sqrt{3}} z_3\right) a\,\hat{\mathbf{y}} + \tfrac{1}{3}\left(x_3 + y_3 + z_3\right) c\,\hat{\mathbf{z}} \quad (12f) \quad \text{B}$$

$$\mathbf{B}_9 = y_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3 = \tfrac{1}{2}\left(-x_3 + y_3\right) a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}} x_3 - \tfrac{1}{2\sqrt{3}} y_3 + \tfrac{1}{\sqrt{3}} z_3\right) a\,\hat{\mathbf{y}} + \tfrac{1}{3}\left(x_3 + y_3 + z_3\right) c\,\hat{\mathbf{z}} \quad (12f) \quad \text{B}$$

$$\mathbf{B}_{10} = \left(\tfrac{1}{2} - z_3\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_3\right)\mathbf{a}_3 = \tfrac{1}{2}\left(x_3 - z_3\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2\sqrt{3}} x_3 - \tfrac{1}{\sqrt{3}} y_3 + \tfrac{1}{2\sqrt{3}} z_3\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - \tfrac{1}{3} x_3 - \tfrac{1}{3} y_3 - \tfrac{1}{3} z_3\right) c\,\hat{\mathbf{z}} \quad (12f) \quad \text{B}$$

$$\mathbf{B}_{11} = \left(\tfrac{1}{2} - y_3\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_3\right)\mathbf{a}_3 = \tfrac{1}{2}\left(-y_3 + z_3\right) a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{\sqrt{3}} x_3 + \tfrac{1}{2\sqrt{3}} y_3 + \tfrac{1}{2\sqrt{3}} z_3\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - \tfrac{1}{3} x_3 - \tfrac{1}{3} y_3 - \tfrac{1}{3} z_3\right) c\,\hat{\mathbf{z}} \quad (12f) \quad \text{B}$$

$$\mathbf{B}_{12} = \left(\tfrac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - z_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - y_3\right)\mathbf{a}_3 = \tfrac{1}{2}\left(-x_3 + y_3\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2\sqrt{3}} x_3 + \tfrac{1}{2\sqrt{3}} y_3 - \tfrac{1}{\sqrt{3}} z_3\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - \tfrac{1}{3} x_3 - \tfrac{1}{3} y_3 - \tfrac{1}{3} z_3\right) c\,\hat{\mathbf{z}} \quad (12f) \quad \text{B}$$

$$\mathbf{B}_{13} = -x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3 = \tfrac{1}{2}\left(-x_3 + z_3\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2\sqrt{3}} x_3 - \tfrac{1}{\sqrt{3}} y_3 + \tfrac{1}{2\sqrt{3}} z_3\right) a\,\hat{\mathbf{y}} - \tfrac{1}{3}\left(x_3 + y_3 + z_3\right) c\,\hat{\mathbf{z}} \quad (12f) \quad \text{B}$$

$$\mathbf{B}_{14} = -z_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - y_3\,\mathbf{a}_3 = \tfrac{1}{2}\left(y_3 - z_3\right) a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{\sqrt{3}} x_3 + \tfrac{1}{2\sqrt{3}} y_3 + \tfrac{1}{2\sqrt{3}} z_3\right) a\,\hat{\mathbf{y}} - \tfrac{1}{3}\left(x_3 + y_3 + z_3\right) c\,\hat{\mathbf{z}} \quad (12f) \quad \text{B}$$

$$\mathbf{B}_{15} = -y_3\,\mathbf{a}_1 - z_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3 = \tfrac{1}{2}\left(x_3 - y_3\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2\sqrt{3}} x_3 + \tfrac{1}{2\sqrt{3}} y_3 - \tfrac{1}{\sqrt{3}} z_3\right) a\,\hat{\mathbf{y}} - \tfrac{1}{3}\left(x_3 + y_3 + z_3\right) c\,\hat{\mathbf{z}} \quad (12f) \quad \text{B}$$

$$\mathbf{B}_{16} = \left(\tfrac{1}{2} + z_3\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_3\right)\mathbf{a}_3 = \tfrac{1}{2}\left(-x_3 + z_3\right) a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}} x_3 + \tfrac{1}{\sqrt{3}} y_3 - \tfrac{1}{2\sqrt{3}} z_3\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + \tfrac{1}{3} x_3 + \tfrac{1}{3} y_3 + \tfrac{1}{3} z_3\right) c\,\hat{\mathbf{z}} \quad (12f) \quad \text{B}$$

$$\mathbf{B}_{17} = \left(\tfrac{1}{2} + y_3\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_3\right)\mathbf{a}_3 = \tfrac{1}{2}\left(y_3 - z_3\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{\sqrt{3}} x_3 - \tfrac{1}{2\sqrt{3}} y_3 - \tfrac{1}{2\sqrt{3}} z_3\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + \tfrac{1}{3} x_3 + \tfrac{1}{3} y_3 + \tfrac{1}{3} z_3\right) c\,\hat{\mathbf{z}} \quad (12f) \quad \text{B}$$

$$\mathbf{B}_{18} = \left(\tfrac{1}{2}+x_3\right)\mathbf{a}_1+\left(\tfrac{1}{2}+z_3\right)\mathbf{a}_2+\left(\tfrac{1}{2}+y_3\right)\mathbf{a}_3 = \tfrac{1}{2}(x_3-y_3)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_3-\tfrac{1}{2\sqrt{3}}y_3+\tfrac{1}{\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+\tfrac{1}{3}x_3+\tfrac{1}{3}y_3+\tfrac{1}{3}z_3\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{B}$$

$$\mathbf{B}_{19} = x_4\,\mathbf{a}_1+y_4\,\mathbf{a}_2+z_4\,\mathbf{a}_3 = \tfrac{1}{2}(x_4-z_4)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_4+\tfrac{1}{\sqrt{3}}y_4-\tfrac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}}+\tfrac{1}{3}(x_4+y_4+z_4)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{O I}$$

$$\mathbf{B}_{20} = z_4\,\mathbf{a}_1+x_4\,\mathbf{a}_2+y_4\,\mathbf{a}_3 = \tfrac{1}{2}(-y_4+z_4)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{\sqrt{3}}x_4-\tfrac{1}{2\sqrt{3}}y_4-\tfrac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}}+\tfrac{1}{3}(x_4+y_4+z_4)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{O I}$$

$$\mathbf{B}_{21} = y_4\,\mathbf{a}_1+z_4\,\mathbf{a}_2+x_4\,\mathbf{a}_3 = \tfrac{1}{2}(-x_4+y_4)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_4-\tfrac{1}{2\sqrt{3}}y_4+\tfrac{1}{\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}}+\tfrac{1}{3}(x_4+y_4+z_4)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{O I}$$

$$\mathbf{B}_{22} = \left(\tfrac{1}{2}-z_4\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_4\right)\mathbf{a}_2+\left(\tfrac{1}{2}-x_4\right)\mathbf{a}_3 = \tfrac{1}{2}(x_4-z_4)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2\sqrt{3}}x_4-\tfrac{1}{\sqrt{3}}y_4+\tfrac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-\tfrac{1}{3}x_4-\tfrac{1}{3}y_4-\tfrac{1}{3}z_4\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{O I}$$

$$\mathbf{B}_{23} = \left(\tfrac{1}{2}-y_4\right)\mathbf{a}_1+\left(\tfrac{1}{2}-x_4\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_4\right)\mathbf{a}_3 = \tfrac{1}{2}(-y_4+z_4)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{\sqrt{3}}x_4+\tfrac{1}{2\sqrt{3}}y_4+\tfrac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-\tfrac{1}{3}x_4-\tfrac{1}{3}y_4-\tfrac{1}{3}z_4\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{O I}$$

$$\mathbf{B}_{24} = \left(\tfrac{1}{2}-x_4\right)\mathbf{a}_1+\left(\tfrac{1}{2}-z_4\right)\mathbf{a}_2+\left(\tfrac{1}{2}-y_4\right)\mathbf{a}_3 = \tfrac{1}{2}(-x_4+y_4)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2\sqrt{3}}x_4+\tfrac{1}{2\sqrt{3}}y_4-\tfrac{1}{\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-\tfrac{1}{3}x_4-\tfrac{1}{3}y_4-\tfrac{1}{3}z_4\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{O I}$$

$$\mathbf{B}_{25} = -x_4\,\mathbf{a}_1-y_4\,\mathbf{a}_2-z_4\,\mathbf{a}_3 = \tfrac{1}{2}(-x_4+z_4)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2\sqrt{3}}x_4-\tfrac{1}{\sqrt{3}}y_4+\tfrac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}}-\tfrac{1}{3}(x_4+y_4+z_4)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{O I}$$

$$\mathbf{B}_{26} = -z_4\,\mathbf{a}_1-x_4\,\mathbf{a}_2-y_4\,\mathbf{a}_3 = \tfrac{1}{2}(y_4-z_4)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{\sqrt{3}}x_4+\tfrac{1}{2\sqrt{3}}y_4+\tfrac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}}-\tfrac{1}{3}(x_4+y_4+z_4)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{O I}$$

$$\mathbf{B}_{27} = -y_4\,\mathbf{a}_1-z_4\,\mathbf{a}_2-x_4\,\mathbf{a}_3 = \tfrac{1}{2}(x_4-y_4)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2\sqrt{3}}x_4+\tfrac{1}{2\sqrt{3}}y_4-\tfrac{1}{\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}}-\tfrac{1}{3}(x_4+y_4+z_4)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{O I}$$

$$\mathbf{B}_{28} = \left(\tfrac{1}{2}+z_4\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_4\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_4\right)\mathbf{a}_3 = \tfrac{1}{2}(-x_4+z_4)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_4+\tfrac{1}{\sqrt{3}}y_4-\tfrac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+\tfrac{1}{3}x_4+\tfrac{1}{3}y_4+\tfrac{1}{3}z_4\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{O I}$$

$$\mathbf{B}_{29} = \left(\tfrac{1}{2}+y_4\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_4\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_4\right)\mathbf{a}_3 = \tfrac{1}{2}(y_4-z_4)\,a\,\hat{\mathbf{x}}+\left(\tfrac{1}{\sqrt{3}}x_4-\tfrac{1}{2\sqrt{3}}y_4-\tfrac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+\tfrac{1}{3}x_4+\tfrac{1}{3}y_4+\tfrac{1}{3}z_4\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{O I}$$

$$\mathbf{B}_{30} = \left(\tfrac{1}{2}+x_4\right)\mathbf{a}_1+\left(\tfrac{1}{2}+z_4\right)\mathbf{a}_2+\left(\tfrac{1}{2}+y_4\right)\mathbf{a}_3 = \tfrac{1}{2}(x_4-y_4)\,a\,\hat{\mathbf{x}}+\left(-\tfrac{1}{2\sqrt{3}}x_4-\tfrac{1}{2\sqrt{3}}y_4+\tfrac{1}{\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+\tfrac{1}{3}x_4+\tfrac{1}{3}y_4+\tfrac{1}{3}z_4\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{O I}$$

$$\mathbf{B}_{31} = x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3 = \tfrac{1}{2}\,(x_5 - z_5)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_5 + \tfrac{1}{\sqrt{3}}y_5 - \tfrac{1}{2\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,(x_5 + y_5 + z_5)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{O II}$$

$$\mathbf{B}_{32} = z_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + y_5\,\mathbf{a}_3 = \tfrac{1}{2}\,(-y_5 + z_5)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{\sqrt{3}}x_5 - \tfrac{1}{2\sqrt{3}}y_5 - \tfrac{1}{2\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,(x_5 + y_5 + z_5)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{O II}$$

$$\mathbf{B}_{33} = y_5\,\mathbf{a}_1 + z_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3 = \tfrac{1}{2}\,(-x_5 + y_5)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_5 - \tfrac{1}{2\sqrt{3}}y_5 + \tfrac{1}{\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,(x_5 + y_5 + z_5)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{O II}$$

$$\mathbf{B}_{34} = \left(\tfrac{1}{2} - z_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_5\right)\mathbf{a}_3 = \tfrac{1}{2}\,(x_5 - z_5)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2\sqrt{3}}x_5 - \tfrac{1}{\sqrt{3}}y_5 + \tfrac{1}{2\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - \tfrac{1}{3}x_5 - \tfrac{1}{3}y_5 - \tfrac{1}{3}z_5\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{O II}$$

$$\mathbf{B}_{35} = \left(\tfrac{1}{2} - y_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_5\right)\mathbf{a}_3 = \tfrac{1}{2}\,(-y_5 + z_5)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{\sqrt{3}}x_5 + \tfrac{1}{2\sqrt{3}}y_5 + \tfrac{1}{2\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - \tfrac{1}{3}x_5 - \tfrac{1}{3}y_5 - \tfrac{1}{3}z_5\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{O II}$$

$$\mathbf{B}_{36} = \left(\tfrac{1}{2} - x_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - z_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - y_5\right)\mathbf{a}_3 = \tfrac{1}{2}\,(-x_5 + y_5)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2\sqrt{3}}x_5 + \tfrac{1}{2\sqrt{3}}y_5 - \tfrac{1}{\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - \tfrac{1}{3}x_5 - \tfrac{1}{3}y_5 - \tfrac{1}{3}z_5\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{O II}$$

$$\mathbf{B}_{37} = -x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3 = \tfrac{1}{2}\,(-x_5 + z_5)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2\sqrt{3}}x_5 - \tfrac{1}{\sqrt{3}}y_5 + \tfrac{1}{2\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}} - \tfrac{1}{3}\,(x_5 + y_5 + z_5)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{O II}$$

$$\mathbf{B}_{38} = -z_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - y_5\,\mathbf{a}_3 = \tfrac{1}{2}\,(y_5 - z_5)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{\sqrt{3}}x_5 + \tfrac{1}{2\sqrt{3}}y_5 + \tfrac{1}{2\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}} - \tfrac{1}{3}\,(x_5 + y_5 + z_5)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{O II}$$

$$\mathbf{B}_{39} = -y_5\,\mathbf{a}_1 - z_5\,\mathbf{a}_2 - x_5\,\mathbf{a}_3 = \tfrac{1}{2}\,(x_5 - y_5)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2\sqrt{3}}x_5 + \tfrac{1}{2\sqrt{3}}y_5 - \tfrac{1}{\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}} - \tfrac{1}{3}\,(x_5 + y_5 + z_5)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{O II}$$

$$\mathbf{B}_{40} = \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_5\right)\mathbf{a}_3 = \tfrac{1}{2}\,(-x_5 + z_5)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_5 + \tfrac{1}{\sqrt{3}}y_5 - \tfrac{1}{2\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + \tfrac{1}{3}x_5 + \tfrac{1}{3}y_5 + \tfrac{1}{3}z_5\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{O II}$$

$$\mathbf{B}_{41} = \left(\tfrac{1}{2} + y_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_3 = \tfrac{1}{2}\,(y_5 - z_5)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{\sqrt{3}}x_5 - \tfrac{1}{2\sqrt{3}}y_5 - \tfrac{1}{2\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + \tfrac{1}{3}x_5 + \tfrac{1}{3}y_5 + \tfrac{1}{3}z_5\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{O II}$$

$$\mathbf{B}_{42} = \left(\tfrac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + y_5\right)\mathbf{a}_3 = \tfrac{1}{2}\,(x_5 - y_5)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_5 - \tfrac{1}{2\sqrt{3}}y_5 + \tfrac{1}{\sqrt{3}}z_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + \tfrac{1}{3}x_5 + \tfrac{1}{3}y_5 + \tfrac{1}{3}z_5\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{O II}$$

References:

- A. D. Mighell, A. Perloff, and S. Block, *The crystal structure of the high temperature form of barium borate, $BaO{\cdot}B_2O_3$*, Acta Cryst. **20**, 819–823 (1966), doi:10.1107/S0365110X66001920.

Found in:

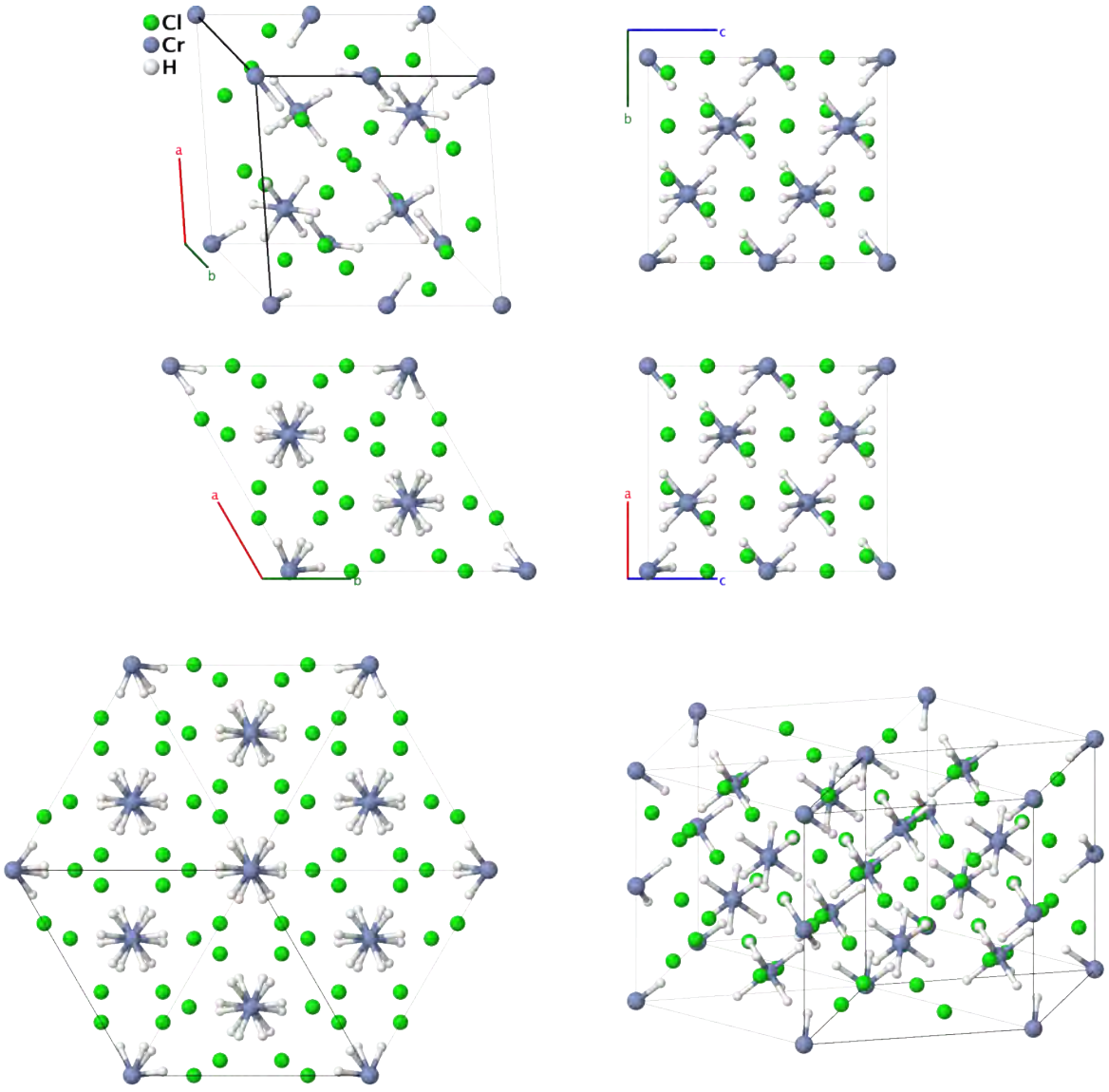
- R. Fröhlich, *Crystal Structure of the low-temperature form of* BaB_2O_4, Zeitschrift für Kristallographie - Crystalline Materials **168**, 109–112 (1984), doi:10.1524/zkri.1984.168.14.109.

Geometry files:

- CIF: pp. 1744
- POSCAR: pp. 1744

$CrCl_3(H_2O)_6$ ($J2_2$) Structure: A3BC6_hR20_167_e_b_f

http://aflow.org/prototype-encyclopedia/A3BC6_hR20_167_e_b_f



Prototype	:	$Cl_3Cr(H_2O)_6$
AFLOW prototype label	:	A3BC6_hR20_167_e_b_f
***Strukturbericht* designation**	:	$J2_2$
Pearson symbol	:	hR20
Space group number	:	167
Space group symbol	:	$R\bar{3}c$
AFLOW prototype command	:	`aflow --proto=A3BC6_hR20_167_e_b_f [--hex] --params=`$a, c/a, x_2, x_3, y_3, z_3$

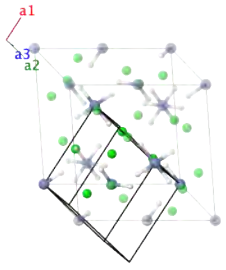
Other compounds with this structure

- $AlCl_3(H_2O)_6$

- The positions of the hydrogen atoms in the water molecules were not determined, so we only provide the positions of the oxygen atoms (labeled as H_2O).

Rhombohedral primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2\sqrt{3}}\,a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,c\,\hat{\mathbf{z}}\\ \mathbf{a}_2 &= \tfrac{1}{\sqrt{3}}\,a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,c\,\hat{\mathbf{z}}\\ \mathbf{a}_3 &= -\tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2\sqrt{3}}\,a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2b)$	Cr
$\mathbf{B}_2$	=	$\tfrac{1}{2}\,\mathbf{a}_1 + \tfrac{1}{2}\,\mathbf{a}_2 + \tfrac{1}{2}\,\mathbf{a}_3$	=	$\tfrac{1}{2}c\,\hat{\mathbf{z}}$	$(2b)$	Cr
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_2\right)\mathbf{a}_2 + \tfrac{1}{4}\,\mathbf{a}_3$	=	$\left(-\tfrac{1}{8} + \tfrac{1}{2}x_2\right)a\,\hat{\mathbf{x}} + \left(\tfrac{\sqrt{3}}{8} - \tfrac{\sqrt{3}}{2}x_2\right)a\,\hat{\mathbf{y}} + \tfrac{1}{4}c\,\hat{\mathbf{z}}$	$(6e)$	Cl
$\mathbf{B}_4$	=	$\tfrac{1}{4}\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_2\right)\mathbf{a}_3$	=	$\left(-\tfrac{1}{8} + \tfrac{1}{2}x_2\right)a\,\hat{\mathbf{x}} + \left(-\tfrac{\sqrt{3}}{8} + \tfrac{\sqrt{3}}{2}x_2\right)a\,\hat{\mathbf{y}} + \tfrac{1}{4}c\,\hat{\mathbf{z}}$	$(6e)$	Cl
$\mathbf{B}_5$	=	$\left(\tfrac{1}{2} - x_2\right)\mathbf{a}_1 + \tfrac{1}{4}\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$\left(\tfrac{1}{4} - x_2\right)a\,\hat{\mathbf{x}} + \tfrac{1}{4}c\,\hat{\mathbf{z}}$	$(6e)$	Cl
$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_2\right)\mathbf{a}_2 + \tfrac{3}{4}\,\mathbf{a}_3$	=	$-\left(\tfrac{1}{2}x_2 + \tfrac{3}{8}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{8\sqrt{3}} + \tfrac{\sqrt{3}}{2}x_2\right)a\,\hat{\mathbf{y}} + \tfrac{5}{12}c\,\hat{\mathbf{z}}$	$(6e)$	Cl
$\mathbf{B}_7$	=	$\tfrac{3}{4}\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_2\right)\mathbf{a}_3$	=	$\left(\tfrac{1}{8} - \tfrac{1}{2}x_2\right)a\,\hat{\mathbf{x}} - \left(\tfrac{\sqrt{3}}{2}x_2 + \tfrac{5}{8\sqrt{3}}\right)a\,\hat{\mathbf{y}} + \tfrac{5}{12}c\,\hat{\mathbf{z}}$	$(6e)$	Cl
$\mathbf{B}_8$	=	$\left(\tfrac{1}{2} + x_2\right)\mathbf{a}_1 + \tfrac{3}{4}\,\mathbf{a}_2 - x_2\,\mathbf{a}_3$	=	$\left(\tfrac{1}{4} + x_2\right)a\,\hat{\mathbf{x}} + \tfrac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \tfrac{5}{12}c\,\hat{\mathbf{z}}$	$(6e)$	Cl
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\tfrac{1}{2}\,(x_3 - z_3)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_3 + \tfrac{1}{\sqrt{3}}y_3 - \tfrac{1}{2\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,(x_3 + y_3 + z_3)\,c\,\hat{\mathbf{z}}$	$(12f)$	H_2O
$\mathbf{B}_{10}$	=	$z_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + y_3\,\mathbf{a}_3$	=	$\tfrac{1}{2}\,(-y_3 + z_3)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{\sqrt{3}}x_3 - \tfrac{1}{2\sqrt{3}}y_3 - \tfrac{1}{2\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,(x_3 + y_3 + z_3)\,c\,\hat{\mathbf{z}}$	$(12f)$	H_2O
$\mathbf{B}_{11}$	=	$y_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$\tfrac{1}{2}\,(-x_3 + y_3)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_3 - \tfrac{1}{2\sqrt{3}}y_3 + \tfrac{1}{\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,(x_3 + y_3 + z_3)\,c\,\hat{\mathbf{z}}$	$(12f)$	H_2O
$\mathbf{B}_{12}$	=	$\left(\tfrac{1}{2} - z_3\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_3\right)\mathbf{a}_3$	=	$\tfrac{1}{2}\,(x_3 - z_3)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2\sqrt{3}}x_3 - \tfrac{1}{\sqrt{3}}y_3 + \tfrac{1}{2\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - \tfrac{1}{3}x_3 - \tfrac{1}{3}y_3 - \tfrac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	$(12f)$	H_2O

$\mathbf{B}_{13}$	=	$\left(\frac{1}{2}-y_3\right)\mathbf{a}_1+\left(\frac{1}{2}-x_3\right)\mathbf{a}_2+\left(\frac{1}{2}-z_3\right)\mathbf{a}_3$	=	$\frac{1}{2}(-y_3+z_3)\,a\,\hat{\mathbf{x}}+\left(-\frac{1}{\sqrt{3}}x_3+\frac{1}{2\sqrt{3}}y_3+\frac{1}{2\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-\frac{1}{3}x_3-\frac{1}{3}y_3-\frac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	$(12f)$	H_2O
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2}-x_3\right)\mathbf{a}_1+\left(\frac{1}{2}-z_3\right)\mathbf{a}_2+\left(\frac{1}{2}-y_3\right)\mathbf{a}_3$	=	$\frac{1}{2}(-x_3+y_3)\,a\,\hat{\mathbf{x}}+\left(\frac{1}{2\sqrt{3}}x_3+\frac{1}{2\sqrt{3}}y_3-\frac{1}{\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-\frac{1}{3}x_3-\frac{1}{3}y_3-\frac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	$(12f)$	H_2O
$\mathbf{B}_{15}$	=	$-x_3\,\mathbf{a}_1-y_3\,\mathbf{a}_2-z_3\,\mathbf{a}_3$	=	$\frac{1}{2}(-x_3+z_3)\,a\,\hat{\mathbf{x}}+\left(\frac{1}{2\sqrt{3}}x_3-\frac{1}{\sqrt{3}}y_3+\frac{1}{2\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}}-\frac{1}{3}(x_3+y_3+z_3)\,c\,\hat{\mathbf{z}}$	$(12f)$	H_2O
$\mathbf{B}_{16}$	=	$-z_3\,\mathbf{a}_1-x_3\,\mathbf{a}_2-y_3\,\mathbf{a}_3$	=	$\frac{1}{2}(y_3-z_3)\,a\,\hat{\mathbf{x}}+\left(-\frac{1}{\sqrt{3}}x_3+\frac{1}{2\sqrt{3}}y_3+\frac{1}{2\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}}-\frac{1}{3}(x_3+y_3+z_3)\,c\,\hat{\mathbf{z}}$	$(12f)$	H_2O
$\mathbf{B}_{17}$	=	$-y_3\,\mathbf{a}_1-z_3\,\mathbf{a}_2-x_3\,\mathbf{a}_3$	=	$\frac{1}{2}(x_3-y_3)\,a\,\hat{\mathbf{x}}+\left(\frac{1}{2\sqrt{3}}x_3+\frac{1}{2\sqrt{3}}y_3-\frac{1}{\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}}-\frac{1}{3}(x_3+y_3+z_3)\,c\,\hat{\mathbf{z}}$	$(12f)$	H_2O
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2}+z_3\right)\mathbf{a}_1+\left(\frac{1}{2}+y_3\right)\mathbf{a}_2+\left(\frac{1}{2}+x_3\right)\mathbf{a}_3$	=	$\frac{1}{2}(-x_3+z_3)\,a\,\hat{\mathbf{x}}+\left(-\frac{1}{2\sqrt{3}}x_3+\frac{1}{\sqrt{3}}y_3-\frac{1}{2\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+\frac{1}{3}x_3+\frac{1}{3}y_3+\frac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	$(12f)$	H_2O
$\mathbf{B}_{19}$	=	$\left(\frac{1}{2}+y_3\right)\mathbf{a}_1+\left(\frac{1}{2}+x_3\right)\mathbf{a}_2+\left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	=	$\frac{1}{2}(y_3-z_3)\,a\,\hat{\mathbf{x}}+\left(\frac{1}{\sqrt{3}}x_3-\frac{1}{2\sqrt{3}}y_3-\frac{1}{2\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+\frac{1}{3}x_3+\frac{1}{3}y_3+\frac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	$(12f)$	H_2O
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2}+x_3\right)\mathbf{a}_1+\left(\frac{1}{2}+z_3\right)\mathbf{a}_2+\left(\frac{1}{2}+y_3\right)\mathbf{a}_3$	=	$\frac{1}{2}(x_3-y_3)\,a\,\hat{\mathbf{x}}+\left(-\frac{1}{2\sqrt{3}}x_3-\frac{1}{2\sqrt{3}}y_3+\frac{1}{\sqrt{3}}z_3\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+\frac{1}{3}x_3+\frac{1}{3}y_3+\frac{1}{3}z_3\right)c\,\hat{\mathbf{z}}$	$(12f)$	H_2O

References:

- K. R. Andress and C. Carpenter, *Kristallhydrate. II. Die Struktur von Chromchlorid- und Aluminiumchloridhexahydrat*, Zeitschrift für Kristallographie - Crystalline Materials **87**, 446–463 (1934), doi:10.1524/zkri.1934.87.1.446.

Found in:

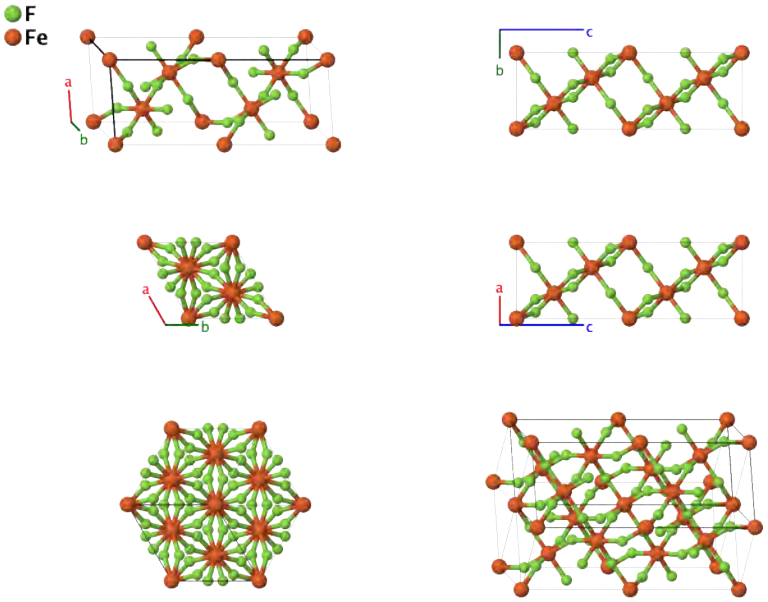
- C. Gottfried and F. Schossberger, eds., *Strukturbericht Band III 1933-1935* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1745
- POSCAR: pp. 1745

FeF_3 ($D0_{12}$) Structure: A3B_hR8_167_e_b

http://aflow.org/prototype-encyclopedia/A3B_hR8_167_e_b



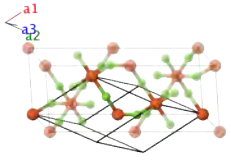
Prototype	:	F_3Fe
AFLOW prototype label	:	A3B_hR8_167_e_b
***Strukturbericht* designation**	:	$D0_{12}$
Pearson symbol	:	hR8
Space group number	:	167
Space group symbol	:	$R\bar{3}c$
AFLOW prototype command	:	`aflow --proto=A3B_hR8_167_e_b [--hex] --params=`$a, c/a, x_2$

Other compounds with this structure

- CoF_3, RuF_3, RhF_3, PdF_3, IrF_3, α-AlF_3, and AlH_3

Rhombohedral primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = -\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($2b$)	Fe
$\mathbf{B}_2$	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	($2b$)	Fe
$\mathbf{B}_3$	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(-\frac{1}{8} + \frac{1}{2}x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{\sqrt{3}}{8} - \frac{\sqrt{3}}{2}x_2\right)a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($6e$)	F

$$\mathbf{B}_4 = \tfrac{1}{4}\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_2\right)\mathbf{a}_3 = \left(-\tfrac{1}{8} + \tfrac{1}{2}x_2\right)a\,\hat{\mathbf{x}} + \left(-\tfrac{\sqrt{3}}{8} + \tfrac{\sqrt{3}}{2}x_2\right)a\,\hat{\mathbf{y}} + \tfrac{1}{4}c\,\hat{\mathbf{z}} \quad (6e) \quad \text{F}$$

$$\mathbf{B}_5 = \left(\tfrac{1}{2} - x_2\right)\mathbf{a}_1 + \tfrac{1}{4}\,\mathbf{a}_2 + x_2\,\mathbf{a}_3 = \left(\tfrac{1}{4} - x_2\right)a\,\hat{\mathbf{x}} + \tfrac{1}{4}c\,\hat{\mathbf{z}} \quad (6e) \quad \text{F}$$

$$\mathbf{B}_6 = -x_2\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_2\right)\mathbf{a}_2 + \tfrac{3}{4}\,\mathbf{a}_3 = -\left(\tfrac{1}{2}x_2 + \tfrac{3}{8}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{8\sqrt{3}} + \tfrac{\sqrt{3}}{2}x_2\right)a\,\hat{\mathbf{y}} + \tfrac{5}{12}c\,\hat{\mathbf{z}} \quad (6e) \quad \text{F}$$

$$\mathbf{B}_7 = \tfrac{3}{4}\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_2\right)\mathbf{a}_3 = \left(\tfrac{1}{8} - \tfrac{1}{2}x_2\right)a\,\hat{\mathbf{x}} - \left(\tfrac{\sqrt{3}}{2}x_2 + \tfrac{5}{8\sqrt{3}}\right)a\,\hat{\mathbf{y}} + \tfrac{5}{12}c\,\hat{\mathbf{z}} \quad (6e) \quad \text{F}$$

$$\mathbf{B}_8 = \left(\tfrac{1}{2} + x_2\right)\mathbf{a}_1 + \tfrac{3}{4}\,\mathbf{a}_2 - x_2\,\mathbf{a}_3 = \left(\tfrac{1}{4} + x_2\right)a\,\hat{\mathbf{x}} + \tfrac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \tfrac{5}{12}c\,\hat{\mathbf{z}} \quad (6e) \quad \text{F}$$

References:

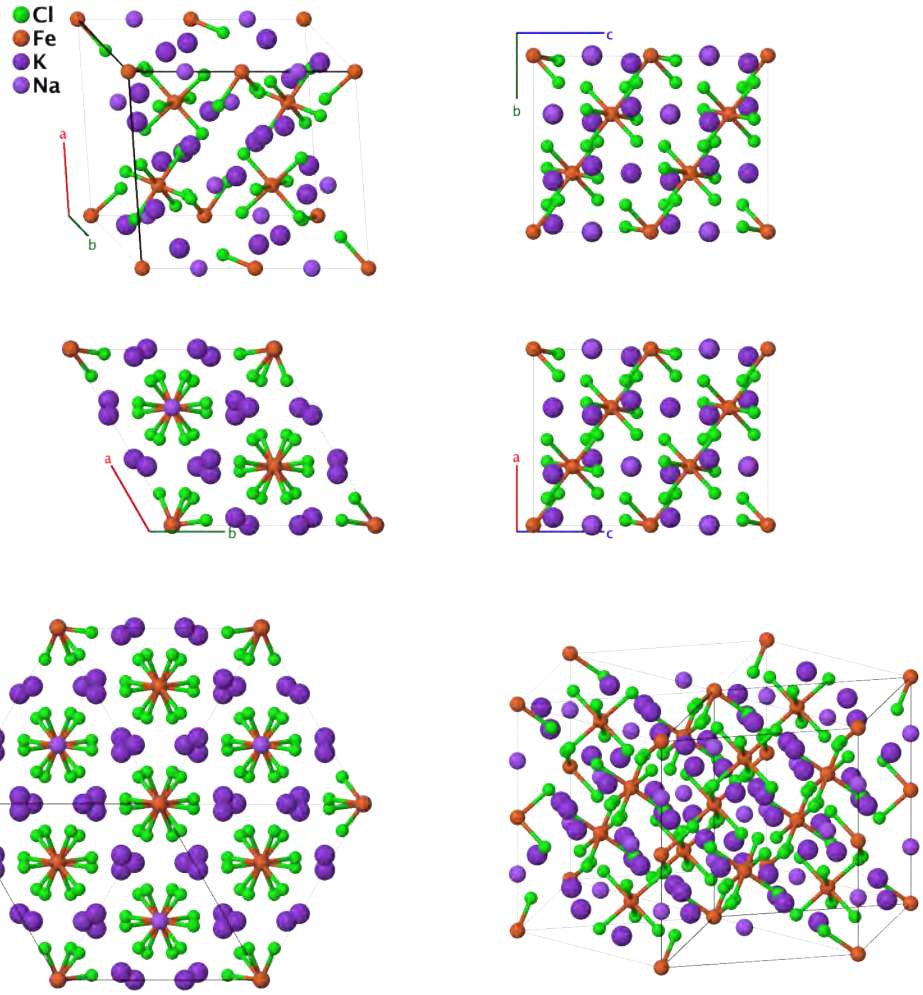
- M. A. Hepworth, K. H. Jack, R. D. Peacock, and G. J. Westland, *The crystal structures of the trifluorides of iron, cobalt, ruthenium, rhodium, palladium and iridium*, Acta Cryst. **10**, 63–69 (1957), doi:10.1107/S0365110X57000158.

Geometry files:

- CIF: pp. 1745
- POSCAR: pp. 1746

Rinneite ($K_3NaFeCl_6$) Structure: A6BC3D_hR22_167_f_b_e_a

http://aflow.org/prototype-encyclopedia/A6BC3D_hR22_167_f_b_e_a



Prototype	:	Cl_6FeK_3Na
AFLOW prototype label	:	A6BC3D_hR22_167_f_b_e_a
***Strukturbericht* designation**	:	None
Pearson symbol	:	hR22
Space group number	:	167
Space group symbol	:	$R\bar{3}c$
AFLOW prototype command	:	`aflow --proto=A6BC3D_hR22_167_f_b_e_a [--hex] --params=`$a, c/a, x_3, x_4, y_4, z_4$

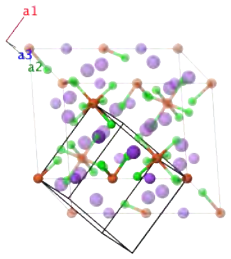
Other compounds with this structure

- Ca_3LiRuO_6, Sr_3NiIrO_6, and K_4CdCl_6

- We use the data taken at 293 K.

Rhombohedral primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2\sqrt{3}}\,a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,c\,\hat{\mathbf{z}}\\ \mathbf{a}_2 &= \tfrac{1}{\sqrt{3}}\,a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,c\,\hat{\mathbf{z}}\\ \mathbf{a}_3 &= -\tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2\sqrt{3}}\,a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\tfrac{1}{4}\,\mathbf{a}_1 + \tfrac{1}{4}\,\mathbf{a}_2 + \tfrac{1}{4}\,\mathbf{a}_3$	=	$\tfrac{1}{4}c\,\hat{\mathbf{z}}$	$(2a)$	Na
$\mathbf{B}_2$	=	$\tfrac{3}{4}\,\mathbf{a}_1 + \tfrac{3}{4}\,\mathbf{a}_2 + \tfrac{3}{4}\,\mathbf{a}_3$	=	$\tfrac{3}{4}c\,\hat{\mathbf{z}}$	$(2a)$	Na
$\mathbf{B}_3$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2b)$	Fe
$\mathbf{B}_4$	=	$\tfrac{1}{2}\,\mathbf{a}_1 + \tfrac{1}{2}\,\mathbf{a}_2 + \tfrac{1}{2}\,\mathbf{a}_3$	=	$\tfrac{1}{2}c\,\hat{\mathbf{z}}$	$(2b)$	Fe
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_3\right)\mathbf{a}_2 + \tfrac{1}{4}\,\mathbf{a}_3$	=	$\left(-\tfrac{1}{8} + \tfrac{1}{2}x_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{\sqrt{3}}{8} - \tfrac{\sqrt{3}}{2}x_3\right)a\,\hat{\mathbf{y}} + \tfrac{1}{4}c\,\hat{\mathbf{z}}$	$(6e)$	K
$\mathbf{B}_6$	=	$\tfrac{1}{4}\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_3\right)\mathbf{a}_3$	=	$\left(-\tfrac{1}{8} + \tfrac{1}{2}x_3\right)a\,\hat{\mathbf{x}} + \left(-\tfrac{\sqrt{3}}{8} + \tfrac{\sqrt{3}}{2}x_3\right)a\,\hat{\mathbf{y}} + \tfrac{1}{4}c\,\hat{\mathbf{z}}$	$(6e)$	K
$\mathbf{B}_7$	=	$\left(\tfrac{1}{2} - x_3\right)\mathbf{a}_1 + \tfrac{1}{4}\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$\left(\tfrac{1}{4} - x_3\right)a\,\hat{\mathbf{x}} + \tfrac{1}{4}c\,\hat{\mathbf{z}}$	$(6e)$	K
$\mathbf{B}_8$	=	$-x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_3\right)\mathbf{a}_2 + \tfrac{3}{4}\,\mathbf{a}_3$	=	$-\left(\tfrac{1}{2}x_3 + \tfrac{3}{8}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{8\sqrt{3}} + \tfrac{\sqrt{3}}{2}x_3\right)a\,\hat{\mathbf{y}} + \tfrac{5}{12}c\,\hat{\mathbf{z}}$	$(6e)$	K
$\mathbf{B}_9$	=	$\tfrac{3}{4}\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_3\right)\mathbf{a}_3$	=	$\left(\tfrac{1}{8} - \tfrac{1}{2}x_3\right)a\,\hat{\mathbf{x}} - \left(\tfrac{\sqrt{3}}{2}x_3 + \tfrac{5}{8\sqrt{3}}\right)a\,\hat{\mathbf{y}} + \tfrac{5}{12}c\,\hat{\mathbf{z}}$	$(6e)$	K
$\mathbf{B}_{10}$	=	$\left(\tfrac{1}{2} + x_3\right)\mathbf{a}_1 + \tfrac{3}{4}\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$\left(\tfrac{1}{4} + x_3\right)a\,\hat{\mathbf{x}} + \tfrac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \tfrac{5}{12}c\,\hat{\mathbf{z}}$	$(6e)$	K
$\mathbf{B}_{11}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\tfrac{1}{2}\,(x_4 - z_4)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_4 + \tfrac{1}{\sqrt{3}}y_4 - \tfrac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,(x_4 + y_4 + z_4)\,c\,\hat{\mathbf{z}}$	$(12f)$	Cl
$\mathbf{B}_{12}$	=	$z_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + y_4\,\mathbf{a}_3$	=	$\tfrac{1}{2}\,(-y_4 + z_4)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{\sqrt{3}}x_4 - \tfrac{1}{2\sqrt{3}}y_4 - \tfrac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,(x_4 + y_4 + z_4)\,c\,\hat{\mathbf{z}}$	$(12f)$	Cl
$\mathbf{B}_{13}$	=	$y_4\,\mathbf{a}_1 + z_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$\tfrac{1}{2}\,(-x_4 + y_4)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_4 - \tfrac{1}{2\sqrt{3}}y_4 + \tfrac{1}{\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,(x_4 + y_4 + z_4)\,c\,\hat{\mathbf{z}}$	$(12f)$	Cl
$\mathbf{B}_{14}$	=	$\left(\tfrac{1}{2} - z_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_4\right)\mathbf{a}_3$	=	$\tfrac{1}{2}\,(x_4 - z_4)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2\sqrt{3}}x_4 - \tfrac{1}{\sqrt{3}}y_4 + \tfrac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - \tfrac{1}{3}x_4 - \tfrac{1}{3}y_4 - \tfrac{1}{3}z_4\right)c\,\hat{\mathbf{z}}$	$(12f)$	Cl
$\mathbf{B}_{15}$	=	$\left(\tfrac{1}{2} - y_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\tfrac{1}{2}\,(-y_4 + z_4)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{\sqrt{3}}x_4 + \tfrac{1}{2\sqrt{3}}y_4 + \tfrac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - \tfrac{1}{3}x_4 - \tfrac{1}{3}y_4 - \tfrac{1}{3}z_4\right)c\,\hat{\mathbf{z}}$	$(12f)$	Cl

$\mathbf{B}_{16}$	=	$\left(\frac{1}{2}-x_4\right)\mathbf{a}_1+\left(\frac{1}{2}-z_4\right)\mathbf{a}_2+\left(\frac{1}{2}-y_4\right)\mathbf{a}_3$	=	$\frac{1}{2}(-x_4+y_4)\,a\,\hat{\mathbf{x}}+\left(\frac{1}{2\sqrt{3}}x_4+\frac{1}{2\sqrt{3}}y_4-\frac{1}{\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-\frac{1}{3}x_4-\frac{1}{3}y_4-\frac{1}{3}z_4\right)c\,\hat{\mathbf{z}}$	$(12f)$	Cl
$\mathbf{B}_{17}$	=	$-x_4\,\mathbf{a}_1-y_4\,\mathbf{a}_2-z_4\,\mathbf{a}_3$	=	$\frac{1}{2}(-x_4+z_4)\,a\,\hat{\mathbf{x}}+\left(\frac{1}{2\sqrt{3}}x_4-\frac{1}{\sqrt{3}}y_4+\frac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}}-\frac{1}{3}(x_4+y_4+z_4)\,c\,\hat{\mathbf{z}}$	$(12f)$	Cl
$\mathbf{B}_{18}$	=	$-z_4\,\mathbf{a}_1-x_4\,\mathbf{a}_2-y_4\,\mathbf{a}_3$	=	$\frac{1}{2}(y_4-z_4)\,a\,\hat{\mathbf{x}}+\left(-\frac{1}{\sqrt{3}}x_4+\frac{1}{2\sqrt{3}}y_4+\frac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}}-\frac{1}{3}(x_4+y_4+z_4)\,c\,\hat{\mathbf{z}}$	$(12f)$	Cl
$\mathbf{B}_{19}$	=	$-y_4\,\mathbf{a}_1-z_4\,\mathbf{a}_2-x_4\,\mathbf{a}_3$	=	$\frac{1}{2}(x_4-y_4)\,a\,\hat{\mathbf{x}}+\left(\frac{1}{2\sqrt{3}}x_4+\frac{1}{2\sqrt{3}}y_4-\frac{1}{\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}}-\frac{1}{3}(x_4+y_4+z_4)\,c\,\hat{\mathbf{z}}$	$(12f)$	Cl
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2}+z_4\right)\mathbf{a}_1+\left(\frac{1}{2}+y_4\right)\mathbf{a}_2+\left(\frac{1}{2}+x_4\right)\mathbf{a}_3$	=	$\frac{1}{2}(-x_4+z_4)\,a\,\hat{\mathbf{x}}+\left(-\frac{1}{2\sqrt{3}}x_4+\frac{1}{\sqrt{3}}y_4-\frac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+\frac{1}{3}x_4+\frac{1}{3}y_4+\frac{1}{3}z_4\right)c\,\hat{\mathbf{z}}$	$(12f)$	Cl
$\mathbf{B}_{21}$	=	$\left(\frac{1}{2}+y_4\right)\mathbf{a}_1+\left(\frac{1}{2}+x_4\right)\mathbf{a}_2+\left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	=	$\frac{1}{2}(y_4-z_4)\,a\,\hat{\mathbf{x}}+\left(\frac{1}{\sqrt{3}}x_4-\frac{1}{2\sqrt{3}}y_4-\frac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+\frac{1}{3}x_4+\frac{1}{3}y_4+\frac{1}{3}z_4\right)c\,\hat{\mathbf{z}}$	$(12f)$	Cl
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2}+x_4\right)\mathbf{a}_1+\left(\frac{1}{2}+z_4\right)\mathbf{a}_2+\left(\frac{1}{2}+y_4\right)\mathbf{a}_3$	=	$\frac{1}{2}(x_4-y_4)\,a\,\hat{\mathbf{x}}+\left(-\frac{1}{2\sqrt{3}}x_4-\frac{1}{2\sqrt{3}}y_4+\frac{1}{\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+\frac{1}{3}x_4+\frac{1}{3}y_4+\frac{1}{3}z_4\right)c\,\hat{\mathbf{z}}$	$(12f)$	Cl

References:

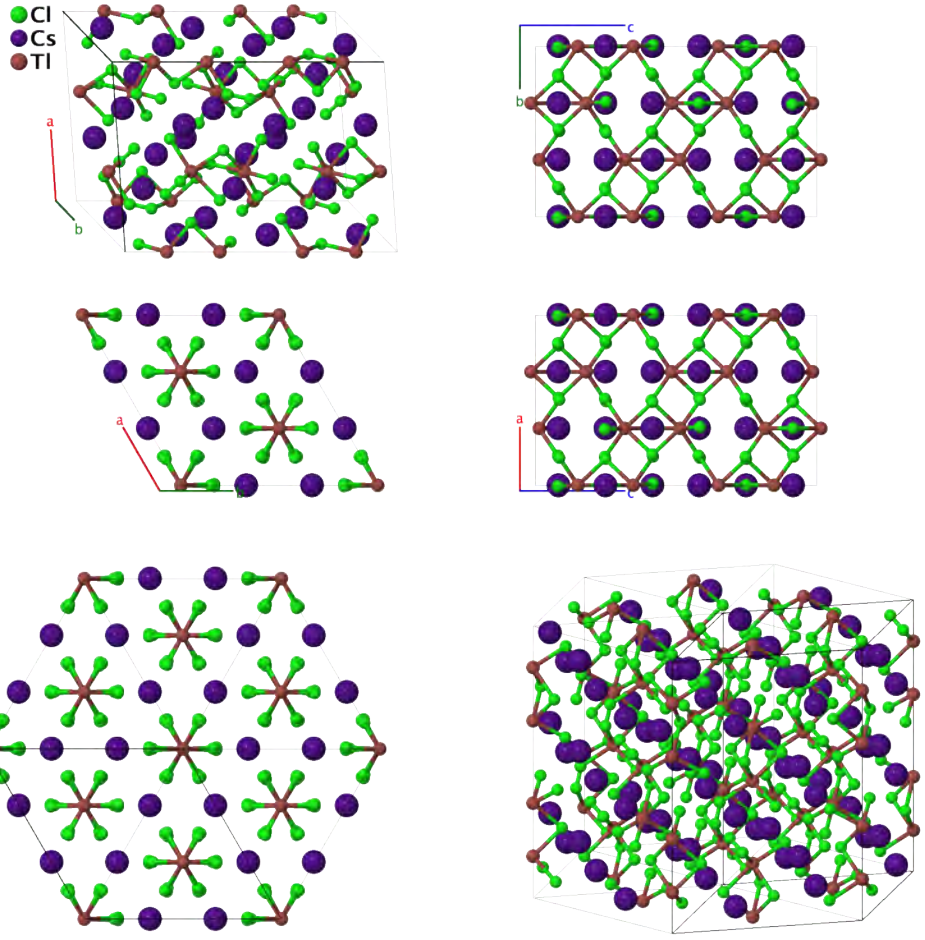
- B. N. Figgis, A. N. Sobolev, E. S. Kucharski, and V. Broughton, *Rinneite, $K_3Na[FeCl_6]$, at 293, 84 and 9.5 K*, Acta Crystallogr. C **56**, e228–e229 (2000), doi:10.1107/S0108270100006053.

Geometry files:

- CIF: pp. 1746
- POSCAR: pp. 1746

$Cs_3Tl_2Cl_9$ ($K7_2$) Structure: A9B3C2_hR28_167_ef_e_c

http://aflow.org/prototype-encyclopedia/A9B3C2_hR28_167_ef_e_c



Prototype	:	$Cl_9Cs_3Tl_2$
AFLOW prototype label	:	A9B3C2_hR28_167_ef_e_c
***Strukturbericht* designation**	:	$K7_2$
Pearson symbol	:	hR28
Space group number	:	167
Space group symbol	:	$R\bar{3}c$
AFLOW prototype command	:	`aflow --proto=A9B3C2_hR28_167_ef_e_c [--hex] --params=`$a, c/a, x_1, x_2, x_3, x_4, y_4, z_4$

Other compounds with this structure

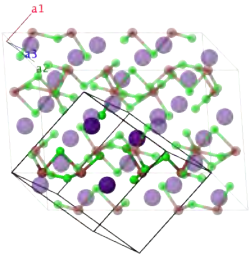
- $Cs_3Dy_2Br_9$, $Cs_3Er_2Br_9$, $Cs_3Ho_2Br_9$, $Cs_3Lu_2Cl_9$, $Cs_3Tb_2Br_9$, $Cs_3Yb_2Br_9$, $Ba_3Os_2O_9$, and $Ba_3W_2O_9$

- (Hoard, 1935) followed by (Downs, 2003), give the atomic coordinates in the style of (Wyckoff, 1922), who lists space group D^6_{3d} (the Schönflies notation for space group $R\bar{3}c$) as space group #203, instead of #167, and uses an origin which

corresponds to (1/4, 1/4, 1/4) in our lattice coordinates. We used FINDSYM to convert this into our standard setting for space group #167.

Rhombohedral primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2\sqrt{3}}\,a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,c\,\hat{\mathbf{z}}\\ \mathbf{a}_2 &= \tfrac{1}{\sqrt{3}}\,a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,c\,\hat{\mathbf{z}}\\ \mathbf{a}_3 &= -\tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2\sqrt{3}}\,a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + x_1\,\mathbf{a}_2 + x_1\,\mathbf{a}_3$	=	$x_1 c\,\hat{\mathbf{z}}$	(4*c*)	Tl
$\mathbf{B}_2$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_1\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	Tl
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 - x_1\,\mathbf{a}_2 - x_1\,\mathbf{a}_3$	=	$-x_1 c\,\hat{\mathbf{z}}$	(4*c*)	Tl
$\mathbf{B}_4$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_1\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right) c\,\hat{\mathbf{z}}$	(4*c*)	Tl
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(-\frac{1}{8} + \frac{1}{2}x_2\right) a\,\hat{\mathbf{x}} + \left(\frac{\sqrt{3}}{8} - \frac{\sqrt{3}}{2}x_2\right) a\,\hat{\mathbf{y}} + \frac{1}{4} c\,\hat{\mathbf{z}}$	(6*e*)	Cl I
$\mathbf{B}_6$	=	$\frac{1}{4}\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_3$	=	$\left(-\frac{1}{8} + \frac{1}{2}x_2\right) a\,\hat{\mathbf{x}} + \left(-\frac{\sqrt{3}}{8} + \frac{\sqrt{3}}{2}x_2\right) a\,\hat{\mathbf{y}} + \frac{1}{4} c\,\hat{\mathbf{z}}$	(6*e*)	Cl I
$\mathbf{B}_7$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_2\right) a\,\hat{\mathbf{x}} + \frac{1}{4} c\,\hat{\mathbf{z}}$	(6*e*)	Cl I
$\mathbf{B}_8$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-\left(\frac{1}{2}x_2 + \frac{3}{8}\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{8\sqrt{3}} + \frac{\sqrt{3}}{2}x_2\right) a\,\hat{\mathbf{y}} + \frac{5}{12} c\,\hat{\mathbf{z}}$	(6*e*)	Cl I
$\mathbf{B}_9$	=	$\frac{3}{4}\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{8} - \frac{1}{2}x_2\right) a\,\hat{\mathbf{x}} - \left(\frac{\sqrt{3}}{2}x_2 + \frac{5}{8\sqrt{3}}\right) a\,\hat{\mathbf{y}} + \frac{5}{12} c\,\hat{\mathbf{z}}$	(6*e*)	Cl I
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - x_2\,\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_2\right) a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}} a\,\hat{\mathbf{y}} + \frac{5}{12} c\,\hat{\mathbf{z}}$	(6*e*)	Cl I
$\mathbf{B}_{11}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(-\frac{1}{8} + \frac{1}{2}x_3\right) a\,\hat{\mathbf{x}} + \left(\frac{\sqrt{3}}{8} - \frac{\sqrt{3}}{2}x_3\right) a\,\hat{\mathbf{y}} + \frac{1}{4} c\,\hat{\mathbf{z}}$	(6*e*)	Cs
$\mathbf{B}_{12}$	=	$\frac{1}{4}\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_3$	=	$\left(-\frac{1}{8} + \frac{1}{2}x_3\right) a\,\hat{\mathbf{x}} + \left(-\frac{\sqrt{3}}{8} + \frac{\sqrt{3}}{2}x_3\right) a\,\hat{\mathbf{y}} + \frac{1}{4} c\,\hat{\mathbf{z}}$	(6*e*)	Cs
$\mathbf{B}_{13}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_3\right) a\,\hat{\mathbf{x}} + \frac{1}{4} c\,\hat{\mathbf{z}}$	(6*e*)	Cs
$\mathbf{B}_{14}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-\left(\frac{1}{2}x_3 + \frac{3}{8}\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{8\sqrt{3}} + \frac{\sqrt{3}}{2}x_3\right) a\,\hat{\mathbf{y}} + \frac{5}{12} c\,\hat{\mathbf{z}}$	(6*e*)	Cs
$\mathbf{B}_{15}$	=	$\frac{3}{4}\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{8} - \frac{1}{2}x_3\right) a\,\hat{\mathbf{x}} - \left(\frac{\sqrt{3}}{2}x_3 + \frac{5}{8\sqrt{3}}\right) a\,\hat{\mathbf{y}} + \frac{5}{12} c\,\hat{\mathbf{z}}$	(6*e*)	Cs
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_3\right) a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}} a\,\hat{\mathbf{y}} + \frac{5}{12} c\,\hat{\mathbf{z}}$	(6*e*)	Cs

$$\mathbf{B}_{17} = x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3 = \tfrac{1}{2}\,(x_4 - z_4)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_4 + \tfrac{1}{\sqrt{3}}y_4 - \tfrac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,(x_4 + y_4 + z_4)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Cl II}$$

$$\mathbf{B}_{18} = z_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + y_4\,\mathbf{a}_3 = \tfrac{1}{2}\,(-y_4 + z_4)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{\sqrt{3}}x_4 - \tfrac{1}{2\sqrt{3}}y_4 - \tfrac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,(x_4 + y_4 + z_4)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Cl II}$$

$$\mathbf{B}_{19} = y_4\,\mathbf{a}_1 + z_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3 = \tfrac{1}{2}\,(-x_4 + y_4)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_4 - \tfrac{1}{2\sqrt{3}}y_4 + \tfrac{1}{\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} + \tfrac{1}{3}\,(x_4 + y_4 + z_4)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Cl II}$$

$$\mathbf{B}_{20} = \left(\tfrac{1}{2} - z_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_4\right)\mathbf{a}_3 = \tfrac{1}{2}\,(x_4 - z_4)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2\sqrt{3}}x_4 - \tfrac{1}{\sqrt{3}}y_4 + \tfrac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - \tfrac{1}{3}x_4 - \tfrac{1}{3}y_4 - \tfrac{1}{3}z_4\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Cl II}$$

$$\mathbf{B}_{21} = \left(\tfrac{1}{2} - y_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_4\right)\mathbf{a}_3 = \tfrac{1}{2}\,(-y_4 + z_4)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{\sqrt{3}}x_4 + \tfrac{1}{2\sqrt{3}}y_4 + \tfrac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - \tfrac{1}{3}x_4 - \tfrac{1}{3}y_4 - \tfrac{1}{3}z_4\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Cl II}$$

$$\mathbf{B}_{22} = \left(\tfrac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - z_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - y_4\right)\mathbf{a}_3 = \tfrac{1}{2}\,(-x_4 + y_4)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2\sqrt{3}}x_4 + \tfrac{1}{2\sqrt{3}}y_4 - \tfrac{1}{\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - \tfrac{1}{3}x_4 - \tfrac{1}{3}y_4 - \tfrac{1}{3}z_4\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Cl II}$$

$$\mathbf{B}_{23} = -x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3 = \tfrac{1}{2}\,(-x_4 + z_4)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2\sqrt{3}}x_4 - \tfrac{1}{\sqrt{3}}y_4 + \tfrac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} - \tfrac{1}{3}\,(x_4 + y_4 + z_4)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Cl II}$$

$$\mathbf{B}_{24} = -z_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - y_4\,\mathbf{a}_3 = \tfrac{1}{2}\,(y_4 - z_4)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{\sqrt{3}}x_4 + \tfrac{1}{2\sqrt{3}}y_4 + \tfrac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} - \tfrac{1}{3}\,(x_4 + y_4 + z_4)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Cl II}$$

$$\mathbf{B}_{25} = -y_4\,\mathbf{a}_1 - z_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3 = \tfrac{1}{2}\,(x_4 - y_4)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2\sqrt{3}}x_4 + \tfrac{1}{2\sqrt{3}}y_4 - \tfrac{1}{\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} - \tfrac{1}{3}\,(x_4 + y_4 + z_4)\,c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Cl II}$$

$$\mathbf{B}_{26} = \left(\tfrac{1}{2} + z_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_4\right)\mathbf{a}_3 = \tfrac{1}{2}\,(-x_4 + z_4)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_4 + \tfrac{1}{\sqrt{3}}y_4 - \tfrac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + \tfrac{1}{3}x_4 + \tfrac{1}{3}y_4 + \tfrac{1}{3}z_4\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Cl II}$$

$$\mathbf{B}_{27} = \left(\tfrac{1}{2} + y_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_4\right)\mathbf{a}_3 = \tfrac{1}{2}\,(y_4 - z_4)\,a\,\hat{\mathbf{x}} + \left(\tfrac{1}{\sqrt{3}}x_4 - \tfrac{1}{2\sqrt{3}}y_4 - \tfrac{1}{2\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + \tfrac{1}{3}x_4 + \tfrac{1}{3}y_4 + \tfrac{1}{3}z_4\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Cl II}$$

$$\mathbf{B}_{28} = \left(\tfrac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + z_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + y_4\right)\mathbf{a}_3 = \tfrac{1}{2}\,(x_4 - y_4)\,a\,\hat{\mathbf{x}} + \left(-\tfrac{1}{2\sqrt{3}}x_4 - \tfrac{1}{2\sqrt{3}}y_4 + \tfrac{1}{\sqrt{3}}z_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + \tfrac{1}{3}x_4 + \tfrac{1}{3}y_4 + \tfrac{1}{3}z_4\right)c\,\hat{\mathbf{z}} \quad (12f) \quad \text{Cl II}$$

References:

- J. L. Hoard and L. Goldstein, *The Crystal Structure of Cesium Enneachlordithalliate, $Cs_3Tl_2Cl_9$*, J. Chem. Phys. **3**, 199–202 (1935), doi:10.1063/1.1749633.

- R. W. G. Wyckoff, *The Analytical Expression of the Results of the Theory of Space-Groups*, vol. 318 (Carnegie Institution of Washington, Washington DC, 1922).

Found in:

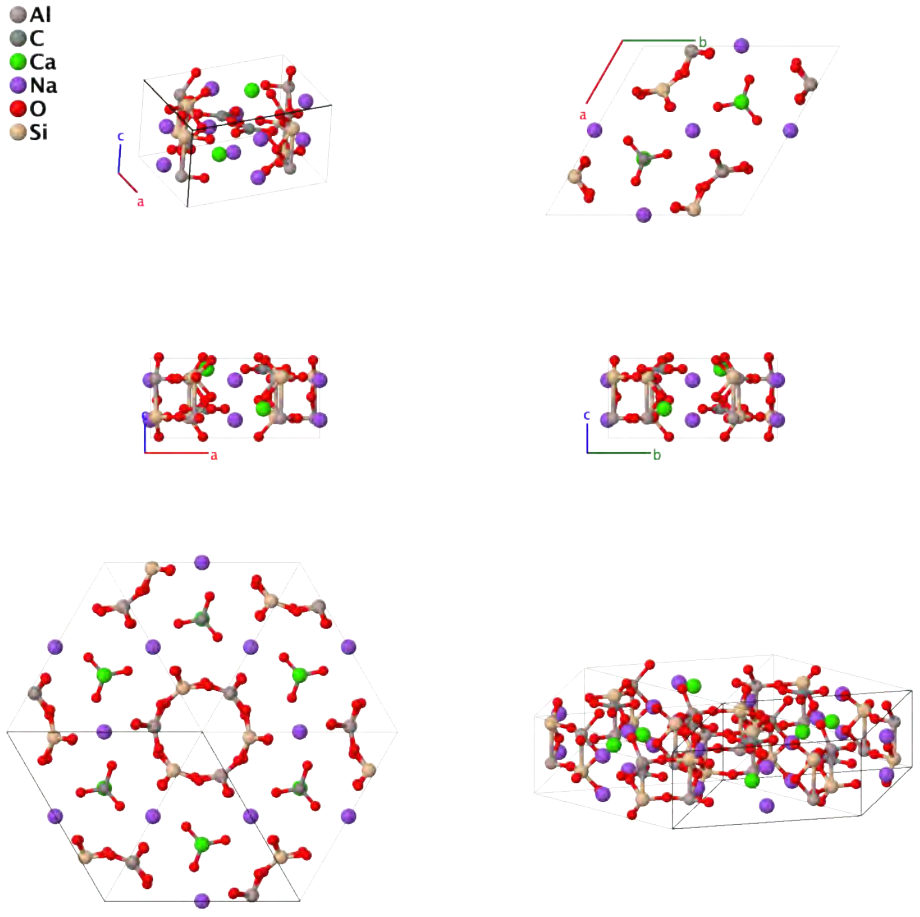
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1747
- POSCAR: pp. 1747

Crancrinite ($Na_6Ca_2Al_6Si_6O_{24}(CO_3)_2$, $S\,3_3$ (I)) Structure: A3BCD3E15F3_hP52_173_c_b_b_c_5c_c

http://aflow.org/prototype-encyclopedia/A3BCD3E15F3_hP52_173_c_b_b_c_5c_c



Prototype	:	$Al_3CCaNa_3O_{15}Si_3$
AFLOW prototype label	:	A3BCD3E15F3_hP52_173_c_b_b_c_5c_c
***Strukturbericht* designation**	:	$S\,3_3$ (I)
Pearson symbol	:	hP52
Space group number	:	173
Space group symbol	:	$P6_3$
AFLOW prototype command	:	`aflow --proto=A3BCD3E15F3_hP52_173_c_b_b_c_5c_c --params=`$a, c/a, z_1, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}$

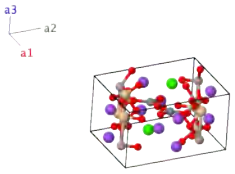
- The term "crancrinite" covers a wide range of mineral compositions, some of which may include OH radicals and H_2O molecules. We use this early determination of the structure as the base.
- There are two major inconsistencies between (Kozu, 1933) and (Gottfried, 1937):
  - (Gottfried, 1937) interchanges the z coordinates of the carbon and calcium ions compared to (Kozu, 1933). Kozu's coordinates are consistent with CO_3 radicals, so we use their version.

  - (Kozu, 1933) gives $x_{10} = 0.33$ for the x coordinate of silicon. This puts the silicon atoms very close to carbon. We follow (Gottfried, 1937) and take $x_{10} = 0.033$, which gives a reasonable distance.

- (Gottfried, 1937) gave this structure the label $S\,3_3$, but later (Gottfried, 1938) gave the same label to parawollastonite, $CaSiO_3$. We will refer to crancrinite as $S\,3_3$ (I), and parawollastonite as $S\,3_3$ (II).

Hexagonal primitive vectors:

$$\begin{aligned}
\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}} \\
\mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}} \\
\mathbf{a}_3 &= c\,\hat{\mathbf{z}}
\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	$(2b)$	C
$\mathbf{B}_2$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\,\hat{\mathbf{z}}$	$(2b)$	C
$\mathbf{B}_3$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(2b)$	Ca
$\mathbf{B}_4$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	$(2b)$	Ca
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_3 + y_3)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_3 + y_3)\,a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(6c)$	Al
$\mathbf{B}_6$	=	$-y_3\,\mathbf{a}_1 + (x_3 - y_3)\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_3 - y_3\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(6c)$	Al
$\mathbf{B}_7$	=	$(-x_3 + y_3)\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\left(-x_3 + \frac{1}{2}y_3\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(6c)$	Al
$\mathbf{B}_8$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$-\frac{1}{2}\,(x_3 + y_3)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(x_3 - y_3)\,a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	$(6c)$	Al
$\mathbf{B}_9$	=	$y_3\,\mathbf{a}_1 + (-x_3 + y_3)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_3 + y_3\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	$(6c)$	Al
$\mathbf{B}_{10}$	=	$(x_3 - y_3)\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(x_3 - \frac{1}{2}y_3\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	$(6c)$	Al
$\mathbf{B}_{11}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_4 + y_4)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_4 + y_4)\,a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(6c)$	Na
$\mathbf{B}_{12}$	=	$-y_4\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_4 - y_4\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(6c)$	Na
$\mathbf{B}_{13}$	=	$(-x_4 + y_4)\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\left(-x_4 + \frac{1}{2}y_4\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_4 a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(6c)$	Na
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$-\frac{1}{2}\,(x_4 + y_4)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(x_4 - y_4)\,a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	$(6c)$	Na
$\mathbf{B}_{15}$	=	$y_4\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_4 + y_4\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	$(6c)$	Na
$\mathbf{B}_{16}$	=	$(x_4 - y_4)\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(x_4 - \frac{1}{2}y_4\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_4 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	$(6c)$	Na
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_5 + y_5)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_5 + y_5)\,a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(6c)$	O I
$\mathbf{B}_{18}$	=	$-y_5\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_5 - y_5\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(6c)$	O I

$$\mathbf{B}_{19} = (-x_5 + y_5)\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3 = \left(-x_5 + \tfrac{1}{2}y_5\right)a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}y_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}} \quad (6c) \quad \text{O I}$$

$$\mathbf{B}_{20} = -x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_3 = -\tfrac{1}{2}(x_5 + y_5)\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}(x_5 - y_5)\,a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_5\right)c\,\hat{\mathbf{z}} \quad (6c) \quad \text{O I}$$

$$\mathbf{B}_{21} = y_5\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_3 = \left(-\tfrac{1}{2}x_5 + y_5\right)a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_5\right)c\,\hat{\mathbf{z}} \quad (6c) \quad \text{O I}$$

$$\mathbf{B}_{22} = (x_5 - y_5)\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_3 = \left(x_5 - \tfrac{1}{2}y_5\right)a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}y_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_5\right)c\,\hat{\mathbf{z}} \quad (6c) \quad \text{O I}$$

$$\mathbf{B}_{23} = x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 = \tfrac{1}{2}(x_6 + y_6)\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}(-x_6 + y_6)\,a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}} \quad (6c) \quad \text{O II}$$

$$\mathbf{B}_{24} = -y_6\,\mathbf{a}_1 + (x_6 - y_6)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 = \left(\tfrac{1}{2}x_6 - y_6\right)a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}} \quad (6c) \quad \text{O II}$$

$$\mathbf{B}_{25} = (-x_6 + y_6)\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 = \left(-x_6 + \tfrac{1}{2}y_6\right)a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}y_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}} \quad (6c) \quad \text{O II}$$

$$\mathbf{B}_{26} = -x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 = -\tfrac{1}{2}(x_6 + y_6)\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}(x_6 - y_6)\,a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right)c\,\hat{\mathbf{z}} \quad (6c) \quad \text{O II}$$

$$\mathbf{B}_{27} = y_6\,\mathbf{a}_1 + (-x_6 + y_6)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 = \left(-\tfrac{1}{2}x_6 + y_6\right)a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right)c\,\hat{\mathbf{z}} \quad (6c) \quad \text{O II}$$

$$\mathbf{B}_{28} = (x_6 - y_6)\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 = \left(x_6 - \tfrac{1}{2}y_6\right)a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}y_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right)c\,\hat{\mathbf{z}} \quad (6c) \quad \text{O II}$$

$$\mathbf{B}_{29} = x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 = \tfrac{1}{2}(x_7 + y_7)\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}(-x_7 + y_7)\,a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}} \quad (6c) \quad \text{O III}$$

$$\mathbf{B}_{30} = -y_7\,\mathbf{a}_1 + (x_7 - y_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 = \left(\tfrac{1}{2}x_7 - y_7\right)a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}x_7 a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}} \quad (6c) \quad \text{O III}$$

$$\mathbf{B}_{31} = (-x_7 + y_7)\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 = \left(-x_7 + \tfrac{1}{2}y_7\right)a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}y_7 a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}} \quad (6c) \quad \text{O III}$$

$$\mathbf{B}_{32} = -x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_7\right)\mathbf{a}_3 = -\tfrac{1}{2}(x_7 + y_7)\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}(x_7 - y_7)\,a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_7\right)c\,\hat{\mathbf{z}} \quad (6c) \quad \text{O III}$$

$$\mathbf{B}_{33} = y_7\,\mathbf{a}_1 + (-x_7 + y_7)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_7\right)\mathbf{a}_3 = \left(-\tfrac{1}{2}x_7 + y_7\right)a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}x_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_7\right)c\,\hat{\mathbf{z}} \quad (6c) \quad \text{O III}$$

$$\mathbf{B}_{34} = (x_7 - y_7)\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_7\right)\mathbf{a}_3 = \left(x_7 - \tfrac{1}{2}y_7\right)a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}y_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_7\right)c\,\hat{\mathbf{z}} \quad (6c) \quad \text{O III}$$

$$\mathbf{B}_{35} = x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 = \tfrac{1}{2}(x_8 + y_8)\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}(-x_8 + y_8)\,a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}} \quad (6c) \quad \text{O IV}$$

$$\mathbf{B}_{36} = -y_8\,\mathbf{a}_1 + (x_8 - y_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 = \left(\tfrac{1}{2}x_8 - y_8\right)a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}x_8 a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}} \quad (6c) \quad \text{O IV}$$

$$\mathbf{B}_{37} = (-x_8 + y_8)\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 = \left(-x_8 + \tfrac{1}{2}y_8\right)a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}y_8 a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}} \quad (6c) \quad \text{O IV}$$

$$\mathbf{B}_{38} = -x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_8\right)\mathbf{a}_3 = -\tfrac{1}{2}(x_8 + y_8)\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}(x_8 - y_8)\,a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_8\right)c\,\hat{\mathbf{z}} \quad (6c) \quad \text{O IV}$$

$$\mathbf{B}_{39} = y_8\,\mathbf{a}_1 + (-x_8 + y_8)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_8\right)\mathbf{a}_3 = \left(-\tfrac{1}{2}x_8 + y_8\right)a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}x_8 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_8\right)c\,\hat{\mathbf{z}} \quad (6c) \quad \text{O IV}$$

$$\mathbf{B}_{40} = (x_8 - y_8)\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_8\right)\mathbf{a}_3 = \left(x_8 - \tfrac{1}{2}y_8\right)a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}y_8 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_8\right)c\,\hat{\mathbf{z}} \quad (6c) \quad \text{O IV}$$

$$\mathbf{B}_{41} = x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 = \tfrac{1}{2}(x_9 + y_9)\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}(-x_9 + y_9)\,a\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}} \quad (6c) \quad \text{O V}$$

$$\mathbf{B}_{42} = -y_9\,\mathbf{a}_1 + (x_9 - y_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 = \left(\tfrac{1}{2}x_9 - y_9\right)a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}x_9 a\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}} \quad (6c) \quad \text{O V}$$

$\mathbf{B}_{43}$	=	$(-x_9 + y_9)\,\mathbf{a}_1 - x_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$\left(-x_9 + \frac{1}{2}y_9\right) a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2} y_9 a\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(6*c*)	O V
$\mathbf{B}_{44}$	=	$-x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$-\frac{1}{2}(x_9 + y_9)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(x_9 - y_9)\,a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right) c\,\hat{\mathbf{z}}$	(6*c*)	O V
$\mathbf{B}_{45}$	=	$y_9\,\mathbf{a}_1 + (-x_9 + y_9)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_9 + y_9\right) a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2} x_9 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right) c\,\hat{\mathbf{z}}$	(6*c*)	O V
$\mathbf{B}_{46}$	=	$(x_9 - y_9)\,\mathbf{a}_1 + x_9\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$\left(x_9 - \frac{1}{2}y_9\right) a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2} y_9 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right) c\,\hat{\mathbf{z}}$	(6*c*)	O V
$\mathbf{B}_{47}$	=	$x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$\frac{1}{2}(x_{10} + y_{10})\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(-x_{10} + y_{10})\,a\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}}$	(6*c*)	Si
$\mathbf{B}_{48}$	=	$-y_{10}\,\mathbf{a}_1 + (x_{10} - y_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_{10} - y_{10}\right) a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2} x_{10} a\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}}$	(6*c*)	Si
$\mathbf{B}_{49}$	=	$(-x_{10} + y_{10})\,\mathbf{a}_1 - x_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$\left(-x_{10} + \frac{1}{2}y_{10}\right) a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2} y_{10} a\,\hat{\mathbf{y}} + z_{10} c\,\hat{\mathbf{z}}$	(6*c*)	Si
$\mathbf{B}_{50}$	=	$-x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{10}\right)\mathbf{a}_3$	=	$-\frac{1}{2}(x_{10} + y_{10})\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(x_{10} - y_{10})\,a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{10}\right) c\,\hat{\mathbf{z}}$	(6*c*)	Si
$\mathbf{B}_{51}$	=	$y_{10}\,\mathbf{a}_1 + (-x_{10} + y_{10})\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{10}\right)\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_{10} + y_{10}\right) a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2} x_{10} a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{10}\right) c\,\hat{\mathbf{z}}$	(6*c*)	Si
$\mathbf{B}_{52}$	=	$(x_{10} - y_{10})\,\mathbf{a}_1 + x_{10}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{10}\right)\mathbf{a}_3$	=	$\left(x_{10} - \frac{1}{2}y_{10}\right) a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2} y_{10} a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{10}\right) c\,\hat{\mathbf{z}}$	(6*c*)	Si

References:

- S. Kôzu and K. Takané, *Crystal Structure of Cancrinite from Dôdô, Korea (Part I)*, Proc. Imp. Acad. Japan **9**, 56–59 (1933), doi:10.2183/pjab1912.9.56.
- S. Kôzu and K. Takané, *Crystal Structure of Cancrinite from Dôdô, Korea (Part II)*, Proc. Imp. Acad. Japan **9**, 105–108 (1933), doi:10.2183/pjab1912.9.105.
- C. Gottfried, ed., *Strukturbericht Band IV 1936* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1938).

Found in:

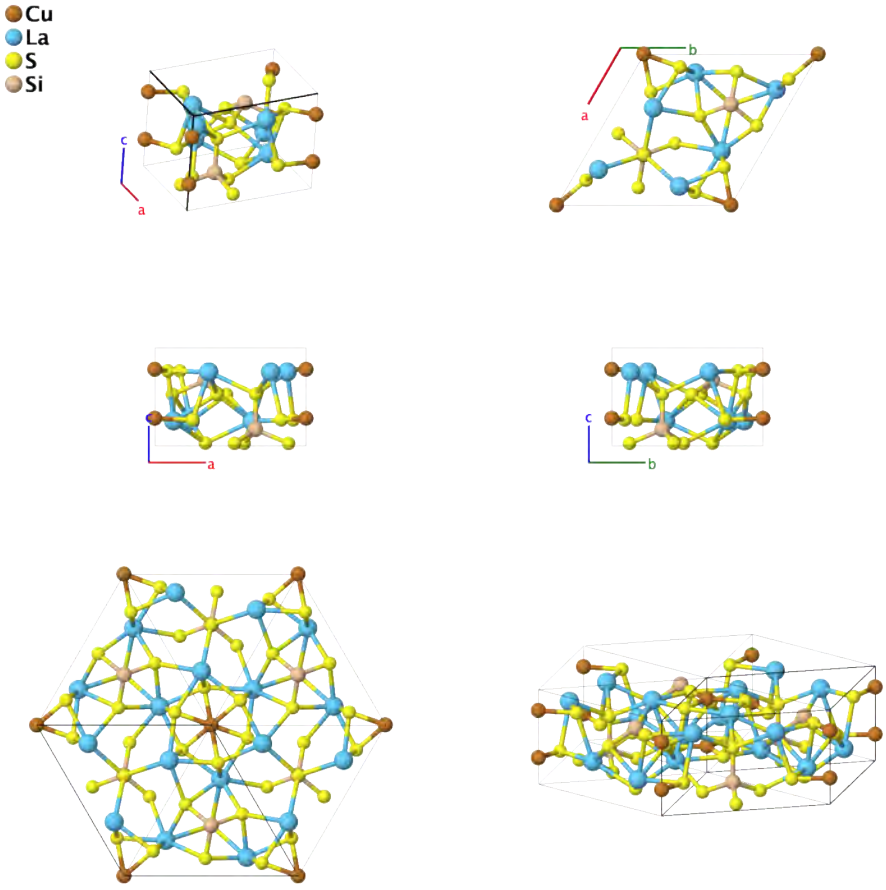
- C. Gottfried and F. Schossberger, eds., *Strukturbericht Band III 1933-1935* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1747
- POSCAR: pp. 1748

La_3CuSiS_7 Structure: AB3C7D_hP24_173_a_c_b2c_b

http://aflow.org/prototype-encyclopedia/AB3C7D_hP24_173_a_c_b2c_b



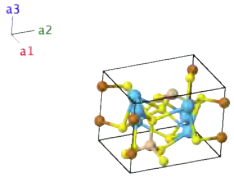
Prototype	:	$CuLa_3S_7Si$
AFLOW prototype label	:	AB3C7D_hP24_173_a_c_b2c_b
***Strukturbericht* designation**	:	None
Pearson symbol	:	hP24
Space group number	:	173
Space group symbol	:	$P6_3$
AFLOW prototype command	:	`aflow --proto=AB3C7D_hP24_173_a_c_b2c_b --params=`$a, c/a, z_1, z_2, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6$

Other compounds with this structure

- Ce_3CuGeS_7, $Ce_3CuGeSe_7$, $Ce_3CuSnSe_7$, Dy_3CuGeS_7, Gd_3CuGeS_7, $Gd_3CuGeSe_7$, $Ho_3CuGeSe_7$, La_3MnFeS_7, Nd_3CuGeS_7, $Nd_3CuGeSe_7$, Pr_3CuGeS_7, $Pr_3CuGeSe_7$, Sm_3CuGeS_7, $Sm_3CuGeSe_7$, Tb_3CuGeS_7, $Tb_3CuGeSe_7$, Y_3CuSiS_7, and $Y_3CuSiSe_7$

Hexagonal primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$z_1\,\mathbf{a}_3$	=	$z_1 c\,\hat{\mathbf{z}}$	$(2a)$	Cu
$\mathbf{B}_2$	=	$\left(\frac{1}{2}+z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+z_1\right)c\,\hat{\mathbf{z}}$	$(2a)$	Cu
$\mathbf{B}_3$	=	$\frac{1}{3}\,\mathbf{a}_1+\frac{2}{3}\,\mathbf{a}_2+z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}+\frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+z_2 c\,\hat{\mathbf{z}}$	$(2b)$	S I
$\mathbf{B}_4$	=	$\frac{2}{3}\,\mathbf{a}_1+\frac{1}{3}\,\mathbf{a}_2+\left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}-\frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_2\right)c\,\hat{\mathbf{z}}$	$(2b)$	S I
$\mathbf{B}_5$	=	$\frac{1}{3}\,\mathbf{a}_1+\frac{2}{3}\,\mathbf{a}_2+z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}+\frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+z_3 c\,\hat{\mathbf{z}}$	$(2b)$	Si
$\mathbf{B}_6$	=	$\frac{2}{3}\,\mathbf{a}_1+\frac{1}{3}\,\mathbf{a}_2+\left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}-\frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_3\right)c\,\hat{\mathbf{z}}$	$(2b)$	Si
$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_1+y_4\,\mathbf{a}_2+z_4\,\mathbf{a}_3$	=	$\frac{1}{2}(x_4+y_4)\,a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}(-x_4+y_4)\,a\,\hat{\mathbf{y}}+z_4 c\,\hat{\mathbf{z}}$	$(6c)$	La
$\mathbf{B}_8$	=	$-y_4\,\mathbf{a}_1+(x_4-y_4)\,\mathbf{a}_2+z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_4-y_4\right)a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}}+z_4 c\,\hat{\mathbf{z}}$	$(6c)$	La
$\mathbf{B}_9$	=	$(-x_4+y_4)\,\mathbf{a}_1-x_4\,\mathbf{a}_2+z_4\,\mathbf{a}_3$	=	$\left(-x_4+\frac{1}{2}y_4\right)a\,\hat{\mathbf{x}}-\frac{\sqrt{3}}{2}y_4 a\,\hat{\mathbf{y}}+z_4 c\,\hat{\mathbf{z}}$	$(6c)$	La
$\mathbf{B}_{10}$	=	$-x_4\,\mathbf{a}_1-y_4\,\mathbf{a}_2+\left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	=	$-\frac{1}{2}(x_4+y_4)\,a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}(x_4-y_4)\,a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_4\right)c\,\hat{\mathbf{z}}$	$(6c)$	La
$\mathbf{B}_{11}$	=	$y_4\,\mathbf{a}_1+(-x_4+y_4)\,\mathbf{a}_2+\left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_4+y_4\right)a\,\hat{\mathbf{x}}-\frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_4\right)c\,\hat{\mathbf{z}}$	$(6c)$	La
$\mathbf{B}_{12}$	=	$(x_4-y_4)\,\mathbf{a}_1+x_4\,\mathbf{a}_2+\left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	=	$\left(x_4-\frac{1}{2}y_4\right)a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}y_4 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_4\right)c\,\hat{\mathbf{z}}$	$(6c)$	La
$\mathbf{B}_{13}$	=	$x_5\,\mathbf{a}_1+y_5\,\mathbf{a}_2+z_5\,\mathbf{a}_3$	=	$\frac{1}{2}(x_5+y_5)\,a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}(-x_5+y_5)\,a\,\hat{\mathbf{y}}+z_5 c\,\hat{\mathbf{z}}$	$(6c)$	S II
$\mathbf{B}_{14}$	=	$-y_5\,\mathbf{a}_1+(x_5-y_5)\,\mathbf{a}_2+z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_5-y_5\right)a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}}+z_5 c\,\hat{\mathbf{z}}$	$(6c)$	S II
$\mathbf{B}_{15}$	=	$(-x_5+y_5)\,\mathbf{a}_1-x_5\,\mathbf{a}_2+z_5\,\mathbf{a}_3$	=	$\left(-x_5+\frac{1}{2}y_5\right)a\,\hat{\mathbf{x}}-\frac{\sqrt{3}}{2}y_5 a\,\hat{\mathbf{y}}+z_5 c\,\hat{\mathbf{z}}$	$(6c)$	S II
$\mathbf{B}_{16}$	=	$-x_5\,\mathbf{a}_1-y_5\,\mathbf{a}_2+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$-\frac{1}{2}(x_5+y_5)\,a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}(x_5-y_5)\,a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	$(6c)$	S II
$\mathbf{B}_{17}$	=	$y_5\,\mathbf{a}_1+(-x_5+y_5)\,\mathbf{a}_2+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_5+y_5\right)a\,\hat{\mathbf{x}}-\frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	$(6c)$	S II
$\mathbf{B}_{18}$	=	$(x_5-y_5)\,\mathbf{a}_1+x_5\,\mathbf{a}_2+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$\left(x_5-\frac{1}{2}y_5\right)a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}y_5 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	$(6c)$	S II
$\mathbf{B}_{19}$	=	$x_6\,\mathbf{a}_1+y_6\,\mathbf{a}_2+z_6\,\mathbf{a}_3$	=	$\frac{1}{2}(x_6+y_6)\,a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}(-x_6+y_6)\,a\,\hat{\mathbf{y}}+z_6 c\,\hat{\mathbf{z}}$	$(6c)$	S III
$\mathbf{B}_{20}$	=	$-y_6\,\mathbf{a}_1+(x_6-y_6)\,\mathbf{a}_2+z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_6-y_6\right)a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}}+z_6 c\,\hat{\mathbf{z}}$	$(6c)$	S III
$\mathbf{B}_{21}$	=	$(-x_6+y_6)\,\mathbf{a}_1-x_6\,\mathbf{a}_2+z_6\,\mathbf{a}_3$	=	$\left(-x_6+\frac{1}{2}y_6\right)a\,\hat{\mathbf{x}}-\frac{\sqrt{3}}{2}y_6 a\,\hat{\mathbf{y}}+z_6 c\,\hat{\mathbf{z}}$	$(6c)$	S III

$$\mathbf{B}_{22} = -x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 = -\tfrac{1}{2}\,(x_6 + y_6)\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}\,(x_6 - y_6)\,a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right)c\,\hat{\mathbf{z}} \quad (6c) \quad \text{S III}$$

$$\mathbf{B}_{23} = y_6\,\mathbf{a}_1 + (-x_6 + y_6)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 = \left(-\tfrac{1}{2}x_6 + y_6\right)a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right)c\,\hat{\mathbf{z}} \quad (6c) \quad \text{S III}$$

$$\mathbf{B}_{24} = (x_6 - y_6)\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 = \left(x_6 - \tfrac{1}{2}y_6\right)a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}y_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right)c\,\hat{\mathbf{z}} \quad (6c) \quad \text{S III}$$

References:

- G. Collin and P. Laruelle, *Structure de $La_6Cu_2Si_2S_{14}$*, Bull. Soc. Fr. Mineral. Cristallogr. **94**, 175–176 (1971), doi:10.3406/bulmi.1971.6576.

Found in:

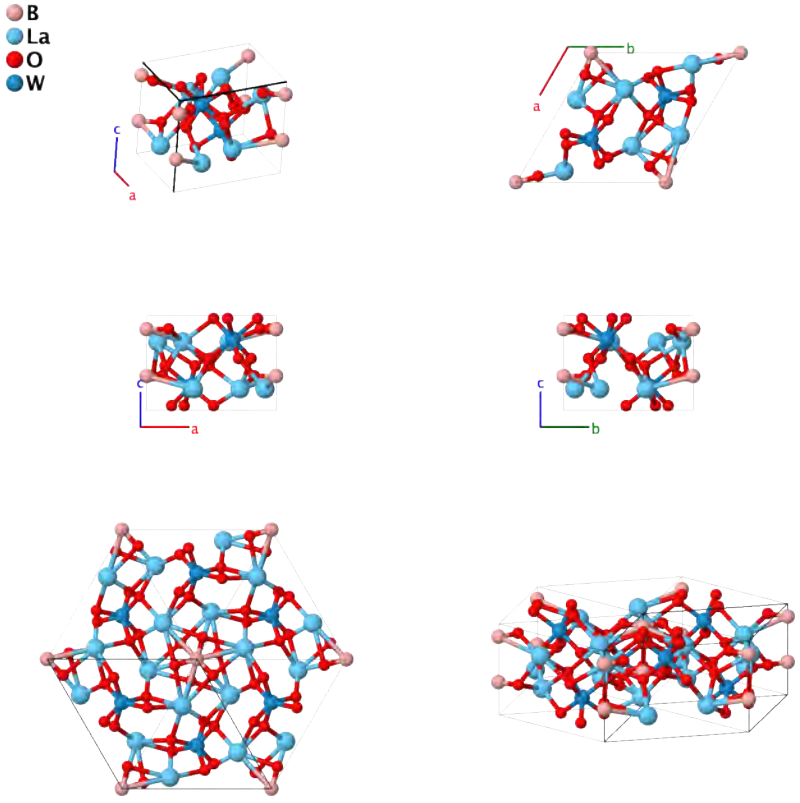
- L. D. Gulay, D. Kaczorowski, and A. Pietraszko, *Crystal structure and magnetic properties of $Ce_3CuSnSe_7$*, J. Alloys Compd. **403**, 49–52 (2005), doi:10.1016/j.jallcom.2005.05.030.

Geometry files:

- CIF: pp. 1748
- POSCAR: pp. 1748

La_3BWO_9 ($P6_3$) Structure: AB3C9D_hP28_173_a_c_3c_b

http://aflow.org/prototype-encyclopedia/AB3C9D_hP28_173_a_c_3c_b



Prototype	:	BLa_3O_9W
AFLOW prototype label	:	AB3C9D_hP28_173_a_c_3c_b
***Strukturbericht* designation**	:	None
Pearson symbol	:	hP28
Space group number	:	173
Space group symbol	:	$P6_3$
AFLOW prototype command	:	`aflow --proto=AB3C9D_hP28_173_a_c_3c_b --params=`$a, c/a, z_1, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5, x_6, y_6, z_6$

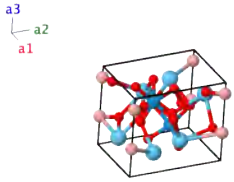
Other compounds with this structure

- La_3BWO_9, Ce_3BWO_9, Nd_3BWO_9, Sm_3BWO_9, Gd_3BWO_9, Tb_3BWO_9, Dy_3BWO_9, and Ho_3BWO_9

- Most refinements of the BLa_3O_9W structure, including (Ashtar, 2020) place it in hexagonal space group $P6_3$ #173. (Han, 2018) find a better fit to the data by refining it in the trigonal $P3$ #143 space group, which places the lanthanum atoms on two independent crystallographic sites. As this may be due to the presence of bismuth impurities, which (Han, 2018) place on the lanthanum site, while (Ashtar, 2020) claim to have very pure samples, we withhold judgment on which structure is correct and present both.
- Space group $P6_3$ does not specify the origin of the z-axis. Here it is set so that the coordinate of the tungsten atom is $z_2 = 1/4$.

Hexagonal primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$z_1\,\mathbf{a}_3$	=	$z_1 c\,\hat{\mathbf{z}}$	$(2a)$	B
$\mathbf{B}_2$	=	$\left(\frac{1}{2}+z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+z_1\right)c\,\hat{\mathbf{z}}$	$(2a)$	B
$\mathbf{B}_3$	=	$\frac{1}{3}\,\mathbf{a}_1+\frac{2}{3}\,\mathbf{a}_2+z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}+\frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+z_2 c\,\hat{\mathbf{z}}$	$(2b)$	W
$\mathbf{B}_4$	=	$\frac{2}{3}\,\mathbf{a}_1+\frac{1}{3}\,\mathbf{a}_2+\left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}-\frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_2\right)c\,\hat{\mathbf{z}}$	$(2b)$	W
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1+y_3\,\mathbf{a}_2+z_3\,\mathbf{a}_3$	=	$\frac{1}{2}(x_3+y_3)\,a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}(-x_3+y_3)\,a\,\hat{\mathbf{y}}+z_3 c\,\hat{\mathbf{z}}$	$(6c)$	La
$\mathbf{B}_6$	=	$-y_3\,\mathbf{a}_1+(x_3-y_3)\,\mathbf{a}_2+z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_3-y_3\right)a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}}+z_3 c\,\hat{\mathbf{z}}$	$(6c)$	La
$\mathbf{B}_7$	=	$(-x_3+y_3)\,\mathbf{a}_1-x_3\,\mathbf{a}_2+z_3\,\mathbf{a}_3$	=	$\left(-x_3+\frac{1}{2}y_3\right)a\,\hat{\mathbf{x}}-\frac{\sqrt{3}}{2}y_3 a\,\hat{\mathbf{y}}+z_3 c\,\hat{\mathbf{z}}$	$(6c)$	La
$\mathbf{B}_8$	=	$-x_3\,\mathbf{a}_1-y_3\,\mathbf{a}_2+\left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	=	$-\frac{1}{2}(x_3+y_3)\,a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}(x_3-y_3)\,a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_3\right)c\,\hat{\mathbf{z}}$	$(6c)$	La
$\mathbf{B}_9$	=	$y_3\,\mathbf{a}_1+(-x_3+y_3)\,\mathbf{a}_2+\left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_3+y_3\right)a\,\hat{\mathbf{x}}-\frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_3\right)c\,\hat{\mathbf{z}}$	$(6c)$	La
$\mathbf{B}_{10}$	=	$(x_3-y_3)\,\mathbf{a}_1+x_3\,\mathbf{a}_2+\left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	=	$\left(x_3-\frac{1}{2}y_3\right)a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}y_3 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_3\right)c\,\hat{\mathbf{z}}$	$(6c)$	La
$\mathbf{B}_{11}$	=	$x_4\,\mathbf{a}_1+y_4\,\mathbf{a}_2+z_4\,\mathbf{a}_3$	=	$\frac{1}{2}(x_4+y_4)\,a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}(-x_4+y_4)\,a\,\hat{\mathbf{y}}+z_4 c\,\hat{\mathbf{z}}$	$(6c)$	O I
$\mathbf{B}_{12}$	=	$-y_4\,\mathbf{a}_1+(x_4-y_4)\,\mathbf{a}_2+z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_4-y_4\right)a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}}+z_4 c\,\hat{\mathbf{z}}$	$(6c)$	O I
$\mathbf{B}_{13}$	=	$(-x_4+y_4)\,\mathbf{a}_1-x_4\,\mathbf{a}_2+z_4\,\mathbf{a}_3$	=	$\left(-x_4+\frac{1}{2}y_4\right)a\,\hat{\mathbf{x}}-\frac{\sqrt{3}}{2}y_4 a\,\hat{\mathbf{y}}+z_4 c\,\hat{\mathbf{z}}$	$(6c)$	O I
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1-y_4\,\mathbf{a}_2+\left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	=	$-\frac{1}{2}(x_4+y_4)\,a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}(x_4-y_4)\,a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_4\right)c\,\hat{\mathbf{z}}$	$(6c)$	O I
$\mathbf{B}_{15}$	=	$y_4\,\mathbf{a}_1+(-x_4+y_4)\,\mathbf{a}_2+\left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_4+y_4\right)a\,\hat{\mathbf{x}}-\frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_4\right)c\,\hat{\mathbf{z}}$	$(6c)$	O I
$\mathbf{B}_{16}$	=	$(x_4-y_4)\,\mathbf{a}_1+x_4\,\mathbf{a}_2+\left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	=	$\left(x_4-\frac{1}{2}y_4\right)a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}y_4 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_4\right)c\,\hat{\mathbf{z}}$	$(6c)$	O I
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1+y_5\,\mathbf{a}_2+z_5\,\mathbf{a}_3$	=	$\frac{1}{2}(x_5+y_5)\,a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}(-x_5+y_5)\,a\,\hat{\mathbf{y}}+z_5 c\,\hat{\mathbf{z}}$	$(6c)$	O II
$\mathbf{B}_{18}$	=	$-y_5\,\mathbf{a}_1+(x_5-y_5)\,\mathbf{a}_2+z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_5-y_5\right)a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}}+z_5 c\,\hat{\mathbf{z}}$	$(6c)$	O II
$\mathbf{B}_{19}$	=	$(-x_5+y_5)\,\mathbf{a}_1-x_5\,\mathbf{a}_2+z_5\,\mathbf{a}_3$	=	$\left(-x_5+\frac{1}{2}y_5\right)a\,\hat{\mathbf{x}}-\frac{\sqrt{3}}{2}y_5 a\,\hat{\mathbf{y}}+z_5 c\,\hat{\mathbf{z}}$	$(6c)$	O II
$\mathbf{B}_{20}$	=	$-x_5\,\mathbf{a}_1-y_5\,\mathbf{a}_2+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$-\frac{1}{2}(x_5+y_5)\,a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}(x_5-y_5)\,a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	$(6c)$	O II

$$\mathbf{B}_{21} = y_5\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_3 = \left(-\tfrac{1}{2}x_5 + y_5\right)a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_5\right)c\,\hat{\mathbf{z}} \quad (6c) \quad \text{O II}$$

$$\mathbf{B}_{22} = (x_5 - y_5)\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_3 = \left(x_5 - \tfrac{1}{2}y_5\right)a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}y_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_5\right)c\,\hat{\mathbf{z}} \quad (6c) \quad \text{O II}$$

$$\mathbf{B}_{23} = x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 = \tfrac{1}{2}(x_6 + y_6)\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}(-x_6 + y_6)\,a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}} \quad (6c) \quad \text{O III}$$

$$\mathbf{B}_{24} = -y_6\,\mathbf{a}_1 + (x_6 - y_6)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 = \left(\tfrac{1}{2}x_6 - y_6\right)a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}} \quad (6c) \quad \text{O III}$$

$$\mathbf{B}_{25} = (-x_6 + y_6)\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 = \left(-x_6 + \tfrac{1}{2}y_6\right)a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}y_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}} \quad (6c) \quad \text{O III}$$

$$\mathbf{B}_{26} = -x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 = -\tfrac{1}{2}(x_6 + y_6)\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}(x_6 - y_6)\,a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right)c\,\hat{\mathbf{z}} \quad (6c) \quad \text{O III}$$

$$\mathbf{B}_{27} = y_6\,\mathbf{a}_1 + (-x_6 + y_6)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 = \left(-\tfrac{1}{2}x_6 + y_6\right)a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right)c\,\hat{\mathbf{z}} \quad (6c) \quad \text{O III}$$

$$\mathbf{B}_{28} = (x_6 - y_6)\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 = \left(x_6 - \tfrac{1}{2}y_6\right)a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}y_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right)c\,\hat{\mathbf{z}} \quad (6c) \quad \text{O III}$$

References:

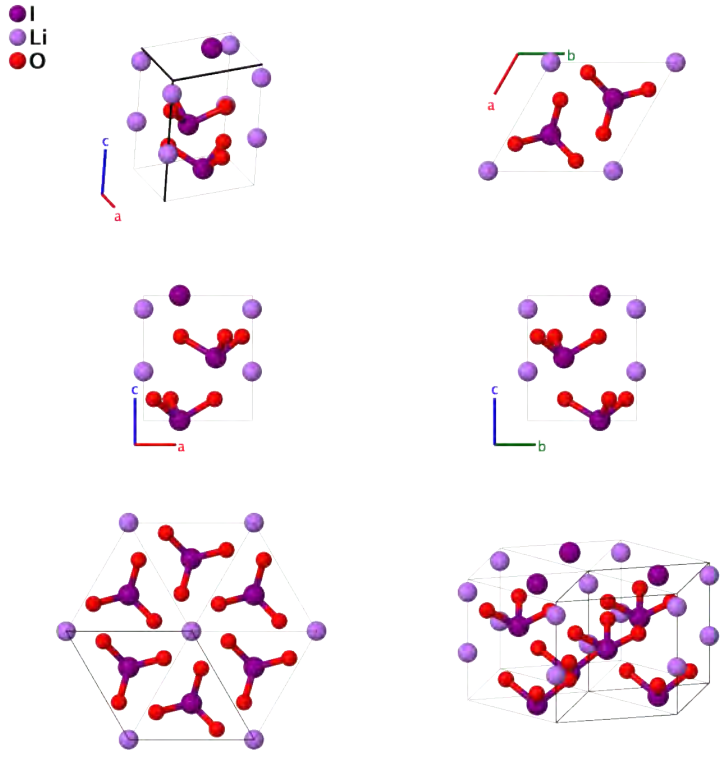
- J. Han, F. Pan, M. S. Molokeev, J. Dai, M. Peng, W. Zhou, and J. Wang, *Redefinition of Crystal Structure and Bi^{3+} Yellow Luminescence with Strong Near-Ultraviolet Excitation in La_3BWO_9:Bi^{3+} Phosphor for White Light-Emitting Diodes*, ACS Appl. Mater. Interfaces **10**, 13660–13668 (2018), doi:10.1021/acsami.8b00808.
- M. Ashtar, J. Guo, Z. Wan, Y. Wang, G. Gong, Y. Liu, Y. Su, and Z. Tian, *A new family of disorder-free Rare-Earth-based kagomé lattice magnets: structure and magnetic characterizations of RE_3BWO_9 (RE = Pr, Nd, Gd-Ho) Boratotungstates*, http://arxiv.org/abs/2002.05420. ArXiv:2002.05420 [cond-mat.mtrl-sci].

Geometry files:

- CIF: pp. 1749
- POSCAR: pp. 1749

α-$LiIO_3$ Structure: ABC3_hP10_173_b_a_c

http://aflow.org/prototype-encyclopedia/ABC3_hP10_173_b_a_c

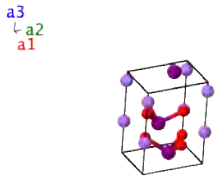


Prototype	:	$ILiO_3$
AFLOW prototype label	:	ABC3_hP10_173_b_a_c
***Strukturbericht* designation**	:	None
Pearson symbol	:	hP10
Space group number	:	173
Space group symbol	:	$P6_3$
AFLOW prototype command	:	`aflow --proto=ABC3_hP10_173_b_a_c --params=`$a, c/a, z_1, z_2, x_3, y_3, z_3$

- $LiIO_3$ is known to exist in three forms:
  - α-$LiIO_3$, stable below 470 K: (Zachariasen, 1931) originally determined that the structure of α-$LiIO_3$ was in space group $P6_322$ #182, which (Hermann, 1937) designated *Strukturbericht* $E2_3$. (Rosenzweig, 1966) subsequently determined that this structure was incorrect because of the small sample size, and determined that the true structure was in space group $P6_3$ #173.
  - β-$LiIO_3$, stable from 573 up to the melting point at 708 K.
  - γ-$LiIO_3$, stable between the α- and β-phases, with an orthorhombic structure in space group $Pna2_1$ #33.

Hexagonal primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$z_1\,\mathbf{a}_3$	=	$z_1 c\,\hat{\mathbf{z}}$	$(2a)$	Li
$\mathbf{B}_2$	=	$\left(\tfrac{1}{2}+z_1\right)\mathbf{a}_3$	=	$\left(\tfrac{1}{2}+z_1\right)c\,\hat{\mathbf{z}}$	$(2a)$	Li
$\mathbf{B}_3$	=	$\tfrac{1}{3}\,\mathbf{a}_1+\tfrac{2}{3}\,\mathbf{a}_2+z_2\,\mathbf{a}_3$	=	$\tfrac{1}{2}a\,\hat{\mathbf{x}}+\tfrac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+z_2 c\,\hat{\mathbf{z}}$	$(2b)$	I
$\mathbf{B}_4$	=	$\tfrac{2}{3}\,\mathbf{a}_1+\tfrac{1}{3}\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_2\right)\mathbf{a}_3$	=	$\tfrac{1}{2}a\,\hat{\mathbf{x}}-\tfrac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_2\right)c\,\hat{\mathbf{z}}$	$(2b)$	I
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1+y_3\,\mathbf{a}_2+z_3\,\mathbf{a}_3$	=	$\tfrac{1}{2}\,(x_3+y_3)\,a\,\hat{\mathbf{x}}+\tfrac{\sqrt{3}}{2}\,(-x_3+y_3)\,a\,\hat{\mathbf{y}}+z_3 c\,\hat{\mathbf{z}}$	$(6c)$	O
$\mathbf{B}_6$	=	$-y_3\,\mathbf{a}_1+(x_3-y_3)\,\mathbf{a}_2+z_3\,\mathbf{a}_3$	=	$\left(\tfrac{1}{2}x_3-y_3\right)a\,\hat{\mathbf{x}}+\tfrac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}}+z_3 c\,\hat{\mathbf{z}}$	$(6c)$	O
$\mathbf{B}_7$	=	$(-x_3+y_3)\,\mathbf{a}_1-x_3\,\mathbf{a}_2+z_3\,\mathbf{a}_3$	=	$\left(-x_3+\tfrac{1}{2}y_3\right)a\,\hat{\mathbf{x}}-\tfrac{\sqrt{3}}{2}y_3 a\,\hat{\mathbf{y}}+z_3 c\,\hat{\mathbf{z}}$	$(6c)$	O
$\mathbf{B}_8$	=	$-x_3\,\mathbf{a}_1-y_3\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_3\right)\mathbf{a}_3$	=	$-\tfrac{1}{2}\,(x_3+y_3)\,a\,\hat{\mathbf{x}}+\tfrac{\sqrt{3}}{2}\,(x_3-y_3)\,a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_3\right)c\,\hat{\mathbf{z}}$	$(6c)$	O
$\mathbf{B}_9$	=	$y_3\,\mathbf{a}_1+(-x_3+y_3)\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_3\right)\mathbf{a}_3$	=	$\left(-\tfrac{1}{2}x_3+y_3\right)a\,\hat{\mathbf{x}}-\tfrac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_3\right)c\,\hat{\mathbf{z}}$	$(6c)$	O
$\mathbf{B}_{10}$	=	$(x_3-y_3)\,\mathbf{a}_1+x_3\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_3\right)\mathbf{a}_3$	=	$\left(x_3-\tfrac{1}{2}y_3\right)a\,\hat{\mathbf{x}}+\tfrac{\sqrt{3}}{2}y_3 a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_3\right)c\,\hat{\mathbf{z}}$	$(6c)$	O

References:

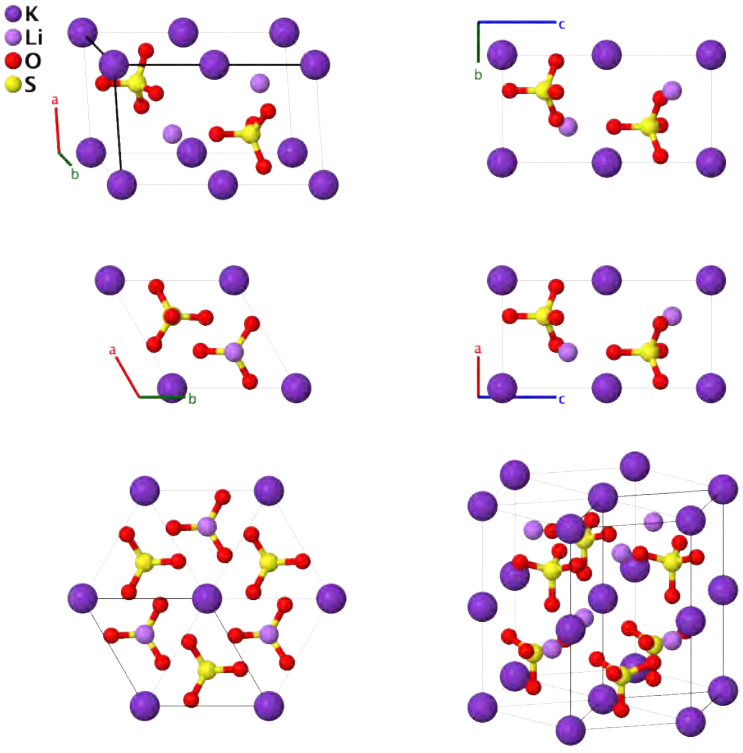
- A. Rosenzweig and B. Morosin, *A reinvestigation of the crystal structure of $LiIO_3$*, Acta Cryst. **20**, 758–761 (1966), doi:10.1107/S0365110X66001804.
- W. H. Zachariasen and F. A. Barta, *Crystal Structure of Lithium Iodate*, Phys. Rev. **37**, 1626–1630 (1931), doi:10.1103/PhysRev.37.1626.
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1749
- POSCAR: pp. 1749

$LiKSO_4$ ($H1_4$) Structure: ABC4D_hP14_173_a_b_bc_b

http://aflow.org/prototype-encyclopedia/ABC4D_hP14_173_a_b_bc_b

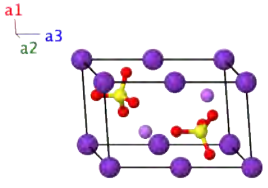


Prototype	:	$KLiO_4S$
AFLOW prototype label	:	ABC4D_hP14_173_a_b_bc_b
***Strukturbericht* designation**	:	$H1_4$
Pearson symbol	:	hP14
Space group number	:	173
Space group symbol	:	$P6_3$
AFLOW prototype command	:	`aflow --proto=ABC4D_hP14_173_a_b_bc_b --params=`$a, c/a, z_1, z_2, z_3, z_4, x_5, y_5, z_5$

- As there is no center of symmetry in the structure, the point $z = 0$ can be set arbitrarily. Following (Bhakay-Tamhane, 1984) we set $z_1 = 0$.
- While *Strukturbericht* symbols $H1_1$, $H1_2$ and $H1_5$ are identical with $S1_1$, $S1_2$ and $S1_5$, respectively, there is no connection between this structure and $S1_4$, garnet (A2B3C12D3_cI160_230_a_c_h_d).

Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$z_1\,\mathbf{a}_3$	=	$z_1 c\,\hat{\mathbf{z}}$	($2a$)	K
$\mathbf{B}_2$	=	$\left(\frac{1}{2}+z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+z_1\right)c\,\hat{\mathbf{z}}$	($2a$)	K
$\mathbf{B}_3$	=	$\frac{1}{3}\,\mathbf{a}_1+\frac{2}{3}\,\mathbf{a}_2+z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}+\frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+z_2 c\,\hat{\mathbf{z}}$	($2b$)	Li
$\mathbf{B}_4$	=	$\frac{2}{3}\,\mathbf{a}_1+\frac{1}{3}\,\mathbf{a}_2+\left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}-\frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_2\right)c\,\hat{\mathbf{z}}$	($2b$)	Li
$\mathbf{B}_5$	=	$\frac{1}{3}\,\mathbf{a}_1+\frac{2}{3}\,\mathbf{a}_2+z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}+\frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+z_3 c\,\hat{\mathbf{z}}$	($2b$)	O I
$\mathbf{B}_6$	=	$\frac{2}{3}\,\mathbf{a}_1+\frac{1}{3}\,\mathbf{a}_2+\left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}-\frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_3\right)c\,\hat{\mathbf{z}}$	($2b$)	O I
$\mathbf{B}_7$	=	$\frac{1}{3}\,\mathbf{a}_1+\frac{2}{3}\,\mathbf{a}_2+z_4\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}+\frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+z_4 c\,\hat{\mathbf{z}}$	($2b$)	S
$\mathbf{B}_8$	=	$\frac{2}{3}\,\mathbf{a}_1+\frac{1}{3}\,\mathbf{a}_2+\left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}-\frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_4\right)c\,\hat{\mathbf{z}}$	($2b$)	S
$\mathbf{B}_9$	=	$x_5\,\mathbf{a}_1+y_5\,\mathbf{a}_2+z_5\,\mathbf{a}_3$	=	$\frac{1}{2}\left(x_5+y_5\right)a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}\left(-x_5+y_5\right)a\,\hat{\mathbf{y}}+z_5 c\,\hat{\mathbf{z}}$	($6c$)	O II
$\mathbf{B}_{10}$	=	$-y_5\,\mathbf{a}_1+\left(x_5-y_5\right)\mathbf{a}_2+z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_5-y_5\right)a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}}+z_5 c\,\hat{\mathbf{z}}$	($6c$)	O II
$\mathbf{B}_{11}$	=	$\left(-x_5+y_5\right)\mathbf{a}_1-x_5\,\mathbf{a}_2+z_5\,\mathbf{a}_3$	=	$\left(-x_5+\frac{1}{2}y_5\right)a\,\hat{\mathbf{x}}-\frac{\sqrt{3}}{2}y_5 a\,\hat{\mathbf{y}}+z_5 c\,\hat{\mathbf{z}}$	($6c$)	O II
$\mathbf{B}_{12}$	=	$-x_5\,\mathbf{a}_1-y_5\,\mathbf{a}_2+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$-\frac{1}{2}\left(x_5+y_5\right)a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}\left(x_5-y_5\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	($6c$)	O II
$\mathbf{B}_{13}$	=	$y_5\,\mathbf{a}_1+\left(-x_5+y_5\right)\mathbf{a}_2+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_5+y_5\right)a\,\hat{\mathbf{x}}-\frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	($6c$)	O II
$\mathbf{B}_{14}$	=	$\left(x_5-y_5\right)\mathbf{a}_1+x_5\,\mathbf{a}_2+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$\left(x_5-\frac{1}{2}y_5\right)a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}y_5 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	($6c$)	O II

References:

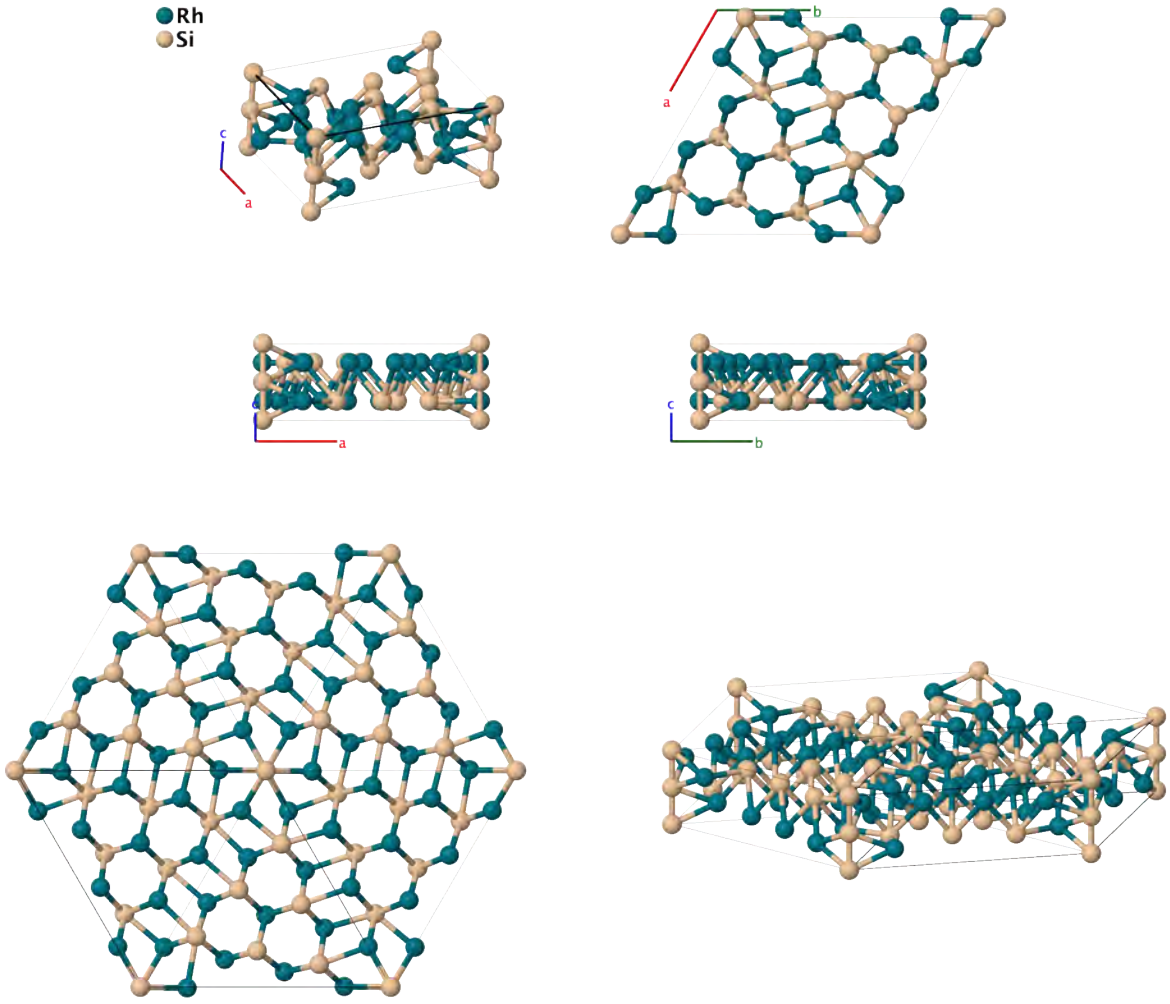
- S. Bhakay-Tamhane, A. Sequiera, and R. Chidambaram, *Structure of lithium potassium sulphate, $LiKSO_4$: a neutron diffraction study*, Acta Crystallogr. C **40**, 1648–1651 (1984), doi:10.1107/S0108270184009045.

Geometry files:

- CIF: pp. 1750
- POSCAR: pp. 1750

$Rh_{20}Si_{13}$ Structure: A10B7_hP34_176_c3h_b2h

http://aflow.org/prototype-encyclopedia/A10B7_hP34_176_c3h_b2h

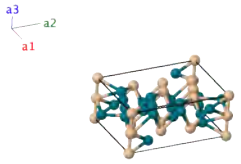


Prototype	:	$Rh_{20}Si_{13}$
AFLOW prototype label	:	A10B7_hP34_176_c3h_b2h
***Strukturbericht* designation**	:	None
Pearson symbol	:	hP34
Space group number	:	176
Space group symbol	:	$P6_3/m$
AFLOW prototype command	:	`aflow --proto=A10B7_hP34_176_c3h_b2h --params=`$a, c/a, x_3, y_3, x_4, y_4, x_5, y_5, x_6, y_6, x_7, y_7$

- The Si-I ($2b$) site is only occupied 50% of the time, which gives the observed stoichiometry.

Hexagonal primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(2*b*)	Si I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	(2*b*)	Si I
$\mathbf{B}_3$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(2*c*)	Rh I
$\mathbf{B}_4$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(2*c*)	Rh I
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_3 + y_3)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_3 + y_3)\,a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Rh II
$\mathbf{B}_6$	=	$-y_3\,\mathbf{a}_1 + (x_3 - y_3)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_3 - y_3\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Rh II
$\mathbf{B}_7$	=	$(-x_3 + y_3)\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(-x_3 + \frac{1}{2}y_3\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_3a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Rh II
$\mathbf{B}_8$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-\frac{1}{2}\,(x_3 + y_3)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(x_3 - y_3)\,a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Rh II
$\mathbf{B}_9$	=	$y_3\,\mathbf{a}_1 + (-x_3 + y_3)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_3 + y_3\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Rh II
$\mathbf{B}_{10}$	=	$(x_3 - y_3)\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(x_3 - \frac{1}{2}y_3\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_3a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Rh II
$\mathbf{B}_{11}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_4 + y_4)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_4 + y_4)\,a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Rh III
$\mathbf{B}_{12}$	=	$-y_4\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_4 - y_4\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Rh III
$\mathbf{B}_{13}$	=	$(-x_4 + y_4)\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(-x_4 + \frac{1}{2}y_4\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_4a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Rh III
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-\frac{1}{2}\,(x_4 + y_4)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(x_4 - y_4)\,a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Rh III
$\mathbf{B}_{15}$	=	$y_4\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_4 + y_4\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Rh III
$\mathbf{B}_{16}$	=	$(x_4 - y_4)\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(x_4 - \frac{1}{2}y_4\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_4a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Rh III
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_5 + y_5)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_5 + y_5)\,a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Rh IV
$\mathbf{B}_{18}$	=	$-y_5\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_5 - y_5\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_5a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Rh IV
$\mathbf{B}_{19}$	=	$(-x_5 + y_5)\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(-x_5 + \frac{1}{2}y_5\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_5a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Rh IV
$\mathbf{B}_{20}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-\frac{1}{2}\,(x_5 + y_5)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(x_5 - y_5)\,a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Rh IV
$\mathbf{B}_{21}$	=	$y_5\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_5 + y_5\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_5a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Rh IV
$\mathbf{B}_{22}$	=	$(x_5 - y_5)\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(x_5 - \frac{1}{2}y_5\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_5a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Rh IV
$\mathbf{B}_{23}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_6 + y_6)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_6 + y_6)\,a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Si II
$\mathbf{B}_{24}$	=	$-y_6\,\mathbf{a}_1 + (x_6 - y_6)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_6 - y_6\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_6a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Si II
$\mathbf{B}_{25}$	=	$(-x_6 + y_6)\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(-x_6 + \frac{1}{2}y_6\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_6a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Si II
$\mathbf{B}_{26}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-\frac{1}{2}\,(x_6 + y_6)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(x_6 - y_6)\,a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Si II
$\mathbf{B}_{27}$	=	$y_6\,\mathbf{a}_1 + (-x_6 + y_6)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_6 + y_6\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_6a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Si II
$\mathbf{B}_{28}$	=	$(x_6 - y_6)\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(x_6 - \frac{1}{2}y_6\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_6a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Si II

$\mathbf{B}_{29}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_7 + y_7)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_7 + y_7)\,a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Si III
$\mathbf{B}_{30}$	=	$-y_7\,\mathbf{a}_1 + (x_7 - y_7)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_7 - y_7\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_7 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Si III
$\mathbf{B}_{31}$	=	$(-x_7 + y_7)\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(-x_7 + \frac{1}{2}y_7\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_7 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Si III
$\mathbf{B}_{32}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-\frac{1}{2}\,(x_7 + y_7)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(x_7 - y_7)\,a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Si III
$\mathbf{B}_{33}$	=	$y_7\,\mathbf{a}_1 + (-x_7 + y_7)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_7 + y_7\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_7 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Si III
$\mathbf{B}_{34}$	=	$(x_7 - y_7)\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(x_7 - \frac{1}{2}y_7\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_7 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Si III

References:

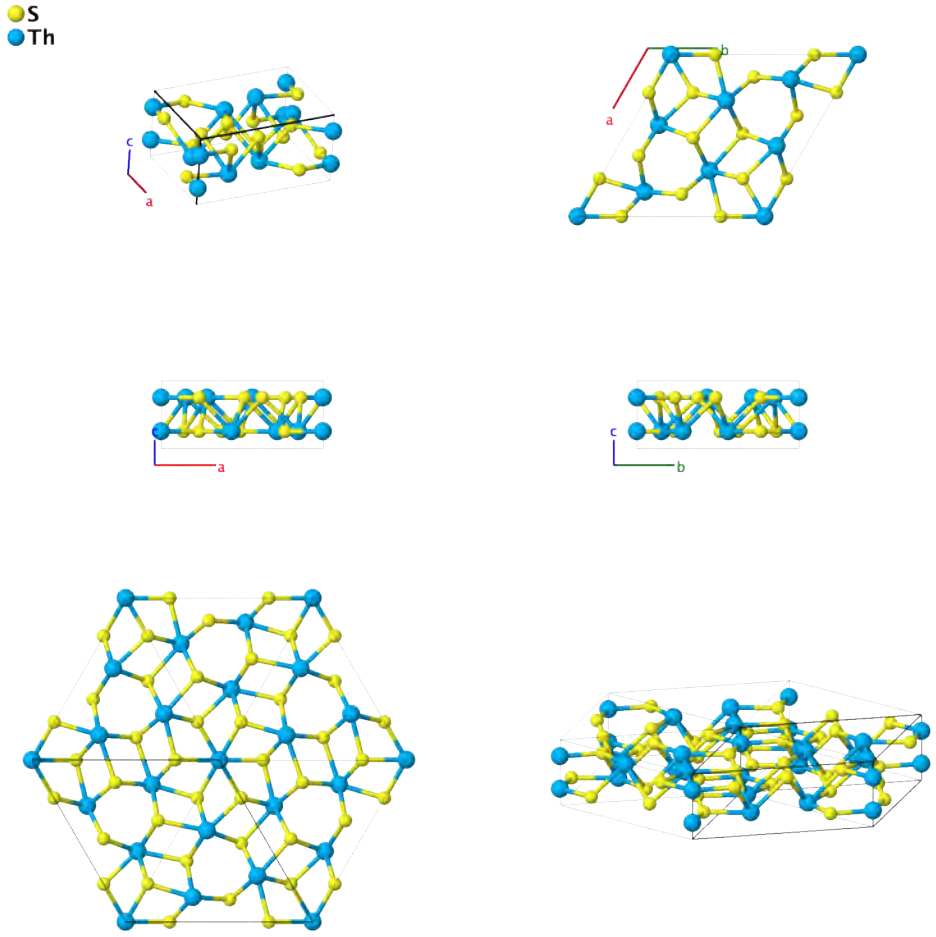
- I. Engström, *The Crystal Structure of $Rh_{20}Si_{13}$*, Acta Chem. Scand. **19**, 1924–1932 (1965), doi:10.3891/acta.chem.scand.19-1924.

Geometry files:

- CIF: pp. 1750
- POSCAR: pp. 1750

Th_7S_{12} ($D8_k$) Structure: A3B2_hP20_176_2h_ah

http://aflow.org/prototype-encyclopedia/A3B2_hP20_176_2h_ah



Prototype	:	$S_{12}Th_7$
AFLOW prototype label	:	A3B2_hP20_176_2h_ah
***Strukturbericht* designation**	:	$D8_k$
Pearson symbol	:	hP20
Space group number	:	176
Space group symbol	:	$P6_3/m$
AFLOW prototype command	:	`aflow --proto=A3B2_hP20_176_2h_ah --params=`$a, c/a, x_2, y_2, x_3, y_3, x_4, y_4$

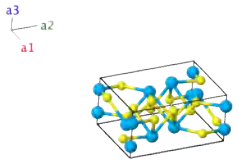
Other compounds with this structure

- Th_7Se_{12} and $(Ga,As)_7Pd_{12}$

- The Th ($2a$) site is half-filled. The stoichiometry for the AFLOW label treats the Th site as fully occupied.

Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$\frac{1}{4}\,\mathbf{a}_3$	$\frac{1}{4}c\,\hat{\mathbf{z}}$	(2*a*)	Th I
$\mathbf{B}_2$	$\frac{3}{4}\,\mathbf{a}_3$	$\frac{3}{4}c\,\hat{\mathbf{z}}$	(2*a*)	Th I
$\mathbf{B}_3$	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	$\frac{1}{2}\,(x_2 + y_2)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_2 + y_2)\,a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	S I
$\mathbf{B}_4$	$-y_2\,\mathbf{a}_1 + (x_2 - y_2)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	$\left(\frac{1}{2}x_2 - y_2\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_2 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	S I
$\mathbf{B}_5$	$(-x_2 + y_2)\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	$\left(-x_2 + \frac{1}{2}y_2\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_2 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	S I
$\mathbf{B}_6$	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	$-\frac{1}{2}\,(x_2 + y_2)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(x_2 - y_2)\,a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	S I
$\mathbf{B}_7$	$y_2\,\mathbf{a}_1 + (-x_2 + y_2)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	$\left(-\frac{1}{2}x_2 + y_2\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_2 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	S I
$\mathbf{B}_8$	$(x_2 - y_2)\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	$\left(x_2 - \frac{1}{2}y_2\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_2 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	S I
$\mathbf{B}_9$	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	$\frac{1}{2}\,(x_3 + y_3)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_3 + y_3)\,a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	S II
$\mathbf{B}_{10}$	$-y_3\,\mathbf{a}_1 + (x_3 - y_3)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	$\left(\frac{1}{2}x_3 - y_3\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	S II
$\mathbf{B}_{11}$	$(-x_3 + y_3)\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	$\left(-x_3 + \frac{1}{2}y_3\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_3 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	S II
$\mathbf{B}_{12}$	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	$-\frac{1}{2}\,(x_3 + y_3)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(x_3 - y_3)\,a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	S II
$\mathbf{B}_{13}$	$y_3\,\mathbf{a}_1 + (-x_3 + y_3)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	$\left(-\frac{1}{2}x_3 + y_3\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	S II
$\mathbf{B}_{14}$	$(x_3 - y_3)\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	$\left(x_3 - \frac{1}{2}y_3\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_3 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	S II
$\mathbf{B}_{15}$	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	$\frac{1}{2}\,(x_4 + y_4)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_4 + y_4)\,a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Th II
$\mathbf{B}_{16}$	$-y_4\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	$\left(\frac{1}{2}x_4 - y_4\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Th II
$\mathbf{B}_{17}$	$(-x_4 + y_4)\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	$\left(-x_4 + \frac{1}{2}y_4\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_4 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Th II
$\mathbf{B}_{18}$	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	$-\frac{1}{2}\,(x_4 + y_4)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(x_4 - y_4)\,a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Th II
$\mathbf{B}_{19}$	$y_4\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	$\left(-\frac{1}{2}x_4 + y_4\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Th II
$\mathbf{B}_{20}$	$(x_4 - y_4)\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	$\left(x_4 - \frac{1}{2}y_4\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_4 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Th II

References:

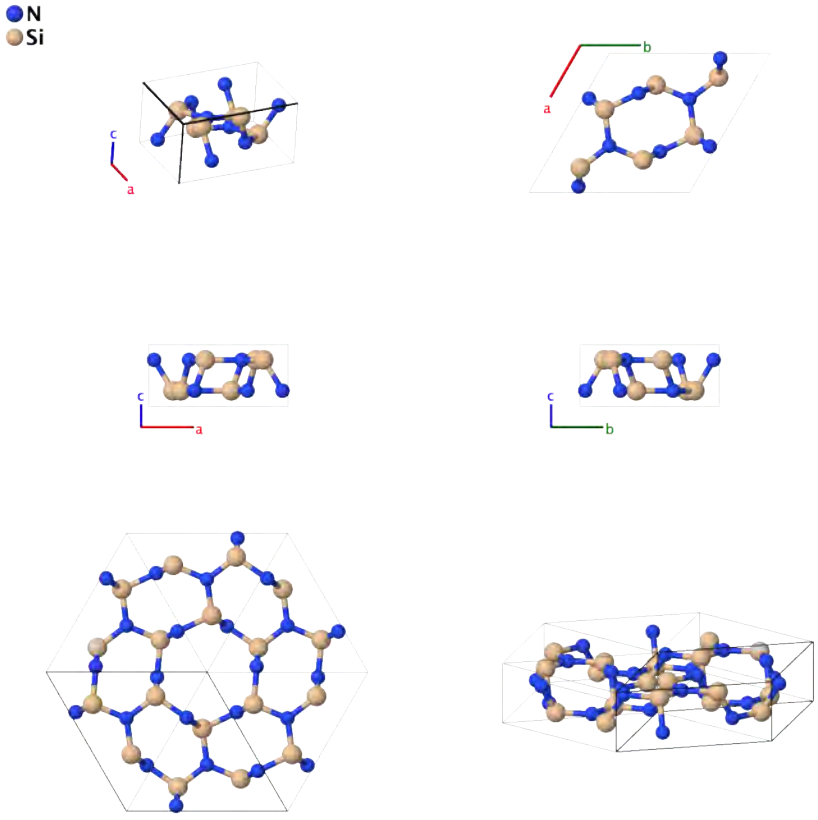
- W. H. Zachariasen, *Crystal chemical studies of the 5f-series of elements. IX. The crystal structure of* Th_7S_{12}, Acta Cryst. **2**, 288–291 (1949), doi:10.1107/S0365110X49000746.

Geometry files:

- CIF: pp. 1751
- POSCAR: pp. 1751

β-Si_3N_4 Structure: A4B3_hP14_176_ch_h

http://aflow.org/prototype-encyclopedia/A4B3_hP14_176_ch_h

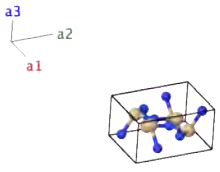


Prototype	:	N_4Si_3
AFLOW prototype label	:	A4B3_hP14_176_ch_h
***Strukturbericht* designation**	:	None
Pearson symbol	:	hP14
Space group number	:	176
Space group symbol	:	$P6_3/m$
AFLOW prototype command	:	`aflow --proto=A4B3_hP14_176_ch_h --params=`$a, c/a, x_2, y_2, x_3, y_3$

- (Grün, 1979) places this structure in space group $P6_3$ #173. His structure is nearly indistinguishable from this one.

Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(2*c*)	N I
$\mathbf{B}_2$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(2*c*)	N I
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_2 + y_2)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_2 + y_2)\,a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	N II
$\mathbf{B}_4$	=	$-y_2\,\mathbf{a}_1 + (x_2 - y_2)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_2 - y_2\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_2 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	N II
$\mathbf{B}_5$	=	$(-x_2 + y_2)\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(-x_2 + \frac{1}{2}y_2\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_2 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	N II
$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-\frac{1}{2}\,(x_2 + y_2)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(x_2 - y_2)\,a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	N II
$\mathbf{B}_7$	=	$y_2\,\mathbf{a}_1 + (-x_2 + y_2)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_2 + y_2\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_2 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	N II
$\mathbf{B}_8$	=	$(x_2 - y_2)\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(x_2 - \frac{1}{2}y_2\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_2 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	N II
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_3 + y_3)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_3 + y_3)\,a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Si
$\mathbf{B}_{10}$	=	$-y_3\,\mathbf{a}_1 + (x_3 - y_3)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_3 - y_3\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Si
$\mathbf{B}_{11}$	=	$(-x_3 + y_3)\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(-x_3 + \frac{1}{2}y_3\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_3 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Si
$\mathbf{B}_{12}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-\frac{1}{2}\,(x_3 + y_3)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(x_3 - y_3)\,a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Si
$\mathbf{B}_{13}$	=	$y_3\,\mathbf{a}_1 + (-x_3 + y_3)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_3 + y_3\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Si
$\mathbf{B}_{14}$	=	$(x_3 - y_3)\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(x_3 - \frac{1}{2}y_3\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_3 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	Si

References:

- P. Yang, H.-K. Fun, I. Ab. Rahman, and M. I. Saleh, *Two phase refinements of the structures of α-Si_3N_4 and β-Si_3N_4 made from rice husk by Rietveld analysis*, Ceram. Int. **21**, 137–142 (1995), doi:10.1016/0272-8842(95)95885-L.
- R. Grün, *The crystal structure of β-Si_3N_4: structural and stability considerations between α- and β-Si_3N_4*, Acta Crystallogr. Sect. B Struct. Sci. **35**, 800–804 (1979), doi:10.1107/S0567740879004933.

Found in:

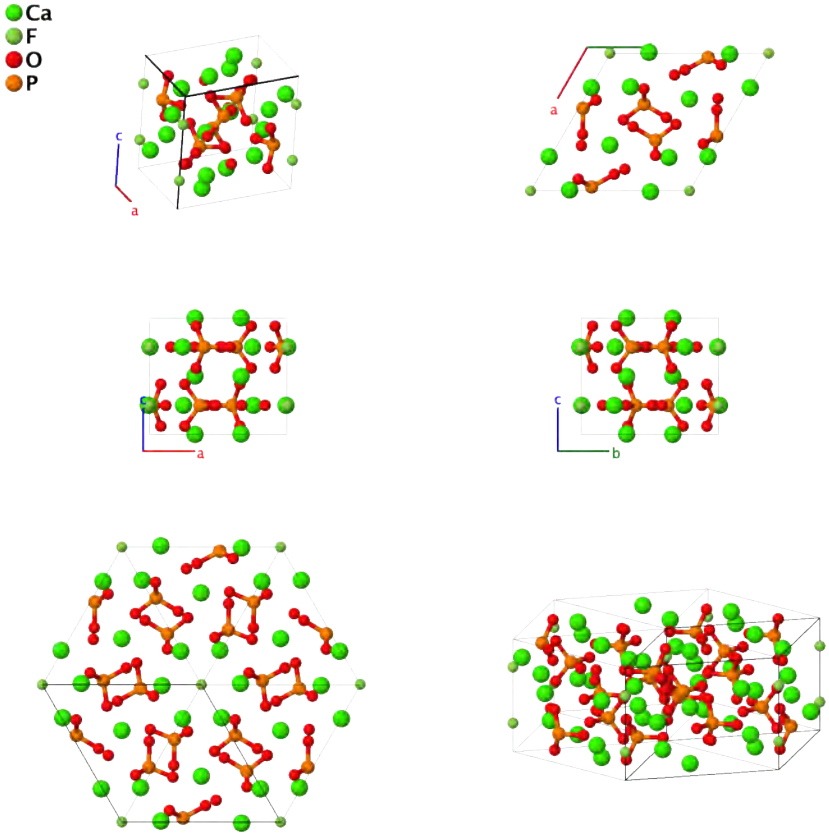
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1751
- POSCAR: pp. 1751

Fluorapatite [$Ca_5F(PO_4)_3$, $H5_7$] Structure: A5BC12D3_hP42_176_fh_a_2hi_h

http://aflow.org/prototype-encyclopedia/A5BC12D3_hP42_176_fh_a_2hi_h



Prototype	:	$Ca_5FO_{12}P_3$
AFLOW prototype label	:	A5BC12D3_hP42_176_fh_a_2hi_h
***Strukturbericht* designation**	:	$H5_7$
Pearson symbol	:	hP42
Space group number	:	176
Space group symbol	:	$P6_3/m$
AFLOW prototype command	:	`aflow --proto=A5BC12D3_hP42_176_fh_a_2hi_h --params=`$a, c/a, z_2, x_3, y_3, x_4, y_4, x_5, y_5, x_6, y_6, x_7, y_7, z_7$

Other compounds with this structure

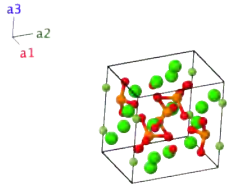
- $Ca_5OH(PO_4)_3$ (hydroxylapatite) and $Ca_5Cl(PO_4)_3$ (chlorapatite)

- Apatite can be formed with most M^{2+} metallic ions replacing the calcium, and many ions (AsO_3, CO_3, Si_3, *etc.*) replacing the phosphate. While these structures are related to the prototype, they may have slight changes in crystal structure. The phosphate apatites are the main source of phosphorus on Earth (Hughes, 2002).

- When OH or Cl replaces F, that ion is displaced from the ($2a$) position to the ($4e$) position, with $z = 0.1979$ for OH and 0.4323 for Cl. The substitute ion fills half of the ($4f$) sites.

Hexagonal primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\,\hat{\mathbf{z}}$	$(2a)$	F
$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\,\hat{\mathbf{z}}$	$(2a)$	F
$\mathbf{B}_3$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(4f)$	Ca I
$\mathbf{B}_4$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	$(4f)$	Ca I
$\mathbf{B}_5$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	$(4f)$	Ca I
$\mathbf{B}_6$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	$(4f)$	Ca I
$\mathbf{B}_7$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_3 + y_3)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_3 + y_3)\,a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	Ca II
$\mathbf{B}_8$	=	$-y_3\,\mathbf{a}_1 + (x_3 - y_3)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_3 - y_3\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	Ca II
$\mathbf{B}_9$	=	$(-x_3 + y_3)\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(-x_3 + \frac{1}{2}y_3\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_3 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	Ca II
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-\frac{1}{2}\,(x_3 + y_3)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(x_3 - y_3)\,a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	Ca II
$\mathbf{B}_{11}$	=	$y_3\,\mathbf{a}_1 + (-x_3 + y_3)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_3 + y_3\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	Ca II
$\mathbf{B}_{12}$	=	$(x_3 - y_3)\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(x_3 - \frac{1}{2}y_3\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_3 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	Ca II
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_4 + y_4)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_4 + y_4)\,a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O I
$\mathbf{B}_{14}$	=	$-y_4\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_4 - y_4\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O I
$\mathbf{B}_{15}$	=	$(-x_4 + y_4)\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(-x_4 + \frac{1}{2}y_4\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_4 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O I
$\mathbf{B}_{16}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-\frac{1}{2}\,(x_4 + y_4)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(x_4 - y_4)\,a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O I
$\mathbf{B}_{17}$	=	$y_4\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_4 + y_4\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O I
$\mathbf{B}_{18}$	=	$(x_4 - y_4)\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(x_4 - \frac{1}{2}y_4\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_4 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O I
$\mathbf{B}_{19}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_5 + y_5)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_5 + y_5)\,a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O II
$\mathbf{B}_{20}$	=	$-y_5\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_5 - y_5\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O II
$\mathbf{B}_{21}$	=	$(-x_5 + y_5)\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(-x_5 + \frac{1}{2}y_5\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_5 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O II
$\mathbf{B}_{22}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-\frac{1}{2}\,(x_5 + y_5)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(x_5 - y_5)\,a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O II
$\mathbf{B}_{23}$	=	$y_5\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_5 + y_5\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O II
$\mathbf{B}_{24}$	=	$(x_5 - y_5)\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(x_5 - \frac{1}{2}y_5\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_5 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O II

$\mathbf{B}_{25}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_6 + y_6)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_6 + y_6)\,a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	P
$\mathbf{B}_{26}$	=	$-y_6\,\mathbf{a}_1 + (x_6 - y_6)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_6 - y_6\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	P
$\mathbf{B}_{27}$	=	$(-x_6 + y_6)\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(-x_6 + \frac{1}{2}y_6\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_6 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	P
$\mathbf{B}_{28}$	=	$-x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-\frac{1}{2}\,(x_6 + y_6)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(x_6 - y_6)\,a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	P
$\mathbf{B}_{29}$	=	$y_6\,\mathbf{a}_1 + (-x_6 + y_6)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_6 + y_6\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	P
$\mathbf{B}_{30}$	=	$(x_6 - y_6)\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(x_6 - \frac{1}{2}y_6\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_6 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	P
$\mathbf{B}_{31}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_7 + y_7)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_7 + y_7)\,a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(12*i*)	O III
$\mathbf{B}_{32}$	=	$-y_7\,\mathbf{a}_1 + (x_7 - y_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_7 - y_7\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_7 a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(12*i*)	O III
$\mathbf{B}_{33}$	=	$(-x_7 + y_7)\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$\left(-x_7 + \frac{1}{2}y_7\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_7 a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(12*i*)	O III
$\mathbf{B}_{34}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$-\frac{1}{2}\,(x_7 + y_7)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(x_7 - y_7)\,a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right)c\,\hat{\mathbf{z}}$	(12*i*)	O III
$\mathbf{B}_{35}$	=	$y_7\,\mathbf{a}_1 + (-x_7 + y_7)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_7 + y_7\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_7 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right)c\,\hat{\mathbf{z}}$	(12*i*)	O III
$\mathbf{B}_{36}$	=	$(x_7 - y_7)\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\mathbf{a}_3$	=	$\left(x_7 - \frac{1}{2}y_7\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_7 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right)c\,\hat{\mathbf{z}}$	(12*i*)	O III
$\mathbf{B}_{37}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$-\frac{1}{2}\,(x_7 + y_7)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(x_7 - y_7)\,a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	(12*i*)	O III
$\mathbf{B}_{38}$	=	$y_7\,\mathbf{a}_1 + (-x_7 + y_7)\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_7 + y_7\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_7 a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	(12*i*)	O III
$\mathbf{B}_{39}$	=	$(x_7 - y_7)\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$\left(x_7 - \frac{1}{2}y_7\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_7 a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	(12*i*)	O III
$\mathbf{B}_{40}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\mathbf{a}_3$	=	$\frac{1}{2}\,(x_7 + y_7)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_7 + y_7)\,a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right)c\,\hat{\mathbf{z}}$	(12*i*)	O III
$\mathbf{B}_{41}$	=	$-y_7\,\mathbf{a}_1 + (x_7 - y_7)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}x_7 - y_7\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_7 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right)c\,\hat{\mathbf{z}}$	(12*i*)	O III
$\mathbf{B}_{42}$	=	$(-x_7 + y_7)\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\mathbf{a}_3$	=	$\left(-x_7 + \frac{1}{2}y_7\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_7 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right)c\,\hat{\mathbf{z}}$	(12*i*)	O III

References:

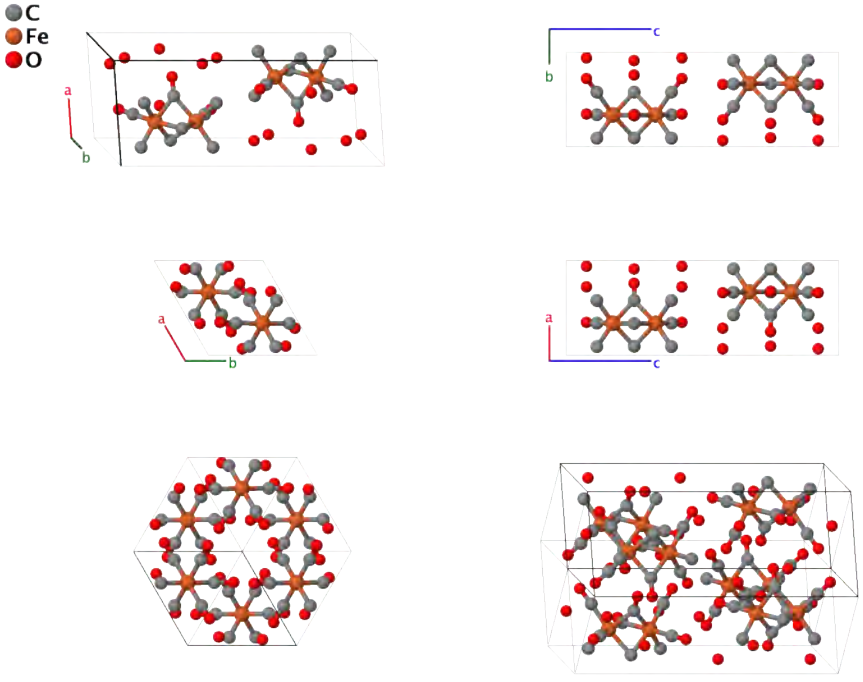
- J. M. Hughes and J. Rakovan, *The Crystal Structure of Apatite, $Ca_5(PO_4)_3(F,OH,Cl)$*, Rev. Mineral. Geochem. **48**, 1–12 (2002), doi:10.2138/rmg.2002.48.1.

Geometry files:

- CIF: pp. 1752
- POSCAR: pp. 1752

$Fe_2(CO)_9$ ($F4_1$) Structure: A9B2C9_hP40_176_hi_f_hi

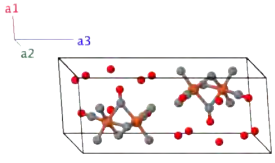
http://aflow.org/prototype-encyclopedia/A9B2C9_hP40_176_hi_f_hi



Prototype	:	$C_9Fe_2O_9$
AFLOW prototype label	:	A9B2C9_hP40_176_hi_f_hi
***Strukturbericht* designation**	:	$F4_1$
Pearson symbol	:	hP40
Space group number	:	176
Space group symbol	:	$P6_3/m$
AFLOW prototype command	:	`aflow --proto=A9B2C9_hP40_176_hi_f_hi --params=`$a, c/a, z_1, x_2, y_2, x_3, y_3, x_4, y_4, z_4, x_5, y_5, z_5$

Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{3}\mathbf{a}_1 + \frac{2}{3}\mathbf{a}_2 + z_1\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	$(4f)$	Fe
$\mathbf{B}_2$	=	$\frac{2}{3}\mathbf{a}_1 + \frac{1}{3}\mathbf{a}_2 + \left(\frac{1}{2}+z_1\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_1\right)c\,\hat{\mathbf{z}}$	$(4f)$	Fe
$\mathbf{B}_3$	=	$\frac{2}{3}\mathbf{a}_1 + \frac{1}{3}\mathbf{a}_2 - z_1\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_1 c\,\hat{\mathbf{z}}$	$(4f)$	Fe

$\mathbf{B}_4$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_1\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_1\right)c\,\hat{\mathbf{z}}$	$(4f)$	Fe
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_2 + y_2)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_2 + y_2)\,a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	C I
$\mathbf{B}_6$	=	$-y_2\,\mathbf{a}_1 + (x_2 - y_2)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_2 - y_2\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_2a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	C I
$\mathbf{B}_7$	=	$(-x_2 + y_2)\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(-x_2 + \frac{1}{2}y_2\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_2a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	C I
$\mathbf{B}_8$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-\frac{1}{2}\,(x_2 + y_2)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(x_2 - y_2)\,a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	C I
$\mathbf{B}_9$	=	$y_2\,\mathbf{a}_1 + (-x_2 + y_2)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_2 + y_2\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_2a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	C I
$\mathbf{B}_{10}$	=	$(x_2 - y_2)\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(x_2 - \frac{1}{2}y_2\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_2a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	C I
$\mathbf{B}_{11}$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_3 + y_3)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_3 + y_3)\,a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O I
$\mathbf{B}_{12}$	=	$-y_3\,\mathbf{a}_1 + (x_3 - y_3)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_3 - y_3\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O I
$\mathbf{B}_{13}$	=	$(-x_3 + y_3)\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(-x_3 + \frac{1}{2}y_3\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_3a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O I
$\mathbf{B}_{14}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-\frac{1}{2}\,(x_3 + y_3)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(x_3 - y_3)\,a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O I
$\mathbf{B}_{15}$	=	$y_3\,\mathbf{a}_1 + (-x_3 + y_3)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_3 + y_3\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O I
$\mathbf{B}_{16}$	=	$(x_3 - y_3)\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(x_3 - \frac{1}{2}y_3\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_3a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O I
$\mathbf{B}_{17}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_4 + y_4)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_4 + y_4)\,a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	$(12i)$	C II
$\mathbf{B}_{18}$	=	$-y_4\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_4 - y_4\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	$(12i)$	C II
$\mathbf{B}_{19}$	=	$(-x_4 + y_4)\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\left(-x_4 + \frac{1}{2}y_4\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_4a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	$(12i)$	C II
$\mathbf{B}_{20}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\,\mathbf{a}_3$	=	$-\frac{1}{2}\,(x_4 + y_4)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(x_4 - y_4)\,a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	$(12i)$	C II
$\mathbf{B}_{21}$	=	$y_4\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\,\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_4 + y_4\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	$(12i)$	C II
$\mathbf{B}_{22}$	=	$(x_4 - y_4)\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\,\mathbf{a}_3$	=	$\left(x_4 - \frac{1}{2}y_4\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_4a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	$(12i)$	C II
$\mathbf{B}_{23}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-\frac{1}{2}\,(x_4 + y_4)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(x_4 - y_4)\,a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	$(12i)$	C II
$\mathbf{B}_{24}$	=	$y_4\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_4 + y_4\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	$(12i)$	C II
$\mathbf{B}_{25}$	=	$(x_4 - y_4)\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\left(x_4 - \frac{1}{2}y_4\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_4a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	$(12i)$	C II
$\mathbf{B}_{26}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_4 + y_4)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_4 + y_4)\,a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	$(12i)$	C II
$\mathbf{B}_{27}$	=	$-y_4\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_4 - y_4\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	$(12i)$	C II
$\mathbf{B}_{28}$	=	$(-x_4 + y_4)\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\,\mathbf{a}_3$	=	$\left(-x_4 + \frac{1}{2}y_4\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_4a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	$(12i)$	C II
$\mathbf{B}_{29}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_5 + y_5)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_5 + y_5)\,a\,\hat{\mathbf{y}} + z_5c\,\hat{\mathbf{z}}$	$(12i)$	O II
$\mathbf{B}_{30}$	=	$-y_5\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_5 - y_5\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_5a\,\hat{\mathbf{y}} + z_5c\,\hat{\mathbf{z}}$	$(12i)$	O II

$$\mathbf{B}_{31} = (-x_5 + y_5)\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3 = \left(-x_5 + \tfrac{1}{2}y_5\right)a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}y_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}} \quad (12i) \quad \text{O II}$$

$$\mathbf{B}_{32} = -x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_3 = -\tfrac{1}{2}(x_5 + y_5)\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}(x_5 - y_5)\,a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_5\right)c\,\hat{\mathbf{z}} \quad (12i) \quad \text{O II}$$

$$\mathbf{B}_{33} = y_5\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_3 = \left(-\tfrac{1}{2}x_5 + y_5\right)a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_5\right)c\,\hat{\mathbf{z}} \quad (12i) \quad \text{O II}$$

$$\mathbf{B}_{34} = (x_5 - y_5)\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_3 = \left(x_5 - \tfrac{1}{2}y_5\right)a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}y_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_5\right)c\,\hat{\mathbf{z}} \quad (12i) \quad \text{O II}$$

$$\mathbf{B}_{35} = -x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3 = -\tfrac{1}{2}(x_5 + y_5)\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}(x_5 - y_5)\,a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}} \quad (12i) \quad \text{O II}$$

$$\mathbf{B}_{36} = y_5\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 - z_5\,\mathbf{a}_3 = \left(-\tfrac{1}{2}x_5 + y_5\right)a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}} \quad (12i) \quad \text{O II}$$

$$\mathbf{B}_{37} = (x_5 - y_5)\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3 = \left(x_5 - \tfrac{1}{2}y_5\right)a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}y_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}} \quad (12i) \quad \text{O II}$$

$$\mathbf{B}_{38} = x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_5\right)\mathbf{a}_3 = \tfrac{1}{2}(x_5 + y_5)\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}(-x_5 + y_5)\,a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_5\right)c\,\hat{\mathbf{z}} \quad (12i) \quad \text{O II}$$

$$\mathbf{B}_{39} = -y_5\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}x_5 - y_5\right)a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_5\right)c\,\hat{\mathbf{z}} \quad (12i) \quad \text{O II}$$

$$\mathbf{B}_{40} = (-x_5 + y_5)\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_5\right)\mathbf{a}_3 = \left(-x_5 + \tfrac{1}{2}y_5\right)a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}y_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_5\right)c\,\hat{\mathbf{z}} \quad (12i) \quad \text{O II}$$

References:

- F. A. Cotton and J. M. Troup, *Accurate determination of a classic structure in the metal carbonyl field: nonacarbonyldi-iron*, J. Chem. Soc., Dalton Trans. pp. 800–802 (1974), doi:10.1039/DT9740000800.

Found in:

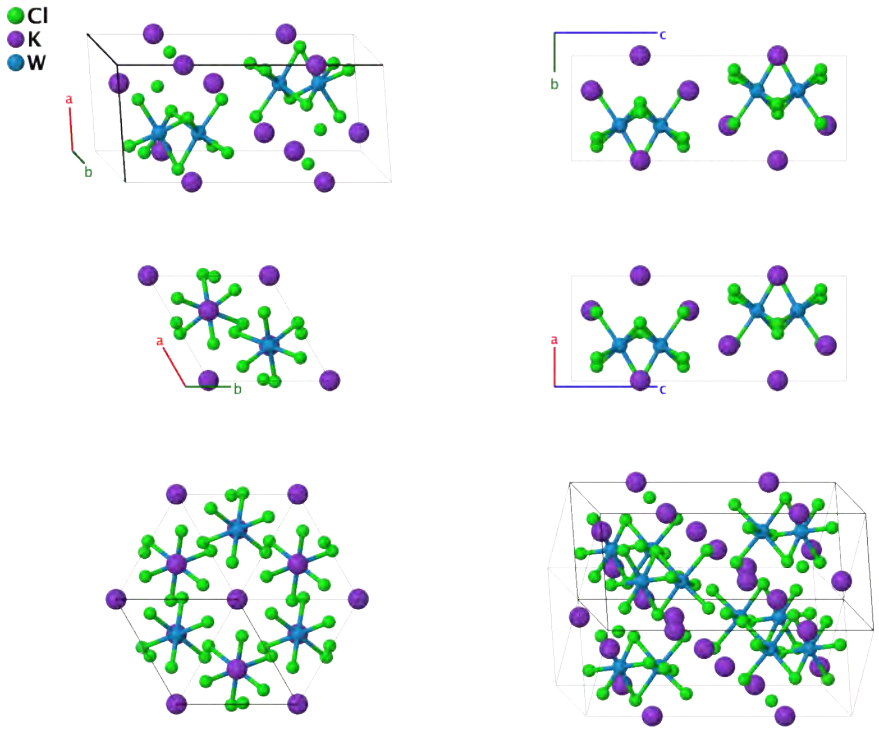
- M. Safa, Z. Dong, Y. Song, and Y. Huang, *Examining the structural changes in $Fe_2(CO)_9$ under high external pressures by Raman spectroscopy*, Can. J. Chem. **85**, 866–872 (2007), doi:10.1139/v07-096.

Geometry files:

- CIF: pp. 1752
- POSCAR: pp. 1753

$K_3W_2Cl_9$ ($K7_1$) Structure: A9B3C2_hP28_176_hi_af_f

http://aflow.org/prototype-encyclopedia/A9B3C2_hP28_176_hi_af_f



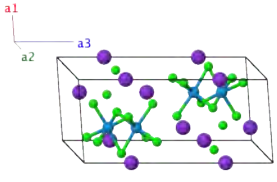
Prototype	:	$Cl_9K_3W_2$
AFLOW prototype label	:	A9B3C2_hP28_176_hi_af_f
***Strukturbericht* designation**	:	$K7_1$
Pearson symbol	:	hP28
Space group number	:	176
Space group symbol	:	$P6_3/m$
AFLOW prototype command	:	`aflow --proto=A9B3C2_hP28_176_hi_af_f --params=`$a, c/a, z_2, z_3, x_4, y_4, x_5, y_5, z_5$

Other compounds with this structure

- $K_3Ti_2Cl_9$, $Rb_3Ti_2Cl_9$, $K_3Mo_2Br_9$, $K_3Ti_2Br_9$, and $Rb_3Os_2Br_9$

Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$\mathbf{B}_1$	=	$\frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\,\hat{\mathbf{z}}$	$(2a)$	K I
$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\,\hat{\mathbf{z}}$	$(2a)$	K I
$\mathbf{B}_3$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(4f)$	K II
$\mathbf{B}_4$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	$(4f)$	K II
$\mathbf{B}_5$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	$(4f)$	K II
$\mathbf{B}_6$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	$(4f)$	K II
$\mathbf{B}_7$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(4f)$	W
$\mathbf{B}_8$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	$(4f)$	W
$\mathbf{B}_9$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(4f)$	W
$\mathbf{B}_{10}$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	$(4f)$	W
$\mathbf{B}_{11}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}(x_4 + y_4)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(-x_4 + y_4)\,a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	Cl I
$\mathbf{B}_{12}$	=	$-y_4\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_4 - y_4\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	Cl I
$\mathbf{B}_{13}$	=	$(-x_4 + y_4)\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(-x_4 + \frac{1}{2}y_4\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_4 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	Cl I
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-\frac{1}{2}(x_4 + y_4)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(x_4 - y_4)\,a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	Cl I
$\mathbf{B}_{15}$	=	$y_4\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_4 + y_4\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	Cl I
$\mathbf{B}_{16}$	=	$(x_4 - y_4)\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(x_4 - \frac{1}{2}y_4\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_4 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	Cl I
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\frac{1}{2}(x_5 + y_5)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(-x_5 + y_5)\,a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(12i)$	Cl II
$\mathbf{B}_{18}$	=	$-y_5\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_5 - y_5\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(12i)$	Cl II
$\mathbf{B}_{19}$	=	$(-x_5 + y_5)\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\left(-x_5 + \frac{1}{2}y_5\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(12i)$	Cl II
$\mathbf{B}_{20}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$-\frac{1}{2}(x_5 + y_5)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(x_5 - y_5)\,a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	$(12i)$	Cl II
$\mathbf{B}_{21}$	=	$y_5\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_5 + y_5\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	$(12i)$	Cl II
$\mathbf{B}_{22}$	=	$(x_5 - y_5)\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\left(x_5 - \frac{1}{2}y_5\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_5 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	$(12i)$	Cl II
$\mathbf{B}_{23}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-\frac{1}{2}(x_5 + y_5)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(x_5 - y_5)\,a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	$(12i)$	Cl II
$\mathbf{B}_{24}$	=	$y_5\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_5 + y_5\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	$(12i)$	Cl II
$\mathbf{B}_{25}$	=	$(x_5 - y_5)\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\left(x_5 - \frac{1}{2}y_5\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	$(12i)$	Cl II
$\mathbf{B}_{26}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\frac{1}{2}(x_5 + y_5)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(-x_5 + y_5)\,a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	$(12i)$	Cl II
$\mathbf{B}_{27}$	=	$-y_5\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}x_5 - y_5\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	$(12i)$	Cl II
$\mathbf{B}_{28}$	=	$(-x_5 + y_5)\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(-x_5 + \frac{1}{2}y_5\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_5 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	$(12i)$	Cl II

References:

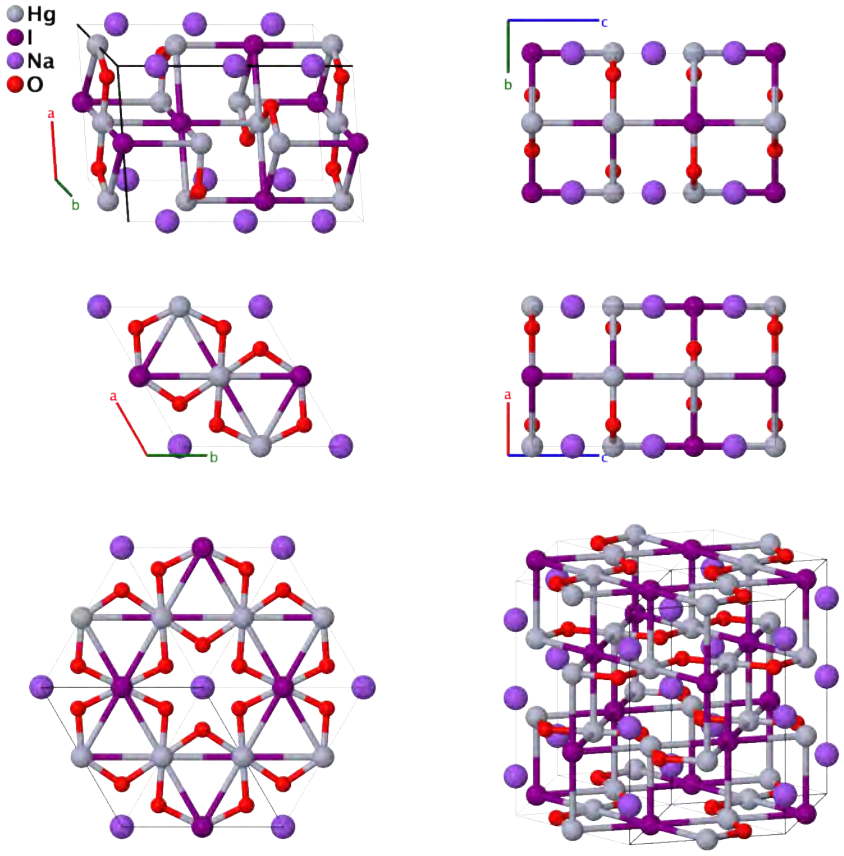
- W. H. Watson, Jr. and J. Waser, *Refinement of the structure of tripotassiumditungsten enneachloride, $K_3W_2Cl_9$*, Acta Cryst. **11**, 689–692 (1958), doi:10.1107/S0365110X58001869.

Geometry files:

- CIF: pp. 1753
- POSCAR: pp. 1753

Hg_2O_2NaI Structure: A2BCD2_hP18_180_f_c_b_i

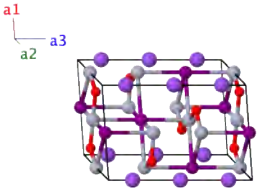
http://aflow.org/prototype-encyclopedia/A2BCD2_hP18_180_f_c_b_i



Prototype	:	Hg_2INaO_2
AFLOW prototype label	:	A2BCD2_hP18_180_f_c_b_i
***Strukturbericht* designation**	:	None
Pearson symbol	:	hP18
Space group number	:	180
Space group symbol	:	$P6_222$
AFLOW prototype command	:	`aflow --proto=A2BCD2_hP18_180_f_c_b_i --params=`$a, c/a, z_3, x_4$

Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$\mathbf{B}_1$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	($3b$)	Na
$\mathbf{B}_2$	=	$\frac{1}{6}\,\mathbf{a}_3$	=	$\frac{1}{6}c\,\hat{\mathbf{z}}$	($3b$)	Na
$\mathbf{B}_3$	=	$\frac{5}{6}\,\mathbf{a}_3$	=	$\frac{5}{6}c\,\hat{\mathbf{z}}$	($3b$)	Na
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{4}a\,\hat{\mathbf{y}}$	($3c$)	I
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{2}{3}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{4}a\,\hat{\mathbf{y}} + \frac{2}{3}c\,\hat{\mathbf{z}}$	($3c$)	I
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{3}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{3}c\,\hat{\mathbf{z}}$	($3c$)	I
$\mathbf{B}_7$	=	$\frac{1}{2}\,\mathbf{a}_1 + z_3\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{4}a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	($6f$)	Hg
$\mathbf{B}_8$	=	$\frac{1}{2}\,\mathbf{a}_2 + \left(\frac{2}{3} + z_3\right)\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{4}a\,\hat{\mathbf{y}} + \left(\frac{2}{3} + z_3\right)c\,\hat{\mathbf{z}}$	($6f$)	Hg
$\mathbf{B}_9$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \left(\frac{1}{3} + z_3\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \left(\frac{1}{3} + z_3\right)c\,\hat{\mathbf{z}}$	($6f$)	Hg
$\mathbf{B}_{10}$	=	$\frac{1}{2}\,\mathbf{a}_2 + \left(\frac{2}{3} - z_3\right)\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{4}a\,\hat{\mathbf{y}} + \left(\frac{2}{3} - z_3\right)c\,\hat{\mathbf{z}}$	($6f$)	Hg
$\mathbf{B}_{11}$	=	$\frac{1}{2}\,\mathbf{a}_1 - z_3\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{4}a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	($6f$)	Hg
$\mathbf{B}_{12}$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \left(\frac{1}{3} - z_3\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \left(\frac{1}{3} - z_3\right)c\,\hat{\mathbf{z}}$	($6f$)	Hg
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + 2x_4\,\mathbf{a}_2$	=	$\frac{3}{2}x_4a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4a\,\hat{\mathbf{y}}$	($6i$)	O
$\mathbf{B}_{14}$	=	$-2x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \frac{2}{3}\,\mathbf{a}_3$	=	$-\frac{3}{2}x_4a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4a\,\hat{\mathbf{y}} + \frac{2}{3}c\,\hat{\mathbf{z}}$	($6i$)	O
$\mathbf{B}_{15}$	=	$x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \frac{1}{3}\,\mathbf{a}_3$	=	$-\sqrt{3}x_4a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$	($6i$)	O
$\mathbf{B}_{16}$	=	$-x_4\,\mathbf{a}_1 - 2x_4\,\mathbf{a}_2$	=	$-\frac{3}{2}x_4a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4a\,\hat{\mathbf{y}}$	($6i$)	O
$\mathbf{B}_{17}$	=	$2x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \frac{2}{3}\,\mathbf{a}_3$	=	$\frac{3}{2}x_4a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4a\,\hat{\mathbf{y}} + \frac{2}{3}c\,\hat{\mathbf{z}}$	($6i$)	O
$\mathbf{B}_{18}$	=	$-x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \frac{1}{3}\,\mathbf{a}_3$	=	$\sqrt{3}x_4a\,\hat{\mathbf{y}} + \frac{1}{3}c\,\hat{\mathbf{z}}$	($6i$)	O

References:

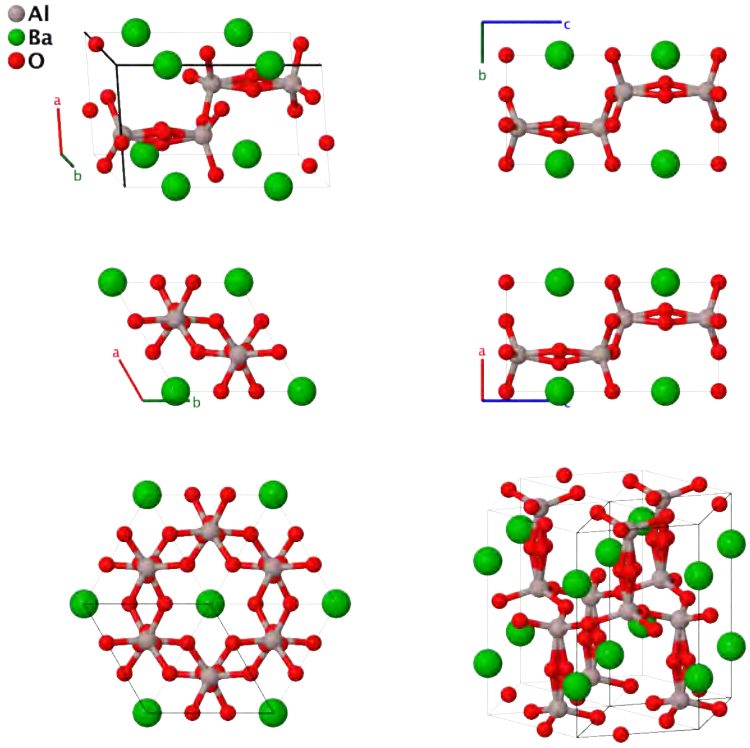
- K. Aurivillius, *Least-Squares Refinement of the Crystal Structures of Orthorhombic HgO and of Hg_2O_2NaI*, Acta Chem. Scand. **18**, 1305–1306 (1964), doi:10.3891/acta.chem.scand.18-1305.

Geometry files:

- CIF: pp. 1753
- POSCAR: pp. 1754

$BaAl_2O_4$ ($H2_8$) Structure: A2BC6_hP18_182_f_b_gh

http://aflow.org/prototype-encyclopedia/A2BC6_hP18_182_f_b_gh

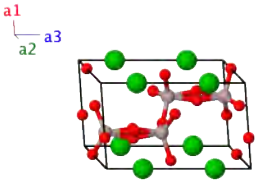


Prototype	:	Al_2BaO_4
AFLOW prototype label	:	A2BC6_hP18_182_f_b_gh
***Strukturbericht* designation**	:	$H2_8$
Pearson symbol	:	hP18
Space group number	:	182
Space group symbol	:	$P6_322$
AFLOW prototype command	:	`aflow --proto=A2BC6_hP18_182_f_b_gh --params=`$a, c/a, z_2, x_3, x_4$

- The structure of $BaAl_2O_4$ was originally determined by (Wallmark, 1937) and designated $H2_8$ by (Gottfried, 1940). This had two oxygen atoms at the ($2c$) Wyckoff position, $\pm(1/3, 2/3, z)$. (Perrotta, 1968) redetermined the structure and found that the best fit to external data required the ($2c$) oxygen atoms to be distributed among the ($6h$) sites. Thus each of the ($6h$) sites in this structure is 1/3 occupied.

Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\,\hat{\mathbf{z}}$	$(2b)$	Ba
$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\,\hat{\mathbf{z}}$	$(2b)$	Ba
$\mathbf{B}_3$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(4f)$	Al
$\mathbf{B}_4$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	$(4f)$	Al
$\mathbf{B}_5$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	$(4f)$	Al
$\mathbf{B}_6$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	$(4f)$	Al
$\mathbf{B}_7$	=	$x_3\,\mathbf{a}_1$	=	$\frac{1}{2}x_3 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}}$	$(6g)$	O I
$\mathbf{B}_8$	=	$x_3\,\mathbf{a}_2$	=	$\frac{1}{2}x_3 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}}$	$(6g)$	O I
$\mathbf{B}_9$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2$	=	$-x_3 a\,\hat{\mathbf{x}}$	$(6g)$	O I
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$-\frac{1}{2}x_3 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(6g)$	O I
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-\frac{1}{2}x_3 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(6g)$	O I
$\mathbf{B}_{12}$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(6g)$	O I
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + 2x_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{3}{2}x_4 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O II
$\mathbf{B}_{14}$	=	$-2x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-\frac{3}{2}x_4 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O II
$\mathbf{B}_{15}$	=	$x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-\sqrt{3}x_4 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O II
$\mathbf{B}_{16}$	=	$-x_4\,\mathbf{a}_1 - 2x_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-\frac{3}{2}x_4 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O II
$\mathbf{B}_{17}$	=	$2x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{2}x_4 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O II
$\mathbf{B}_{18}$	=	$-x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\sqrt{3}x_4 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O II

References:

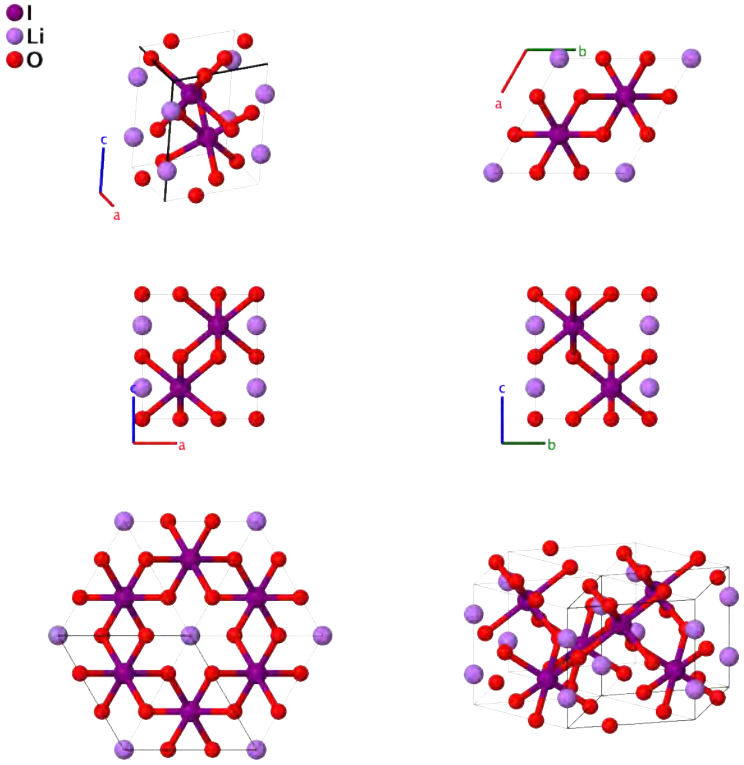
- A. J. Perrotta and J. V. Smith, *The Crystal Structure of $BaAl_2O_4$*, Bull. Soc. Fr. Mineral. Cristallogr. **91**, 85–87 (1968), doi:10.3406/bulmi.1968.6190.
- S. Wallmark and A. Westgren, *X-Ray Analysis of Barium Aluminates*, Ark. Kem. Mineral. Geol. B **12** (1937).
- C. Gottfried, ed., *Strukturbericht Band V 1937* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1940).

Geometry files:

- CIF: pp. 1754
- POSCAR: pp. 1754

$E2_3$ ($LiIO_3$) (*obsolete*) Structure: ABC3_hP10_182_c_b_g

http://aflow.org/prototype-encyclopedia/ABC3_hP10_182_c_b_g

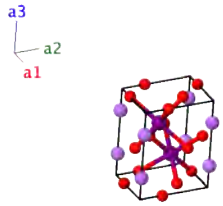


Prototype	:	$ILiO_3$
AFLOW prototype label	:	ABC3_hP10_182_c_b_g
***Strukturbericht* designation**	:	$E2_3$
Pearson symbol	:	hP10
Space group number	:	182
Space group symbol	:	$P6_322$
AFLOW prototype command	:	`aflow --proto=ABC3_hP10_182_c_b_g --params=`$a, c/a, x_3$

- $LiIO_3$ is known to exist in three forms:
  - α-$LiIO_3$, stable below 273 K: (Zachariasen, 1931) originally determined that this was the structure of α-$LiIO_3$, and (Hermann, 1937) designated it as *Strukturbericht* $E2_3$. (Rosenzweig, 1966) subsequently determined that this structure was incorrect because of the small sample size, and determined that the true structure was in space group $P6_3$ #173.
  - β-$LiIO_3$, stable from 573 K up to the melting point at 708 K.
  - γ-$LiIO_3$, stable between the α- and β-phases, with an orthorhombic structure.

Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\,\hat{\mathbf{z}}$	(2*b*)	Li
$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\,\hat{\mathbf{z}}$	(2*b*)	Li
$\mathbf{B}_3$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(2*c*)	I
$\mathbf{B}_4$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(2*c*)	I
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1$	=	$\frac{1}{2}x_3a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}}$	(6*g*)	O
$\mathbf{B}_6$	=	$x_3\,\mathbf{a}_2$	=	$\frac{1}{2}x_3a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}}$	(6*g*)	O
$\mathbf{B}_7$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2$	=	$-x_3a\,\hat{\mathbf{x}}$	(6*g*)	O
$\mathbf{B}_8$	=	$-x_3\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$-\frac{1}{2}x_3a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(6*g*)	O
$\mathbf{B}_9$	=	$-x_3\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-\frac{1}{2}x_3a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(6*g*)	O
$\mathbf{B}_{10}$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(6*g*)	O

References:

- W. H. Zachariasen and F. A. Barta, *Crystal Structure of Lithium Iodate*, Phys. Rev. **37**, 1626–1630 (1931), doi:10.1103/PhysRev.37.1626.
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Found in:

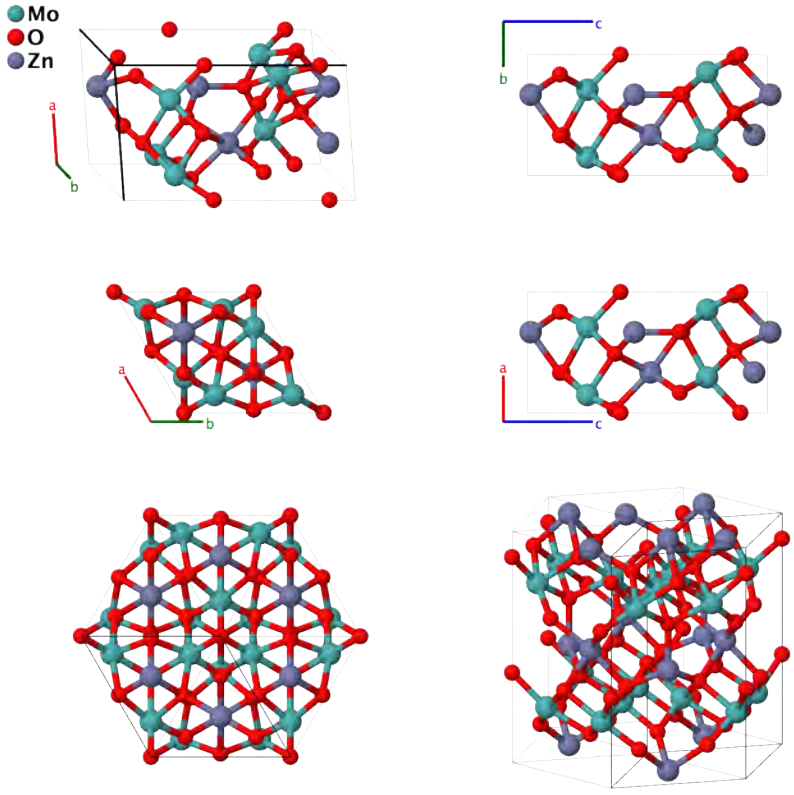
- A. Rosenzweig and B. Morosin, *A reinvestigation of the crystal structure of $LiIO_3$*, Acta Cryst. **20**, 758–761 (1966), doi:10.1107/S0365110X66001804.

Geometry files:

- CIF: pp. 1754
- POSCAR: pp. 1755

$Zn_2Mo_3O_8$ Structure: A3B8C2_hP26_186_c_ab2c_2b

http://aflow.org/prototype-encyclopedia/A3B8C2_hP26_186_c_ab2c_2b



Prototype	:	$Mo_3O_8Zn_2$
AFLOW prototype label	:	A3B8C2_hP26_186_c_ab2c_2b
***Strukturbericht* designation**	:	None
Pearson symbol	:	hP26
Space group number	:	186
Space group symbol	:	$P6_3mc$
AFLOW prototype command	:	`aflow --proto=A3B8C2_hP26_186_c_ab2c_2b --params=`$a, c/a, z_1, z_2, z_3, z_4, x_5, z_5, x_6, z_6, x_7, z_7$

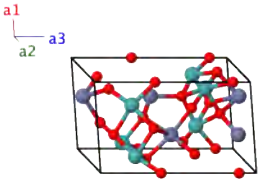
Other compounds with this structure

- $Cd_2Mo_3O_8$, $Co_2Mo_3O_8$, $Fe_2Mo_3O_8$ (kamiokite), $Mg_2Mo_3O_8$, $Mn_2Mo_3O_8$, and $Ni_2Mo_3O_8$

- Space group $P6_3mc$ #186 does not specify the origin of the z-coordinate. Here we fix it by setting the z-coordinate of the molybdenum atom at $z_5 = 1/4$.

Hexagonal primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$z_1\,\mathbf{a}_3$	$z_1 c\,\hat{\mathbf{z}}$	$(2a)$	O I
$\mathbf{B}_2$	$\left(\frac{1}{2}+z_1\right)\mathbf{a}_3$	$\left(\frac{1}{2}+z_1\right)c\,\hat{\mathbf{z}}$	$(2a)$	O I
$\mathbf{B}_3$	$\frac{1}{3}\,\mathbf{a}_1+\frac{2}{3}\,\mathbf{a}_2+z_2\,\mathbf{a}_3$	$\frac{1}{2}a\,\hat{\mathbf{x}}+\frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+z_2 c\,\hat{\mathbf{z}}$	$(2b)$	O II
$\mathbf{B}_4$	$\frac{2}{3}\,\mathbf{a}_1+\frac{1}{3}\,\mathbf{a}_2+\left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	$\frac{1}{2}a\,\hat{\mathbf{x}}-\frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_2\right)c\,\hat{\mathbf{z}}$	$(2b)$	O II
$\mathbf{B}_5$	$\frac{1}{3}\,\mathbf{a}_1+\frac{2}{3}\,\mathbf{a}_2+z_3\,\mathbf{a}_3$	$\frac{1}{2}a\,\hat{\mathbf{x}}+\frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+z_3 c\,\hat{\mathbf{z}}$	$(2b)$	Zn I
$\mathbf{B}_6$	$\frac{2}{3}\,\mathbf{a}_1+\frac{1}{3}\,\mathbf{a}_2+\left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	$\frac{1}{2}a\,\hat{\mathbf{x}}-\frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_3\right)c\,\hat{\mathbf{z}}$	$(2b)$	Zn I
$\mathbf{B}_7$	$\frac{1}{3}\,\mathbf{a}_1+\frac{2}{3}\,\mathbf{a}_2+z_4\,\mathbf{a}_3$	$\frac{1}{2}a\,\hat{\mathbf{x}}+\frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+z_4 c\,\hat{\mathbf{z}}$	$(2b)$	Zn II
$\mathbf{B}_8$	$\frac{2}{3}\,\mathbf{a}_1+\frac{1}{3}\,\mathbf{a}_2+\left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	$\frac{1}{2}a\,\hat{\mathbf{x}}-\frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_4\right)c\,\hat{\mathbf{z}}$	$(2b)$	Zn II
$\mathbf{B}_9$	$x_5\,\mathbf{a}_1-x_5\,\mathbf{a}_2+z_5\,\mathbf{a}_3$	$-\sqrt{3}x_5 a\,\hat{\mathbf{y}}+z_5 c\,\hat{\mathbf{z}}$	$(6c)$	Mo
$\mathbf{B}_{10}$	$x_5\,\mathbf{a}_1+2x_5\,\mathbf{a}_2+z_5\,\mathbf{a}_3$	$\frac{3}{2}x_5 a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}}+z_5 c\,\hat{\mathbf{z}}$	$(6c)$	Mo
$\mathbf{B}_{11}$	$-2x_5\,\mathbf{a}_1-x_5\,\mathbf{a}_2+z_5\,\mathbf{a}_3$	$-\frac{3}{2}x_5 a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}}+z_5 c\,\hat{\mathbf{z}}$	$(6c)$	Mo
$\mathbf{B}_{12}$	$-x_5\,\mathbf{a}_1+x_5\,\mathbf{a}_2+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	$\sqrt{3}x_5 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	$(6c)$	Mo
$\mathbf{B}_{13}$	$-x_5\,\mathbf{a}_1-2x_5\,\mathbf{a}_2+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	$-\frac{3}{2}x_5 a\,\hat{\mathbf{x}}-\frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	$(6c)$	Mo
$\mathbf{B}_{14}$	$2x_5\,\mathbf{a}_1+x_5\,\mathbf{a}_2+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	$\frac{3}{2}x_5 a\,\hat{\mathbf{x}}-\frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	$(6c)$	Mo
$\mathbf{B}_{15}$	$x_6\,\mathbf{a}_1-x_6\,\mathbf{a}_2+z_6\,\mathbf{a}_3$	$-\sqrt{3}x_6 a\,\hat{\mathbf{y}}+z_6 c\,\hat{\mathbf{z}}$	$(6c)$	O III
$\mathbf{B}_{16}$	$x_6\,\mathbf{a}_1+2x_6\,\mathbf{a}_2+z_6\,\mathbf{a}_3$	$\frac{3}{2}x_6 a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}}+z_6 c\,\hat{\mathbf{z}}$	$(6c)$	O III
$\mathbf{B}_{17}$	$-2x_6\,\mathbf{a}_1-x_6\,\mathbf{a}_2+z_6\,\mathbf{a}_3$	$-\frac{3}{2}x_6 a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}}+z_6 c\,\hat{\mathbf{z}}$	$(6c)$	O III
$\mathbf{B}_{18}$	$-x_6\,\mathbf{a}_1+x_6\,\mathbf{a}_2+\left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	$\sqrt{3}x_6 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	$(6c)$	O III
$\mathbf{B}_{19}$	$-x_6\,\mathbf{a}_1-2x_6\,\mathbf{a}_2+\left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	$-\frac{3}{2}x_6 a\,\hat{\mathbf{x}}-\frac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	$(6c)$	O III
$\mathbf{B}_{20}$	$2x_6\,\mathbf{a}_1+x_6\,\mathbf{a}_2+\left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	$\frac{3}{2}x_6 a\,\hat{\mathbf{x}}-\frac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	$(6c)$	O III
$\mathbf{B}_{21}$	$x_7\,\mathbf{a}_1-x_7\,\mathbf{a}_2+z_7\,\mathbf{a}_3$	$-\sqrt{3}x_7 a\,\hat{\mathbf{y}}+z_7 c\,\hat{\mathbf{z}}$	$(6c)$	O IV
$\mathbf{B}_{22}$	$x_7\,\mathbf{a}_1+2x_7\,\mathbf{a}_2+z_7\,\mathbf{a}_3$	$\frac{3}{2}x_7 a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}x_7 a\,\hat{\mathbf{y}}+z_7 c\,\hat{\mathbf{z}}$	$(6c)$	O IV
$\mathbf{B}_{23}$	$-2x_7\,\mathbf{a}_1-x_7\,\mathbf{a}_2+z_7\,\mathbf{a}_3$	$-\frac{3}{2}x_7 a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}x_7 a\,\hat{\mathbf{y}}+z_7 c\,\hat{\mathbf{z}}$	$(6c)$	O IV
$\mathbf{B}_{24}$	$-x_7\,\mathbf{a}_1+x_7\,\mathbf{a}_2+\left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	$\sqrt{3}x_7 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	$(6c)$	O IV
$\mathbf{B}_{25}$	$-x_7\,\mathbf{a}_1-2x_7\,\mathbf{a}_2+\left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	$-\frac{3}{2}x_7 a\,\hat{\mathbf{x}}-\frac{\sqrt{3}}{2}x_7 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	$(6c)$	O IV
$\mathbf{B}_{26}$	$2x_7\,\mathbf{a}_1+x_7\,\mathbf{a}_2+\left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	$\frac{3}{2}x_7 a\,\hat{\mathbf{x}}-\frac{\sqrt{3}}{2}x_7 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	$(6c)$	O IV

References:

- G. B. Ansell and L. Katz, *A Refinement of the Crystal Structure of Zinc Molybdenum(IV) Oxide,* $Zn_2Mo_3O_8$, Acta Cryst. **21**, 482–485 (1966), doi:10.1107/S0365110X66003359.

Found in:

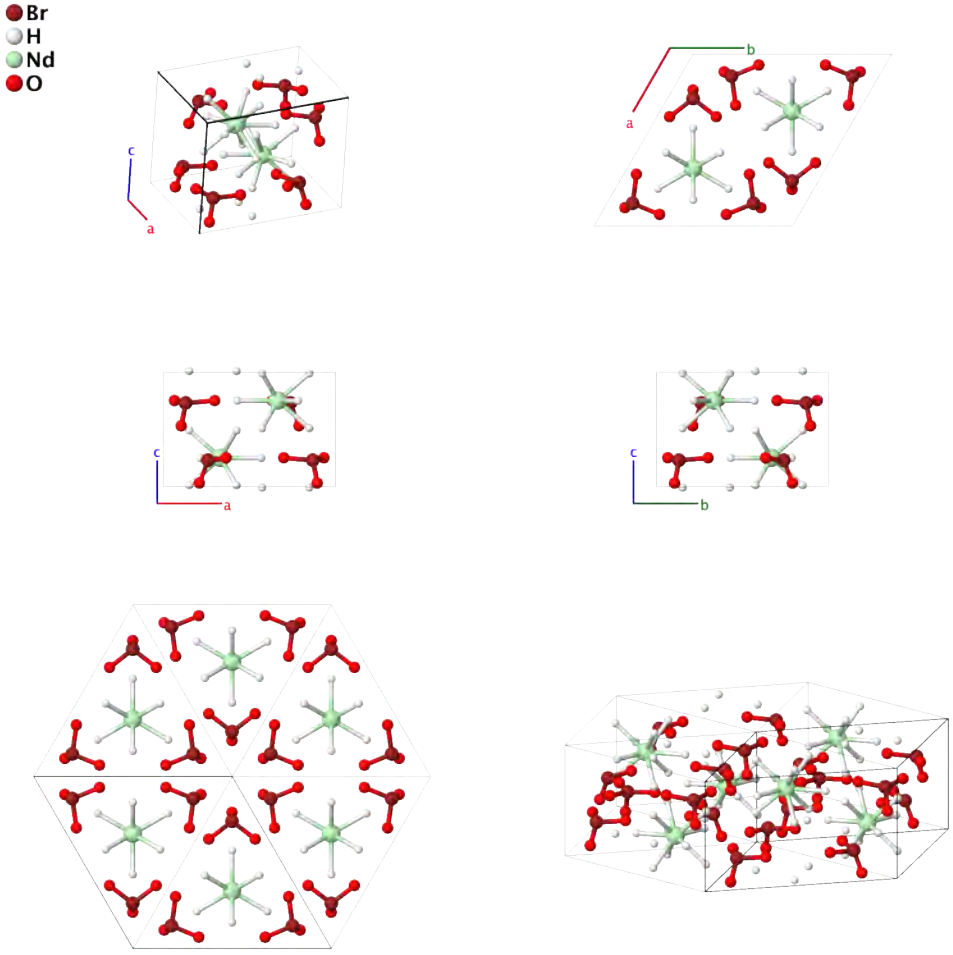
- Y. Kanazawa and A. Sasaki, *Structure of Kamiokite*, Acta Crystallogr. C **42**, 9–11 (1986), doi:10.1107/S0108270186097500.

Geometry files:

- CIF: pp. 1755
- POSCAR: pp. 1755

$Nd(BrO_3)_3{\cdot}9H_2O$ ($G2_2$) Structure: A3B9CD9_hP44_186_c_3c_b_cd

http://aflow.org/prototype-encyclopedia/A3B9CD9_hP44_186_c_3c_b_cd

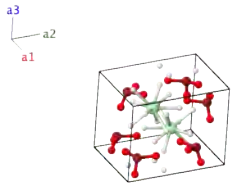


Prototype	:	$Br_3(H_2O)_9NdO_9$
AFLOW prototype label	:	A3B9CD9_hP44_186_c_3c_b_cd
***Strukturbericht* designation**	:	$G2_2$
Pearson symbol	:	hP44
Space group number	:	186
Space group symbol	:	$P6_3mc$
AFLOW prototype command	:	`aflow --proto=A3B9CD9_hP44_186_c_3c_b_cd --params=`$a, c/a, z_1, x_2, z_2, x_3, z_3, x_4, z_4, x_5, z_5, x_6, z_6, x_7, y_7, z_7$

- The positions of the hydrogen atoms in the water molecules were not determined, so we only provide the oxygen atom positions (labeled as H_2O).

Hexagonal primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_1\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_1 c\,\hat{\mathbf{z}}$	(2*b*)	Nd
$\mathbf{B}_2$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_1\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_1\right)c\,\hat{\mathbf{z}}$	(2*b*)	Nd
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$-\sqrt{3}x_2 a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(6*c*)	Br
$\mathbf{B}_4$	=	$x_2\,\mathbf{a}_1 + 2x_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{3}{2}x_2 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_2 a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(6*c*)	Br
$\mathbf{B}_5$	=	$-2x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$-\frac{3}{2}x_2 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_2 a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	(6*c*)	Br
$\mathbf{B}_6$	=	$-x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\sqrt{3}x_2 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	(6*c*)	Br
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 - 2x_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$-\frac{3}{2}x_2 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_2 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	(6*c*)	Br
$\mathbf{B}_8$	=	$2x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\frac{3}{2}x_2 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_2 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	(6*c*)	Br
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$-\sqrt{3}x_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(6*c*)	H_2O I
$\mathbf{B}_{10}$	=	$x_3\,\mathbf{a}_1 + 2x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{3}{2}x_3 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(6*c*)	H_2O I
$\mathbf{B}_{11}$	=	$-2x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$-\frac{3}{2}x_3 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	(6*c*)	H_2O I
$\mathbf{B}_{12}$	=	$-x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\sqrt{3}x_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(6*c*)	H_2O I
$\mathbf{B}_{13}$	=	$-x_3\,\mathbf{a}_1 - 2x_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$-\frac{3}{2}x_3 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(6*c*)	H_2O I
$\mathbf{B}_{14}$	=	$2x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\frac{3}{2}x_3 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	(6*c*)	H_2O I
$\mathbf{B}_{15}$	=	$x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$-\sqrt{3}x_4 a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(6*c*)	H_2O II
$\mathbf{B}_{16}$	=	$x_4\,\mathbf{a}_1 + 2x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\frac{3}{2}x_4 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(6*c*)	H_2O II
$\mathbf{B}_{17}$	=	$-2x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$-\frac{3}{2}x_4 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	(6*c*)	H_2O II
$\mathbf{B}_{18}$	=	$-x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\sqrt{3}x_4 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(6*c*)	H_2O II
$\mathbf{B}_{19}$	=	$-x_4\,\mathbf{a}_1 - 2x_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$-\frac{3}{2}x_4 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(6*c*)	H_2O II
$\mathbf{B}_{20}$	=	$2x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\frac{3}{2}x_4 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(6*c*)	H_2O II
$\mathbf{B}_{21}$	=	$x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$-\sqrt{3}x_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(6*c*)	H_2O III
$\mathbf{B}_{22}$	=	$x_5\,\mathbf{a}_1 + 2x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\frac{3}{2}x_5 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(6*c*)	H_2O III
$\mathbf{B}_{23}$	=	$-2x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$-\frac{3}{2}x_5 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(6*c*)	H_2O III
$\mathbf{B}_{24}$	=	$-x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\sqrt{3}x_5 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	(6*c*)	H_2O III
$\mathbf{B}_{25}$	=	$-x_5\,\mathbf{a}_1 - 2x_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$-\frac{3}{2}x_5 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	(6*c*)	H_2O III
$\mathbf{B}_{26}$	=	$2x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\frac{3}{2}x_5 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	(6*c*)	H_2O III
$\mathbf{B}_{27}$	=	$x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$-\sqrt{3}x_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(6*c*)	O I
$\mathbf{B}_{28}$	=	$x_6\,\mathbf{a}_1 + 2x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\frac{3}{2}x_6 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(6*c*)	O I

$\mathbf{B}_{29}$	=	$-2x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$-\frac{3}{2}x_6 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(6*c*)	O I
$\mathbf{B}_{30}$	=	$-x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + \left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	=	$\sqrt{3}x_6 a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	(6*c*)	O I
$\mathbf{B}_{31}$	=	$-x_6\,\mathbf{a}_1 - 2x_6\,\mathbf{a}_2 + \left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	=	$-\frac{3}{2}x_6 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	(6*c*)	O I
$\mathbf{B}_{32}$	=	$2x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + \left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	=	$\frac{3}{2}x_6 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	(6*c*)	O I
$\mathbf{B}_{33}$	=	$x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$\frac{1}{2}(x_7+y_7)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(-x_7+y_7)\,a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(12*d*)	O II
$\mathbf{B}_{34}$	=	$-y_7\,\mathbf{a}_1 + (x_7-y_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_7 - y_7\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_7 a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(12*d*)	O II
$\mathbf{B}_{35}$	=	$(-x_7+y_7)\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$\left(-x_7 + \frac{1}{2}y_7\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_7 a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(12*d*)	O II
$\mathbf{B}_{36}$	=	$-x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	=	$-\frac{1}{2}(x_7+y_7)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(x_7-y_7)\,a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	(12*d*)	O II
$\mathbf{B}_{37}$	=	$y_7\,\mathbf{a}_1 + (-x_7+y_7)\,\mathbf{a}_2 + \left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_7 + y_7\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_7 a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	(12*d*)	O II
$\mathbf{B}_{38}$	=	$(x_7-y_7)\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + \left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	=	$\left(x_7 - \frac{1}{2}y_7\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_7 a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	(12*d*)	O II
$\mathbf{B}_{39}$	=	$-y_7\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$-\frac{1}{2}(x_7+y_7)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(-x_7+y_7)\,a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(12*d*)	O II
$\mathbf{B}_{40}$	=	$(-x_7+y_7)\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_7 + y_7\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_7 a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(12*d*)	O II
$\mathbf{B}_{41}$	=	$x_7\,\mathbf{a}_1 + (x_7-y_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$\left(x_7 - \frac{1}{2}y_7\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_7 a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(12*d*)	O II
$\mathbf{B}_{42}$	=	$y_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + \left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	=	$\frac{1}{2}(x_7+y_7)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(x_7-y_7)\,a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	(12*d*)	O II
$\mathbf{B}_{43}$	=	$(x_7-y_7)\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}x_7 - y_7\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_7 a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	(12*d*)	O II
$\mathbf{B}_{44}$	=	$-x_7\,\mathbf{a}_1 + (-x_7+y_7)\,\mathbf{a}_2 + \left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	=	$\left(-x_7 + \frac{1}{2}y_7\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_7 a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	(12*d*)	O II

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- L. Helmholz, *The Crystal Structure of Neodymium Bromate Enneahydrate, $Nd(BrO_3)_3{\cdot}9H_2O$*, J. Am. Chem. Soc. **61**, 1544–1550 (1939), doi:10.1021/ja01875a062.

Found in:

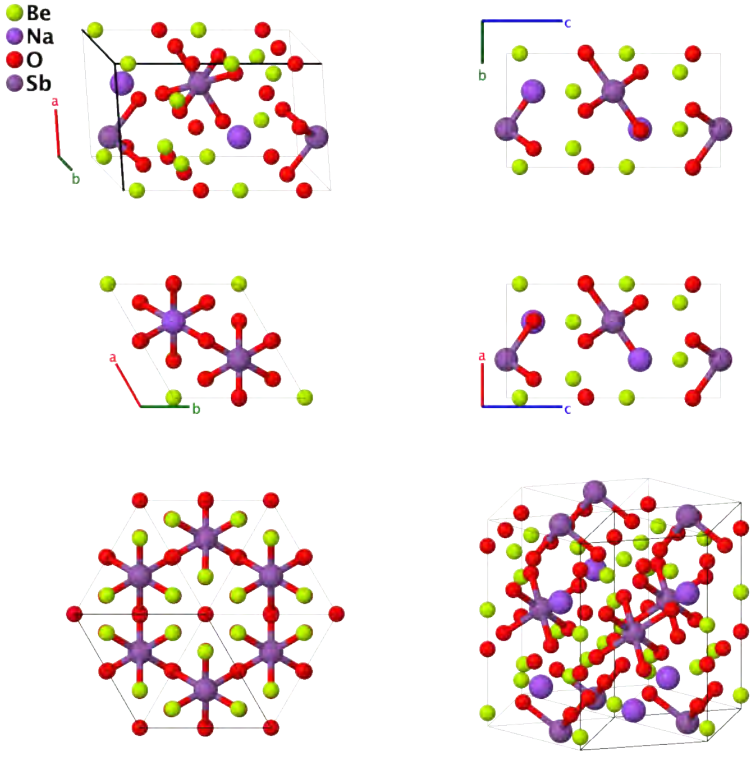
- K. Herrmann, ed., *Strukturbericht Band VII 1939* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1943).

Geometry files:

- CIF: pp. 1756
- POSCAR: pp. 1756

Swedenborgite ($NaBe_4SbO_7$, $E9_2$) Structure: A4BC7D_hP26_186_ac_b_a2c_b

http://aflow.org/prototype-encyclopedia/A4BC7D_hP26_186_ac_b_a2c_b



Prototype	:	Be_4NaO_7Sb
AFLOW prototype label	:	A4BC7D_hP26_186_ac_b_a2c_b
***Strukturbericht* designation**	:	$E9_2$
Pearson symbol	:	hP26
Space group number	:	186
Space group symbol	:	$P6_3mc$
AFLOW prototype command	:	`aflow --proto=A4BC7D_hP26_186_ac_b_a2c_b --params=`$a, c/a, z_1, z_2, z_3, z_4, x_5, z_5, x_6, z_6, x_7, z_7$

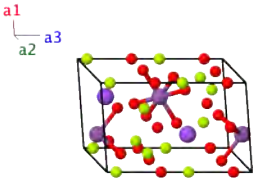
Other compounds with this structure

- $CaBa(Fe_2Mn_2)O_7$, $CaBaCo_3AlO_7$, $CaBaCo_3FeO_7$, $CaBaCo_3ZnO_7$, $CaBaCo_4O_7$, $CaBaFe_4O_7$, $Er_2Si_4N_6C$, $Ho_2Si_4N_6C$, $Tb_2Si_4N_6C$, $Y_2Si_4N_6C$, $YBaCo_3AlO_7$, $YBaCo_3FeO_7$, $YBaCo_4O_7$, $YBaMn_3AlO_7$, $Yb_2Si_4N_6C$, and $YbBaSi_4N_7$

- The actual composition of the studied sample is ($Na_{0.89}Ca_{0.04}Be_4SbO_7$), with the sodium (Na-I) site containing 89% sodium, 4% calcium, and 7% vacancies.
- Space group $P6_3mc$ #186 does not fix the origin of the z-coordinate. Here the position of the antimony atom (Sb-I) is set so that $z_4 = 0$.

Hexagonal primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$z_1\,\mathbf{a}_3$	=	$z_1 c\,\hat{\mathbf{z}}$	$(2a)$	Be I
$\mathbf{B}_2$	=	$\left(\frac{1}{2}+z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+z_1\right)c\,\hat{\mathbf{z}}$	$(2a)$	Be I
$\mathbf{B}_3$	=	$z_2\,\mathbf{a}_3$	=	$z_2 c\,\hat{\mathbf{z}}$	$(2a)$	O I
$\mathbf{B}_4$	=	$\left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+z_2\right)c\,\hat{\mathbf{z}}$	$(2a)$	O I
$\mathbf{B}_5$	=	$\frac{1}{3}\,\mathbf{a}_1+\frac{2}{3}\,\mathbf{a}_2+z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}+\frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+z_3 c\,\hat{\mathbf{z}}$	$(2b)$	Na
$\mathbf{B}_6$	=	$\frac{2}{3}\,\mathbf{a}_1+\frac{1}{3}\,\mathbf{a}_2+\left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}-\frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_3\right)c\,\hat{\mathbf{z}}$	$(2b)$	Na
$\mathbf{B}_7$	=	$\frac{1}{3}\,\mathbf{a}_1+\frac{2}{3}\,\mathbf{a}_2+z_4\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}+\frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+z_4 c\,\hat{\mathbf{z}}$	$(2b)$	Sb
$\mathbf{B}_8$	=	$\frac{2}{3}\,\mathbf{a}_1+\frac{1}{3}\,\mathbf{a}_2+\left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}-\frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_4\right)c\,\hat{\mathbf{z}}$	$(2b)$	Sb
$\mathbf{B}_9$	=	$x_5\,\mathbf{a}_1-x_5\,\mathbf{a}_2+z_5\,\mathbf{a}_3$	=	$-\sqrt{3}x_5 a\,\hat{\mathbf{y}}+z_5 c\,\hat{\mathbf{z}}$	$(6c)$	Be II
$\mathbf{B}_{10}$	=	$x_5\,\mathbf{a}_1+2x_5\,\mathbf{a}_2+z_5\,\mathbf{a}_3$	=	$\frac{3}{2}x_5 a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}}+z_5 c\,\hat{\mathbf{z}}$	$(6c)$	Be II
$\mathbf{B}_{11}$	=	$-2x_5\,\mathbf{a}_1-x_5\,\mathbf{a}_2+z_5\,\mathbf{a}_3$	=	$-\frac{3}{2}x_5 a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}}+z_5 c\,\hat{\mathbf{z}}$	$(6c)$	Be II
$\mathbf{B}_{12}$	=	$-x_5\,\mathbf{a}_1+x_5\,\mathbf{a}_2+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$\sqrt{3}x_5 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	$(6c)$	Be II
$\mathbf{B}_{13}$	=	$-x_5\,\mathbf{a}_1-2x_5\,\mathbf{a}_2+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$-\frac{3}{2}x_5 a\,\hat{\mathbf{x}}-\frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	$(6c)$	Be II
$\mathbf{B}_{14}$	=	$2x_5\,\mathbf{a}_1+x_5\,\mathbf{a}_2+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$\frac{3}{2}x_5 a\,\hat{\mathbf{x}}-\frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	$(6c)$	Be II
$\mathbf{B}_{15}$	=	$x_6\,\mathbf{a}_1-x_6\,\mathbf{a}_2+z_6\,\mathbf{a}_3$	=	$-\sqrt{3}x_6 a\,\hat{\mathbf{y}}+z_6 c\,\hat{\mathbf{z}}$	$(6c)$	O II
$\mathbf{B}_{16}$	=	$x_6\,\mathbf{a}_1+2x_6\,\mathbf{a}_2+z_6\,\mathbf{a}_3$	=	$\frac{3}{2}x_6 a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}}+z_6 c\,\hat{\mathbf{z}}$	$(6c)$	O II
$\mathbf{B}_{17}$	=	$-2x_6\,\mathbf{a}_1-x_6\,\mathbf{a}_2+z_6\,\mathbf{a}_3$	=	$-\frac{3}{2}x_6 a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}}+z_6 c\,\hat{\mathbf{z}}$	$(6c)$	O II
$\mathbf{B}_{18}$	=	$-x_6\,\mathbf{a}_1+x_6\,\mathbf{a}_2+\left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	=	$\sqrt{3}x_6 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	$(6c)$	O II
$\mathbf{B}_{19}$	=	$-x_6\,\mathbf{a}_1-2x_6\,\mathbf{a}_2+\left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	=	$-\frac{3}{2}x_6 a\,\hat{\mathbf{x}}-\frac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	$(6c)$	O II
$\mathbf{B}_{20}$	=	$2x_6\,\mathbf{a}_1+x_6\,\mathbf{a}_2+\left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	=	$\frac{3}{2}x_6 a\,\hat{\mathbf{x}}-\frac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	$(6c)$	O II
$\mathbf{B}_{21}$	=	$x_7\,\mathbf{a}_1-x_7\,\mathbf{a}_2+z_7\,\mathbf{a}_3$	=	$-\sqrt{3}x_7 a\,\hat{\mathbf{y}}+z_7 c\,\hat{\mathbf{z}}$	$(6c)$	O III
$\mathbf{B}_{22}$	=	$x_7\,\mathbf{a}_1+2x_7\,\mathbf{a}_2+z_7\,\mathbf{a}_3$	=	$\frac{3}{2}x_7 a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}x_7 a\,\hat{\mathbf{y}}+z_7 c\,\hat{\mathbf{z}}$	$(6c)$	O III
$\mathbf{B}_{23}$	=	$-2x_7\,\mathbf{a}_1-x_7\,\mathbf{a}_2+z_7\,\mathbf{a}_3$	=	$-\frac{3}{2}x_7 a\,\hat{\mathbf{x}}+\frac{\sqrt{3}}{2}x_7 a\,\hat{\mathbf{y}}+z_7 c\,\hat{\mathbf{z}}$	$(6c)$	O III
$\mathbf{B}_{24}$	=	$-x_7\,\mathbf{a}_1+x_7\,\mathbf{a}_2+\left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	=	$\sqrt{3}x_7 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	$(6c)$	O III
$\mathbf{B}_{25}$	=	$-x_7\,\mathbf{a}_1-2x_7\,\mathbf{a}_2+\left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	=	$-\frac{3}{2}x_7 a\,\hat{\mathbf{x}}-\frac{\sqrt{3}}{2}x_7 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	$(6c)$	O III
$\mathbf{B}_{26}$	=	$2x_7\,\mathbf{a}_1+x_7\,\mathbf{a}_2+\left(\frac{1}{2}+z_7\right)\mathbf{a}_3$	=	$\frac{3}{2}x_7 a\,\hat{\mathbf{x}}-\frac{\sqrt{3}}{2}x_7 a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_7\right)c\,\hat{\mathbf{z}}$	$(6c)$	O III

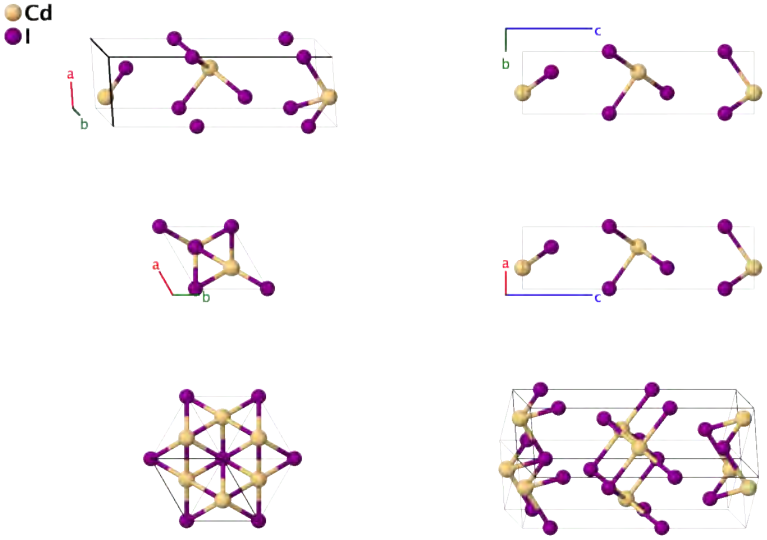
References:

- D. M. C. Huminicki and F. C. Hawthorne, *Refinement of the Crystal Structure of Swedenborgite*, Can. Mineral. **39**, 153–158 (2001), doi:10.2113/gscanmin.39.1.153.

Geometry files:
- CIF: pp. 1756
- POSCAR: pp. 1757

*C*27 (CdI_2) (*questionable*) Structure: AB2_hP6_186_b_ab

http://aflow.org/prototype-encyclopedia/AB2_hP6_186_b_ab

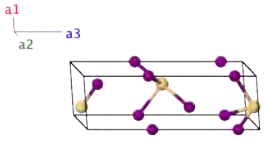


Prototype	:	CdI_2
AFLOW prototype label	:	AB2_hP6_186_b_ab
***Strukturbericht* designation**	:	*C*27
Pearson symbol	:	hP6
Space group number	:	186
Space group symbol	:	$P6_3mc$
AFLOW prototype command	:	`aflow --proto=AB2_hP6_186_b_ab --params=`$a, c/a, z_1, z_2, z_3$

- This is a modification of the *C*6 (ω phase) CdI_2 structure proposed by (Hassel, 1933) to explain extra lines found in the x-ray spectra. This structure was never accepted by most researchers and does not appear in modern lists of *Strukturbericht* symbols, although (Parthé, 1993) lists it as a stacking varient of CdI_2. We include it here as part of the historical record.
- Hassel originally placed a Cd atom at the origin, however both (Gottfried, 1937) and FINDSYM shift the origin as shown.

Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$\mathbf{B}_1$	=	$z_1\,\mathbf{a}_3$	=	$z_1 c\,\hat{\mathbf{z}}$	(2*a*)	I I
$\mathbf{B}_2$	=	$\left(\frac{1}{2}+z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+z_1\right)c\,\hat{\mathbf{z}}$	(2*a*)	I I
$\mathbf{B}_3$	=	$\frac{1}{3}\,\mathbf{a}_1+\frac{2}{3}\,\mathbf{a}_2+z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}+\frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+z_2 c\,\hat{\mathbf{z}}$	(2*b*)	Cd
$\mathbf{B}_4$	=	$\frac{2}{3}\,\mathbf{a}_1+\frac{1}{3}\,\mathbf{a}_2+\left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}-\frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_2\right)c\,\hat{\mathbf{z}}$	(2*b*)	Cd
$\mathbf{B}_5$	=	$\frac{1}{3}\,\mathbf{a}_1+\frac{2}{3}\,\mathbf{a}_2+z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}+\frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+z_3 c\,\hat{\mathbf{z}}$	(2*b*)	I II
$\mathbf{B}_6$	=	$\frac{2}{3}\,\mathbf{a}_1+\frac{1}{3}\,\mathbf{a}_2+\left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}-\frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_3\right)c\,\hat{\mathbf{z}}$	(2*b*)	I II

References:

- O. Hassel, *Zur Kristallstruktur des Cadmiumjodids CdJ_2*, Z. Physik. Chem. **22B**, 333–334 (1933), doi:10.1515/zpch-1933-2228.
- E. Parthé, L. Gelato, B. Chabot, M. Penso, K. Cenzual, and R. Gladyshevskii, in *Standardized Data and Crystal Chemical Characterization of Inorganic Structure Types* (Springer-Verlag, Berlin, Heidelberg, 1993), *Gmelin Handbook of Inorganic and Organometallic Chemistry*, vol. 2, chap. Crystal Chemical Characterization of Inorganic Structure Types, 8 edn., doi:10.1007/978-3-662-02909-1_3.

Found in:

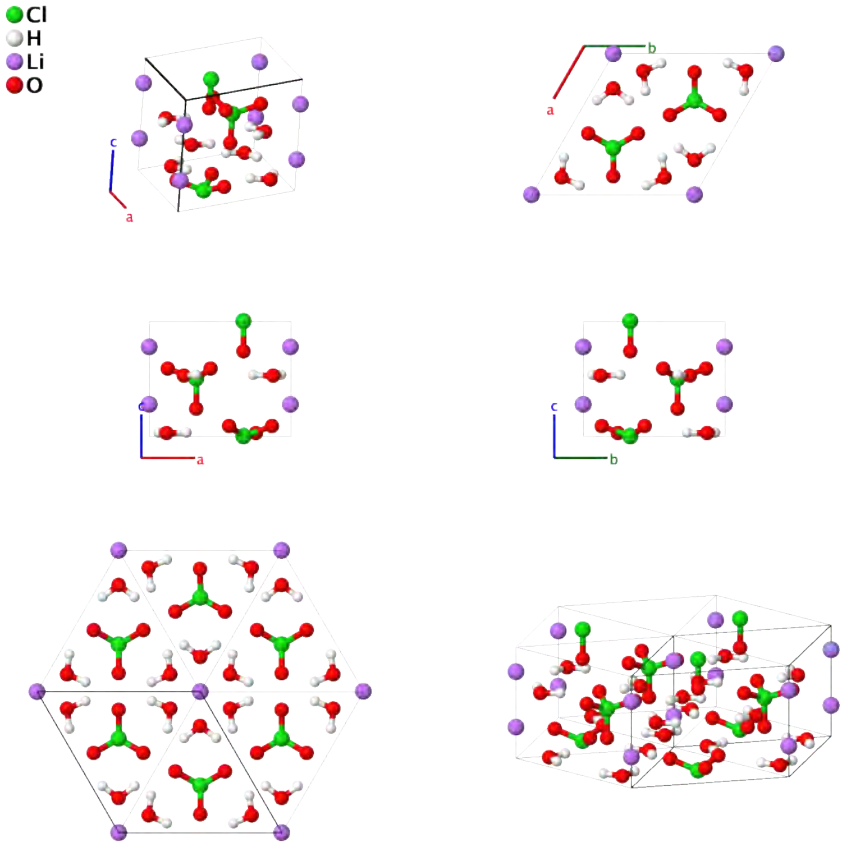
- C. Gottfried and F. Schossberger, eds., *Strukturbericht Band III 1933-1935* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1757
- POSCAR: pp. 1757

$LiClO_4 \cdot 3H_2O$ ($H4_{18}$) Structure: AB6CD7_hP30_186_b_d_a_b2c

http://aflow.org/prototype-encyclopedia/AB6CD7_hP30_186_b_d_a_b2c



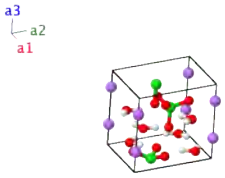
Prototype	:	ClH_6LiO_7
AFLOW prototype label	:	AB6CD7_hP30_186_b_d_a_b2c
***Strukturbericht* designation**	:	$H4_{18}$
Pearson symbol	:	hP30
Space group number	:	186
Space group symbol	:	$P6_3mc$
AFLOW prototype command	:	`aflow --proto=AB6CD7_hP30_186_b_d_a_b2c --params=`$a, c/a, z_1, z_2, z_3, x_4, z_4, x_5, z_5, x_6, y_6, z_6$

Other compounds with this structure

- $LiTcO_4 \cdot 3H_2O$

Hexagonal primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$z_1\,\mathbf{a}_3$	=	$z_1 c\,\hat{\mathbf{z}}$	$(2a)$	Li
$\mathbf{B}_2$	=	$\left(\tfrac{1}{2}+z_1\right)\mathbf{a}_3$	=	$\left(\tfrac{1}{2}+z_1\right)c\,\hat{\mathbf{z}}$	$(2a)$	Li
$\mathbf{B}_3$	=	$\tfrac{1}{3}\,\mathbf{a}_1+\tfrac{2}{3}\,\mathbf{a}_2+z_2\,\mathbf{a}_3$	=	$\tfrac{1}{2}a\,\hat{\mathbf{x}}+\tfrac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+z_2 c\,\hat{\mathbf{z}}$	$(2b)$	Cl
$\mathbf{B}_4$	=	$\tfrac{2}{3}\,\mathbf{a}_1+\tfrac{1}{3}\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_2\right)\mathbf{a}_3$	=	$\tfrac{1}{2}a\,\hat{\mathbf{x}}-\tfrac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_2\right)c\,\hat{\mathbf{z}}$	$(2b)$	Cl
$\mathbf{B}_5$	=	$\tfrac{1}{3}\,\mathbf{a}_1+\tfrac{2}{3}\,\mathbf{a}_2+z_3\,\mathbf{a}_3$	=	$\tfrac{1}{2}a\,\hat{\mathbf{x}}+\tfrac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+z_3 c\,\hat{\mathbf{z}}$	$(2b)$	O I
$\mathbf{B}_6$	=	$\tfrac{2}{3}\,\mathbf{a}_1+\tfrac{1}{3}\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_3\right)\mathbf{a}_3$	=	$\tfrac{1}{2}a\,\hat{\mathbf{x}}-\tfrac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_3\right)c\,\hat{\mathbf{z}}$	$(2b)$	O I
$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_1-x_4\,\mathbf{a}_2+z_4\,\mathbf{a}_3$	=	$-\sqrt{3}x_4 a\,\hat{\mathbf{y}}+z_4 c\,\hat{\mathbf{z}}$	$(6c)$	O II
$\mathbf{B}_8$	=	$x_4\,\mathbf{a}_1+2x_4\,\mathbf{a}_2+z_4\,\mathbf{a}_3$	=	$\tfrac{3}{2}x_4 a\,\hat{\mathbf{x}}+\tfrac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}}+z_4 c\,\hat{\mathbf{z}}$	$(6c)$	O II
$\mathbf{B}_9$	=	$-2x_4\,\mathbf{a}_1-x_4\,\mathbf{a}_2+z_4\,\mathbf{a}_3$	=	$-\tfrac{3}{2}x_4 a\,\hat{\mathbf{x}}+\tfrac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}}+z_4 c\,\hat{\mathbf{z}}$	$(6c)$	O II
$\mathbf{B}_{10}$	=	$-x_4\,\mathbf{a}_1+x_4\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_4\right)\mathbf{a}_3$	=	$\sqrt{3}x_4 a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_4\right)c\,\hat{\mathbf{z}}$	$(6c)$	O II
$\mathbf{B}_{11}$	=	$-x_4\,\mathbf{a}_1-2x_4\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_4\right)\mathbf{a}_3$	=	$-\tfrac{3}{2}x_4 a\,\hat{\mathbf{x}}-\tfrac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_4\right)c\,\hat{\mathbf{z}}$	$(6c)$	O II
$\mathbf{B}_{12}$	=	$2x_4\,\mathbf{a}_1+x_4\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_4\right)\mathbf{a}_3$	=	$\tfrac{3}{2}x_4 a\,\hat{\mathbf{x}}-\tfrac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_4\right)c\,\hat{\mathbf{z}}$	$(6c)$	O II
$\mathbf{B}_{13}$	=	$x_5\,\mathbf{a}_1-x_5\,\mathbf{a}_2+z_5\,\mathbf{a}_3$	=	$-\sqrt{3}x_5 a\,\hat{\mathbf{y}}+z_5 c\,\hat{\mathbf{z}}$	$(6c)$	O III
$\mathbf{B}_{14}$	=	$x_5\,\mathbf{a}_1+2x_5\,\mathbf{a}_2+z_5\,\mathbf{a}_3$	=	$\tfrac{3}{2}x_5 a\,\hat{\mathbf{x}}+\tfrac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}}+z_5 c\,\hat{\mathbf{z}}$	$(6c)$	O III
$\mathbf{B}_{15}$	=	$-2x_5\,\mathbf{a}_1-x_5\,\mathbf{a}_2+z_5\,\mathbf{a}_3$	=	$-\tfrac{3}{2}x_5 a\,\hat{\mathbf{x}}+\tfrac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}}+z_5 c\,\hat{\mathbf{z}}$	$(6c)$	O III
$\mathbf{B}_{16}$	=	$-x_5\,\mathbf{a}_1+x_5\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_5\right)\mathbf{a}_3$	=	$\sqrt{3}x_5 a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	$(6c)$	O III
$\mathbf{B}_{17}$	=	$-x_5\,\mathbf{a}_1-2x_5\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_5\right)\mathbf{a}_3$	=	$-\tfrac{3}{2}x_5 a\,\hat{\mathbf{x}}-\tfrac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	$(6c)$	O III
$\mathbf{B}_{18}$	=	$2x_5\,\mathbf{a}_1+x_5\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_5\right)\mathbf{a}_3$	=	$\tfrac{3}{2}x_5 a\,\hat{\mathbf{x}}-\tfrac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_5\right)c\,\hat{\mathbf{z}}$	$(6c)$	O III
$\mathbf{B}_{19}$	=	$x_6\,\mathbf{a}_1+y_6\,\mathbf{a}_2+z_6\,\mathbf{a}_3$	=	$\tfrac{1}{2}(x_6+y_6)\,a\,\hat{\mathbf{x}}+\tfrac{\sqrt{3}}{2}(-x_6+y_6)\,a\,\hat{\mathbf{y}}+z_6 c\,\hat{\mathbf{z}}$	$(12d)$	H
$\mathbf{B}_{20}$	=	$-y_6\,\mathbf{a}_1+(x_6-y_6)\,\mathbf{a}_2+z_6\,\mathbf{a}_3$	=	$\left(\tfrac{1}{2}x_6-y_6\right)a\,\hat{\mathbf{x}}+\tfrac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}}+z_6 c\,\hat{\mathbf{z}}$	$(12d)$	H
$\mathbf{B}_{21}$	=	$(-x_6+y_6)\,\mathbf{a}_1-x_6\,\mathbf{a}_2+z_6\,\mathbf{a}_3$	=	$\left(-x_6+\tfrac{1}{2}y_6\right)a\,\hat{\mathbf{x}}-\tfrac{\sqrt{3}}{2}y_6 a\,\hat{\mathbf{y}}+z_6 c\,\hat{\mathbf{z}}$	$(12d)$	H
$\mathbf{B}_{22}$	=	$-x_6\,\mathbf{a}_1-y_6\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_6\right)\mathbf{a}_3$	=	$-\tfrac{1}{2}(x_6+y_6)\,a\,\hat{\mathbf{x}}+\tfrac{\sqrt{3}}{2}(x_6-y_6)\,a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	$(12d)$	H
$\mathbf{B}_{23}$	=	$y_6\,\mathbf{a}_1+(-x_6+y_6)\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_6\right)\mathbf{a}_3$	=	$\left(-\tfrac{1}{2}x_6+y_6\right)a\,\hat{\mathbf{x}}-\tfrac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	$(12d)$	H
$\mathbf{B}_{24}$	=	$(x_6-y_6)\,\mathbf{a}_1+x_6\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_6\right)\mathbf{a}_3$	=	$\left(x_6-\tfrac{1}{2}y_6\right)a\,\hat{\mathbf{x}}+\tfrac{\sqrt{3}}{2}y_6 a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_6\right)c\,\hat{\mathbf{z}}$	$(12d)$	H

$\mathbf{B}_{25}$	=	$-y_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$-\frac{1}{2}\,(x_6 + y_6)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_6 + y_6)\,a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(12d)$	H
$\mathbf{B}_{26}$	=	$(-x_6 + y_6)\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_6 + y_6\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(12d)$	H
$\mathbf{B}_{27}$	=	$x_6\,\mathbf{a}_1 + (x_6 - y_6)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\left(x_6 - \frac{1}{2}y_6\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(12d)$	H
$\mathbf{B}_{28}$	=	$y_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\frac{1}{2}\,(x_6 + y_6)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(x_6 - y_6)\,a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\,\hat{\mathbf{z}}$	$(12d)$	H
$\mathbf{B}_{29}$	=	$(x_6 - y_6)\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}x_6 - y_6\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\,\hat{\mathbf{z}}$	$(12d)$	H
$\mathbf{B}_{30}$	=	$-x_6\,\mathbf{a}_1 + (-x_6 + y_6)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\left(-x_6 + \frac{1}{2}y_6\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_6 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\,\hat{\mathbf{z}}$	$(12d)$	H

References:

- J.-O. Lundgren, R. Liminga, and R. Tellgren, *Neutron diffraction refinement of pyroelectric lithium perchlorate trihydrate*, Acta Crystallogr. Sect. B Struct. Sci. **38**, 15–20 (1982), doi:10.1107/S0567740882001940.

Found in:

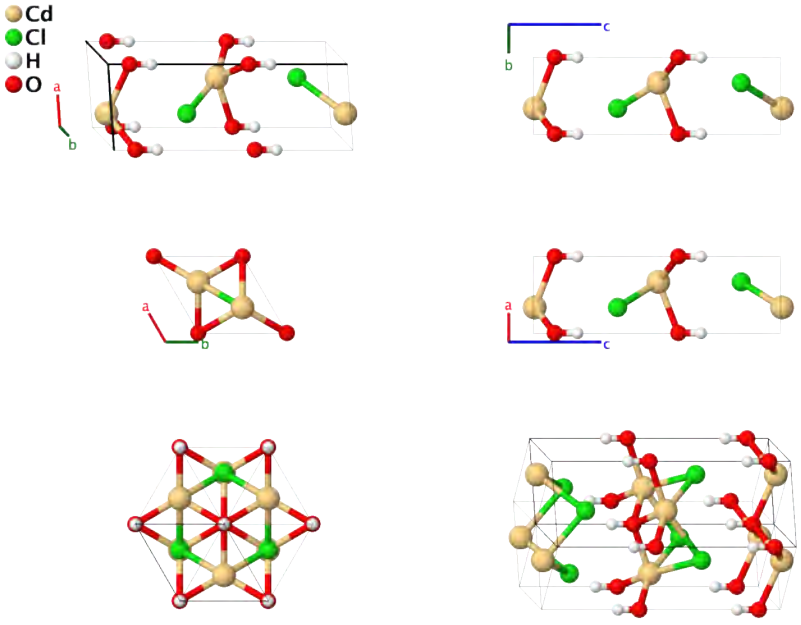
- L. Ojamäe and K. Hermansson, *The OH stretching frequency in $LiClO_4{\cdot}3H_2O(s)$ from ab initio and model potential calculations*, Chem. Phys. **161**, 87–98 (1992), doi:10.1016/0301-0104(92)80179-Y.

Geometry files:

- CIF: pp. 1757
- POSCAR: pp. 1758

Cd(OH)Cl ($E0_3$) Structure: ABCD_hP8_186_b_b_a_a

http://aflow.org/prototype-encyclopedia/ABCD_hP8_186_b_b_a_a

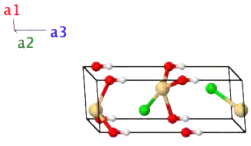


Prototype	:	CdClHO
AFLOW prototype label	:	ABCD_hP8_186_b_b_a_a
***Strukturbericht* designation**	:	$E0_3$
Pearson symbol	:	hP8
Space group number	:	186
Space group symbol	:	$P6_3mc$
AFLOW prototype command	:	`aflow --proto=ABCD_hP8_186_b_b_a_a --params=`$a, c/a, z_1, z_2, z_3, z_4$

- Early determinations of this structure did not locate the hydrogen atoms. (Cudennec, 2000) state that "H atom coordinates were calculated for Cd(OH)Cl."

Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$z_1\,\mathbf{a}_3$	=	$z_1 c\,\hat{\mathbf{z}}$	$(2a)$	H
$\mathbf{B}_2$	=	$\left(\frac{1}{2}+z_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+z_1\right)c\,\hat{\mathbf{z}}$	$(2a)$	H

$\mathbf{B}_3$	=	$z_2\,\mathbf{a}_3$	=	$z_2 c\,\hat{\mathbf{z}}$	($2a$)	O
$\mathbf{B}_4$	=	$\left(\frac{1}{2}+z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+z_2\right)c\,\hat{\mathbf{z}}$	($2a$)	O
$\mathbf{B}_5$	=	$\frac{1}{3}\mathbf{a}_1+\frac{2}{3}\mathbf{a}_2+z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}+\frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+z_3 c\,\hat{\mathbf{z}}$	($2b$)	Cd
$\mathbf{B}_6$	=	$\frac{2}{3}\mathbf{a}_1+\frac{1}{3}\mathbf{a}_2+\left(\frac{1}{2}+z_3\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}-\frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_3\right)c\,\hat{\mathbf{z}}$	($2b$)	Cd
$\mathbf{B}_7$	=	$\frac{1}{3}\mathbf{a}_1+\frac{2}{3}\mathbf{a}_2+z_4\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}+\frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+z_4 c\,\hat{\mathbf{z}}$	($2b$)	Cl
$\mathbf{B}_8$	=	$\frac{2}{3}\mathbf{a}_1+\frac{1}{3}\mathbf{a}_2+\left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}-\frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_4\right)c\,\hat{\mathbf{z}}$	($2b$)	Cl

References:

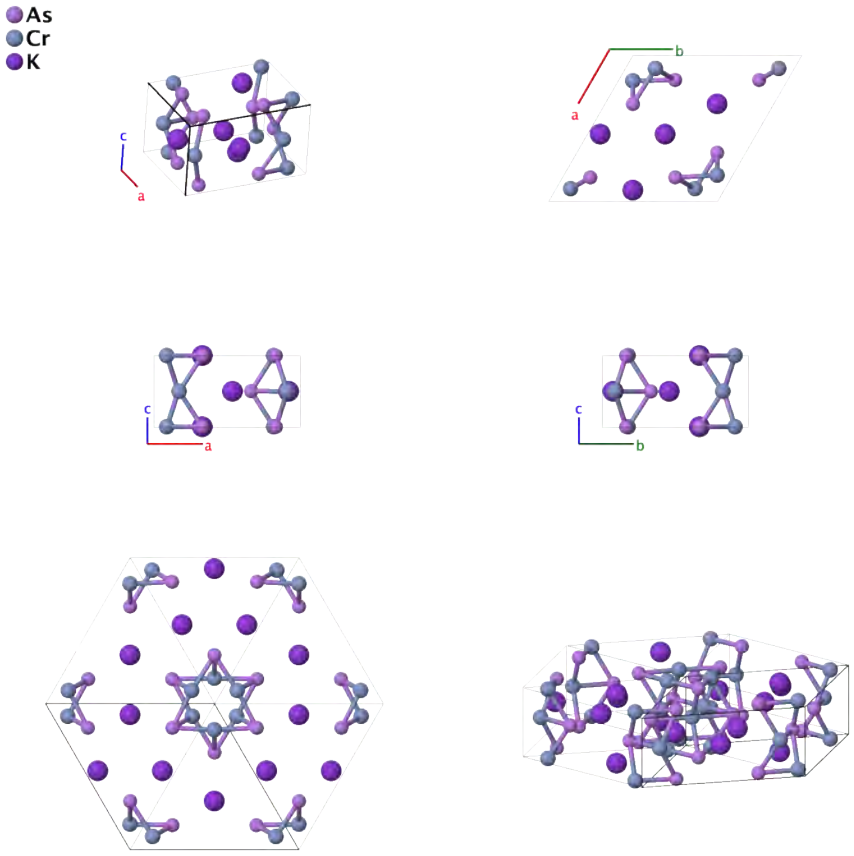
- Y. Cudennec, A. Riou, Y. Gérault, and A. Lecerf, *Synthesis and Crystal Structures of Cd(OH)Cl and Cu(OH)Cl and Relationship to Brucite Type*, J. Solid State Chem. **151**, 308–312 (2000), doi:10.1006/jssc.2000.8659.

Geometry files:

- CIF: pp. 1758
- POSCAR: pp. 1758

Cr-233 Quasi-One-Dimensional Superconductor ($K_2Cr_3As_3$) Structure: A3B3C2_hP16_187_jk_jk_ck

http://aflow.org/prototype-encyclopedia/A3B3C2_hP16_187_jk_jk_ck



Prototype	:	$As_3Cr_3K_2$
AFLOW prototype label	:	A3B3C2_hP16_187_jk_jk_ck
***Strukturbericht* designation**	:	None
Pearson symbol	:	hP16
Space group number	:	187
Space group symbol	:	$P\bar{6}m2$
AFLOW prototype command	:	`aflow --proto=A3B3C2_hP16_187_jk_jk_ck --params=`$a, c/a, x_2, x_3, x_4, x_5, x_6$

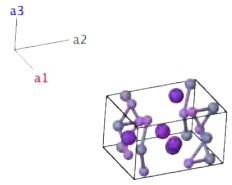
Other compounds with this structure

- $Cs_2Cr_3As_3$, $Rb_2Cr_3As_3$, and $K_2Mo_3As_3$

- Cr-233 designates a class of structures of the form $A_2B_2As_3$, where the 'A' atoms form one-dimensional chains. Several of these compounds have been found to superconduct at temperatures on the order of 5-10 K.

Hexagonal primitive vectors:

$$\mathbf{a}_1 = \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}$	$(1c)$	K I
$\mathbf{B}_2$	=	$x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2$	=	$-\sqrt{3}x_2a\,\hat{\mathbf{y}}$	$(3j)$	As I
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + 2x_2\,\mathbf{a}_2$	=	$\frac{3}{2}x_2a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_2a\,\hat{\mathbf{y}}$	$(3j)$	As I
$\mathbf{B}_4$	=	$-2x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2$	=	$-\frac{3}{2}x_2a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_2a\,\hat{\mathbf{y}}$	$(3j)$	As I
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2$	=	$-\sqrt{3}x_3a\,\hat{\mathbf{y}}$	$(3j)$	Cr I
$\mathbf{B}_6$	=	$x_3\,\mathbf{a}_1 + 2x_3\,\mathbf{a}_2$	=	$\frac{3}{2}x_3a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}}$	$(3j)$	Cr I
$\mathbf{B}_7$	=	$-2x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2$	=	$-\frac{3}{2}x_3a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}}$	$(3j)$	Cr I
$\mathbf{B}_8$	=	$x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-\sqrt{3}x_4a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(3k)$	As II
$\mathbf{B}_9$	=	$x_4\,\mathbf{a}_1 + 2x_4\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{3}{2}x_4a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(3k)$	As II
$\mathbf{B}_{10}$	=	$-2x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-\frac{3}{2}x_4a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(3k)$	As II
$\mathbf{B}_{11}$	=	$x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-\sqrt{3}x_5a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(3k)$	Cr II
$\mathbf{B}_{12}$	=	$x_5\,\mathbf{a}_1 + 2x_5\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{3}{2}x_5a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_5a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(3k)$	Cr II
$\mathbf{B}_{13}$	=	$-2x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-\frac{3}{2}x_5a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_5a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(3k)$	Cr II
$\mathbf{B}_{14}$	=	$x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-\sqrt{3}x_6a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(3k)$	K II
$\mathbf{B}_{15}$	=	$x_6\,\mathbf{a}_1 + 2x_6\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{3}{2}x_6a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_6a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(3k)$	K II
$\mathbf{B}_{16}$	=	$-2x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-\frac{3}{2}x_6a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_6a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(3k)$	K II

References:

- J.-K. Bao, J.-Y. Liu, C.-W. Ma, Z.-H. Meng, Z.-T. Tang, Y.-L. Sun, H.-F. Zhai, H. Jiang, H. Bai, C.-M. Feng, Z.-A. Xu, and G.-H. Cao, *Superconductivity in Quasi-One-Dimensional $K_2Cr_3As_3$ with Significant Electron Correlations*, Phys. Rev. X **5**, 011013 (2015), doi:10.1103/PhysRevX.5.011013.

Found in:

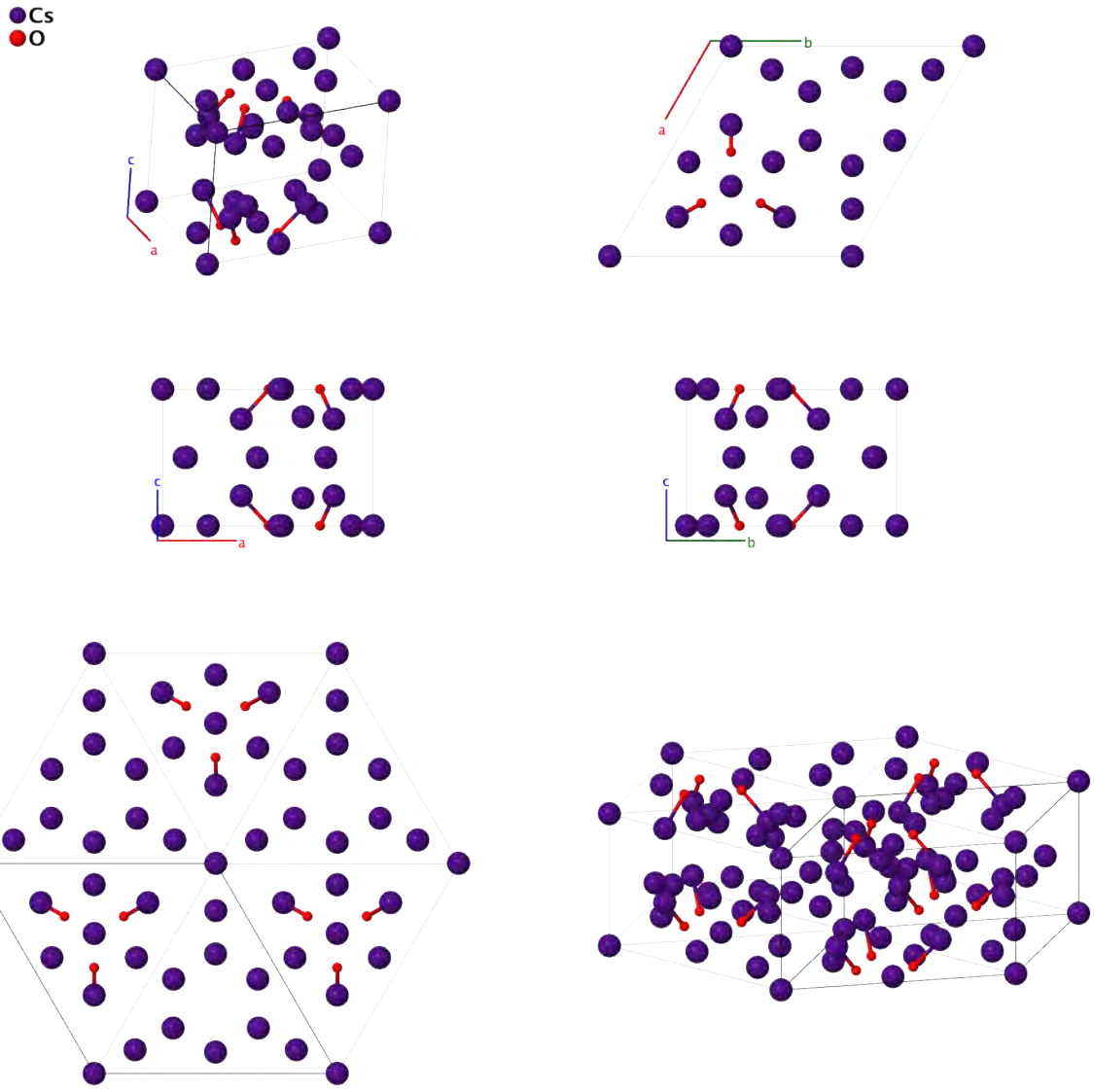
- Q.-G. Mu, B.-B. Ruan, K. Zhao, B.-J. Pan, T. Liu, L. Shan, G.-F. Chen, and Z.-A. Ren, *Superconductivity at 10.4 K in a novel quasi-one-dimensional ternary molybdenum pnictide $K_2Mo_3As_3$*, Sci. Bull. **63**, 952–956 (2018), doi:10.1016/j.scib.2018.06.011.

Geometry files:

- CIF: pp. 1758
- POSCAR: pp. 1759

Cs_7O Structure: A7B_hP24_187_ai2j2kn_j

http://aflow.org/prototype-encyclopedia/A7B_hP24_187_ai2j2kn_j



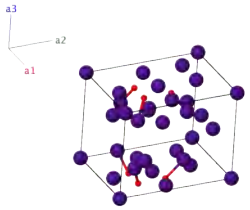
Prototype	:	Cs_7O
AFLOW prototype label	:	A7B_hP24_187_ai2j2kn_j
***Strukturbericht* designation**	:	None
Pearson symbol	:	hP24
Space group number	:	187
Space group symbol	:	$P\bar{6}m2$
AFLOW prototype command	:	`aflow --proto=A7B_hP24_187_ai2j2kn_j --params=`$a, c/a, z_2, x_3, x_4, x_5, x_6, x_7, x_8, z_8$

- This structure is composed of $Cs_{11}O_3$ clusters, similar to the building blocks of the $Cs_{11}O_3$ structure, interlaced with cesium atoms which have approximately the same spacing as in bcc-Cs.

- Lattice constant data was given at -150 °C, while the atomic positions were given at -175 °C.

Hexagonal primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(1a)$	Cs I
$\mathbf{B}_2$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(2i)$	Cs II
$\mathbf{B}_3$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	$(2i)$	Cs II
$\mathbf{B}_4$	=	$x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2$	=	$-\sqrt{3}x_3 a\,\hat{\mathbf{y}}$	$(3j)$	Cs III
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + 2x_3\,\mathbf{a}_2$	=	$\frac{3}{2}x_3 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}}$	$(3j)$	Cs III
$\mathbf{B}_6$	=	$-2x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2$	=	$-\frac{3}{2}x_3 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}}$	$(3j)$	Cs III
$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2$	=	$-\sqrt{3}x_4 a\,\hat{\mathbf{y}}$	$(3j)$	Cs IV
$\mathbf{B}_8$	=	$x_4\,\mathbf{a}_1 + 2x_4\,\mathbf{a}_2$	=	$\frac{3}{2}x_4 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}}$	$(3j)$	Cs IV
$\mathbf{B}_9$	=	$-2x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2$	=	$-\frac{3}{2}x_4 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}}$	$(3j)$	Cs IV
$\mathbf{B}_{10}$	=	$x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2$	=	$-\sqrt{3}x_5 a\,\hat{\mathbf{y}}$	$(3j)$	O
$\mathbf{B}_{11}$	=	$x_5\,\mathbf{a}_1 + 2x_5\,\mathbf{a}_2$	=	$\frac{3}{2}x_5 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}}$	$(3j)$	O
$\mathbf{B}_{12}$	=	$-2x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2$	=	$-\frac{3}{2}x_5 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}}$	$(3j)$	O
$\mathbf{B}_{13}$	=	$x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-\sqrt{3}x_6 a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(3k)$	Cs V
$\mathbf{B}_{14}$	=	$x_6\,\mathbf{a}_1 + 2x_6\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{3}{2}x_6 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(3k)$	Cs V
$\mathbf{B}_{15}$	=	$-2x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-\frac{3}{2}x_6 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(3k)$	Cs V
$\mathbf{B}_{16}$	=	$x_7\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-\sqrt{3}x_7 a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(3k)$	Cs VI
$\mathbf{B}_{17}$	=	$x_7\,\mathbf{a}_1 + 2x_7\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{3}{2}x_7 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_7 a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(3k)$	Cs VI
$\mathbf{B}_{18}$	=	$-2x_7\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-\frac{3}{2}x_7 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_7 a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(3k)$	Cs VI
$\mathbf{B}_{19}$	=	$x_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$-\sqrt{3}x_8 a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	$(6n)$	Cs VII
$\mathbf{B}_{20}$	=	$x_8\,\mathbf{a}_1 + 2x_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$\frac{3}{2}x_8 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_8 a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	$(6n)$	Cs VII
$\mathbf{B}_{21}$	=	$-2x_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$-\frac{3}{2}x_8 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_8 a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	$(6n)$	Cs VII
$\mathbf{B}_{22}$	=	$x_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$-\sqrt{3}x_8 a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	$(6n)$	Cs VII
$\mathbf{B}_{23}$	=	$x_8\,\mathbf{a}_1 + 2x_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$\frac{3}{2}x_8 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_8 a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	$(6n)$	Cs VII
$\mathbf{B}_{24}$	=	$-2x_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$-\frac{3}{2}x_8 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_8 a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	$(6n)$	Cs VII

References:

- A. Simon, *Über Alkalimetall-Suboxide. VII. Das metallreichste Cäsiumoxid-Cs_7O*, Z. Anorg. Allg. Chem. **422**, 208–218 (1976), doi:10.1002/zaac.19764220303.

Found in:

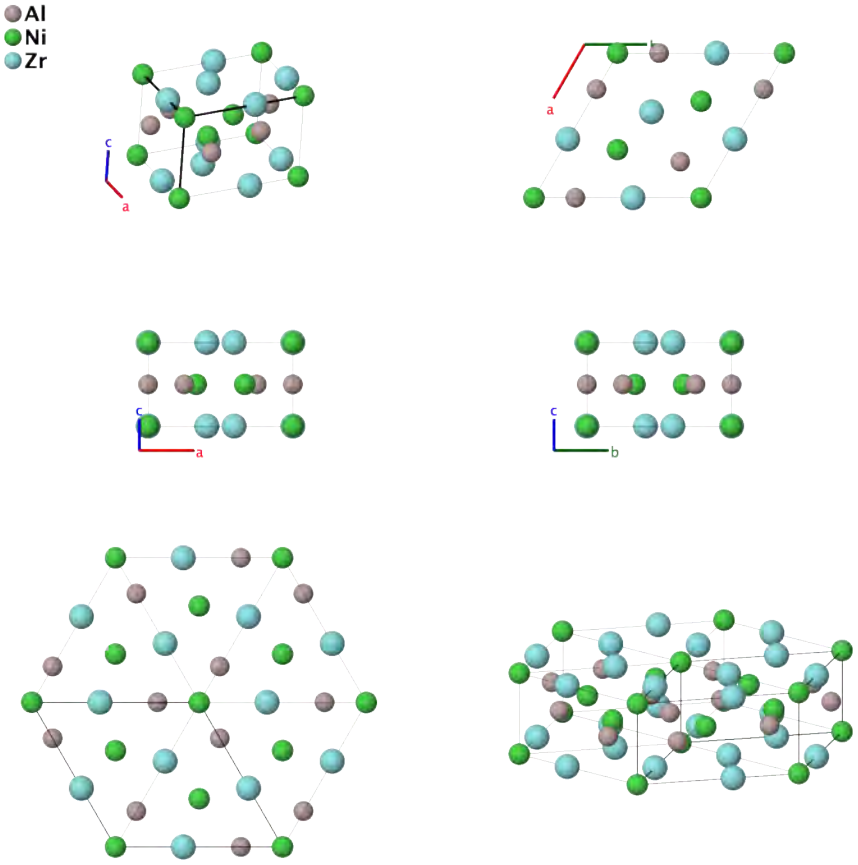
- T. B. Massalski, H. Okamoto, P. R. Subramanian, and L. Kacprzak, eds., *Binary Alloy Phase Diagrams*, vol. 2 (ASM International, Materials Park, Ohio, USA, 1990), 2^{nd} edn. Cd-Ce to Hf-Rb.

Geometry files:

- CIF: pp. 1759
- POSCAR: pp. 1759

ZrNiAl Structure: ABC_hP9_189_g_ad_f

http://aflow.org/prototype-encyclopedia/ABC_hP9_189_g_ad_f



Prototype	:	AlNiZr
AFLOW prototype label	:	ABC_hP9_189_g_ad_f
***Strukturbericht* designation**	:	None
Pearson symbol	:	hP9
Space group number	:	189
Space group symbol	:	$P\bar{6}2m$
AFLOW prototype command	:	`aflow --proto=ABC_hP9_189_g_ad_f --params=`$a, c/a, x_3, x_4$

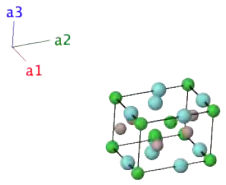
Other compounds with this structure

- AgAsCa, AgSiYb, AlCoPu, AlCuTm, AlNiTb, DyNiIn, DyNiSn, ErNiAl, FeGaU, FeNiP, GdNiIn, GdNiSn, HoNiIn, RhSnZr, RuSiZr, ScIrP, TbNiIn, and BSi_2Ni_6

- This is the ternary form of the Fe_2P structure. In the former case the origin was placed so that the potassium atoms were on the (1*b*) and (2*c*) Wyckoff positions, while here we follow (Shved, 2019) and place the nickel atoms on the (1*a*) and (2*d*) sites. This is merely a shift in the origin by $1/2\, c\, \hat{z}$. If the structures have a common origin then they are nearly identical.
- (Shved, 2019) found evidence of 6-10% mixing between the Ni-II (2*d*) and Al (3*g*) sites.

Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(1a)$	Ni I
$\mathbf{B}_2$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(2d)$	Ni II
$\mathbf{B}_3$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(2d)$	Ni II
$\mathbf{B}_4$	=	$x_3\,\mathbf{a}_1$	=	$\frac{1}{2}x_3a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}}$	$(3f)$	Zr
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_2$	=	$\frac{1}{2}x_3a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}}$	$(3f)$	Zr
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2$	=	$-x_3a\,\hat{\mathbf{x}}$	$(3f)$	Zr
$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}x_4a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(3g)$	Al
$\mathbf{B}_8$	=	$x_4\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}x_4a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(3g)$	Al
$\mathbf{B}_9$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(3g)$	Al

References:

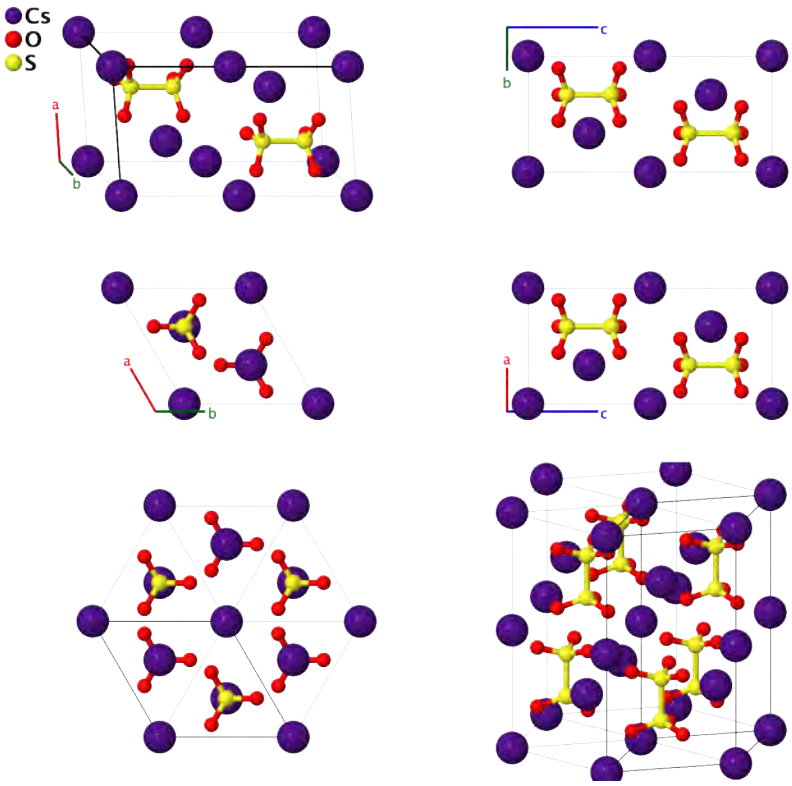
- O. Shved, L. P. Salamakha, S. Mudry, O. Sologub, P. F. Rogl, and E. Bauer, *Zr-based nickel aluminides: crystal structure and electronic properties*, J. Alloys Compd. **821**, 153326 (2020), doi:10.1016/j.jallcom.2019.153326.

Geometry files:

- CIF: pp. 1759
- POSCAR: pp. 1760

$CsSO_3$ ($K1_2$) Structure: AB3C_hP20_190_ac_i_f

http://aflow.org/prototype-encyclopedia/AB3C_hP20_190_ac_i_f

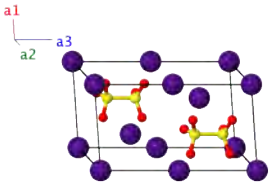


Prototype	:	CsO_3S
AFLOW prototype label	:	AB3C_hP20_190_ac_i_f
***Strukturbericht* designation**	:	$K1_2$
Pearson symbol	:	hP20
Space group number	:	190
Space group symbol	:	$P\bar{6}2c$
AFLOW prototype command	:	`aflow --proto=AB3C_hP20_190_ac_i_f --params=`$a, c/a, z_3, x_4, y_4, z_4$

- (Hägg, 1932) states that the space group might be $P6_322$ #182 or $P\bar{6}2c$ #190. We follow (Downs, 2003) and use the latter. (Hägg, 1932) and (Downs, 2003) set the sulfur coordinate $z_3 = 0.73$. As noted by (Hermann, 1937), this gives a very short S-S bond distance, on the order of 0.5Å. We use the value suggested by (Hermann, 1937), $z_3 = 0.63$, "which gives the S_2O_6 group exactly the same shape it has in $K_2S_2O_6$ [*Strukturbericht* $K1_1$]."

Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	Cs I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	$(2a)$	Cs I
$\mathbf{B}_3$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(2c)$	Cs II
$\mathbf{B}_4$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(2c)$	Cs II
$\mathbf{B}_5$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_3c\,\hat{\mathbf{z}}$	$(4f)$	S
$\mathbf{B}_6$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	$(4f)$	S
$\mathbf{B}_7$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_3c\,\hat{\mathbf{z}}$	$(4f)$	S
$\mathbf{B}_8$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	$(4f)$	S
$\mathbf{B}_9$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_4 + y_4)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_4 + y_4)\,a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	$(12i)$	O
$\mathbf{B}_{10}$	=	$-y_4\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_4 - y_4\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	$(12i)$	O
$\mathbf{B}_{11}$	=	$(-x_4 + y_4)\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\left(-x_4 + \frac{1}{2}y_4\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_4a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	$(12i)$	O
$\mathbf{B}_{12}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_4 + y_4)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_4 + y_4)\,a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	$(12i)$	O
$\mathbf{B}_{13}$	=	$-y_4\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_4 - y_4\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	$(12i)$	O
$\mathbf{B}_{14}$	=	$(-x_4 + y_4)\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\,\mathbf{a}_3$	=	$\left(-x_4 + \frac{1}{2}y_4\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_4a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	$(12i)$	O
$\mathbf{B}_{15}$	=	$y_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_4 + y_4)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(x_4 - y_4)\,a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	$(12i)$	O
$\mathbf{B}_{16}$	=	$(x_4 - y_4)\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_4 - y_4\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	$(12i)$	O
$\mathbf{B}_{17}$	=	$-x_4\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\left(-x_4 + \frac{1}{2}y_4\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_4a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	$(12i)$	O
$\mathbf{B}_{18}$	=	$y_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_4 + y_4)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(x_4 - y_4)\,a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	$(12i)$	O
$\mathbf{B}_{19}$	=	$(x_4 - y_4)\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_4 - y_4\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	$(12i)$	O
$\mathbf{B}_{20}$	=	$-x_4\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\,\mathbf{a}_3$	=	$\left(-x_4 + \frac{1}{2}y_4\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_4a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	$(12i)$	O

References:

- G. Hägg, *Die Kristallstruktur von Caesiumdithionat, $Cs_2S_2O_6$*, Z. Physik. Chem. B **18**, 327–342 (1932), doi:10.1515/zpch-1932-1825.
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Found in:

- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

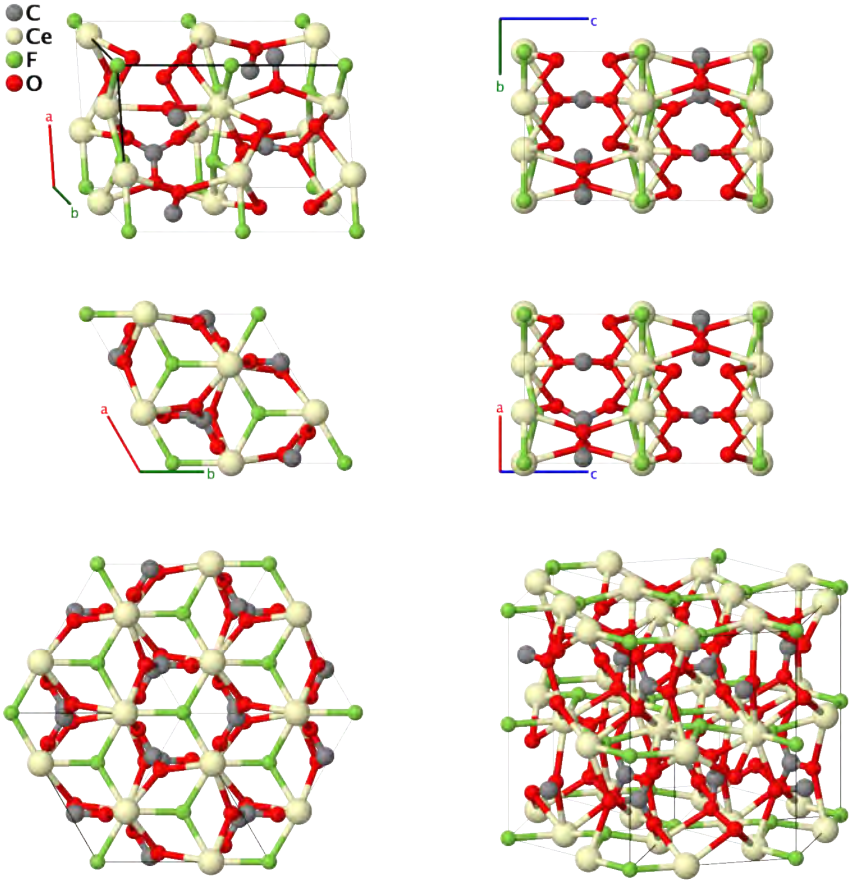
Geometry files:

- CIF: pp. 1760

- POSCAR: pp. 1760

Bastnäsite [$CeF(CO_3)$] Structure: ABCD3_hP36_190_h_g_af_hi

http://aflow.org/prototype-encyclopedia/ABCD3_hP36_190_h_g_af_hi



Prototype	:	$CCeFO_3$
AFLOW prototype label	:	ABCD3_hP36_190_h_g_af_hi
***Strukturbericht* designation**	:	$G7_1$
Pearson symbol	:	hP36
Space group number	:	190
Space group symbol	:	$P\bar{6}2c$
AFLOW prototype command	:	`aflow --proto=ABCD3_hP36_190_h_g_af_hi --params=`$a, c/a, z_2, x_3, x_4, y_4, x_5, y_5, x_6, y_6, z_6$

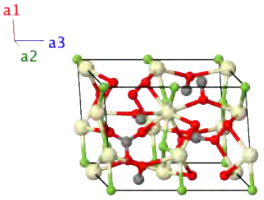
Other compounds with this structure

- $LaF(CO_3)$

- Technically the name bastnäsite belongs to both cerium and lanthanum compounds, and the are usually designated bastnäsite-(Ce) and bastnäsite-(La), respectively.

Hexagonal primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	F I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	$(2a)$	F I
$\mathbf{B}_3$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(4f)$	F II
$\mathbf{B}_4$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	$(4f)$	F II
$\mathbf{B}_5$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	$(4f)$	F II
$\mathbf{B}_6$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	$(4f)$	F II
$\mathbf{B}_7$	=	$x_3\,\mathbf{a}_1$	=	$\frac{1}{2}x_3 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}}$	$(6g)$	Ce
$\mathbf{B}_8$	=	$x_3\,\mathbf{a}_2$	=	$\frac{1}{2}x_3 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}}$	$(6g)$	Ce
$\mathbf{B}_9$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2$	=	$-x_3 a\,\hat{\mathbf{x}}$	$(6g)$	Ce
$\mathbf{B}_{10}$	=	$x_3\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}x_3 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(6g)$	Ce
$\mathbf{B}_{11}$	=	$x_3\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}x_3 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(6g)$	Ce
$\mathbf{B}_{12}$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(6g)$	Ce
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}(x_4 + y_4)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(-x_4 + y_4)\,a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	C
$\mathbf{B}_{14}$	=	$-y_4\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_4 - y_4\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	C
$\mathbf{B}_{15}$	=	$(-x_4 + y_4)\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(-x_4 + \frac{1}{2}y_4\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_4 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	C
$\mathbf{B}_{16}$	=	$y_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}(x_4 + y_4)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(x_4 - y_4)\,a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	C
$\mathbf{B}_{17}$	=	$(x_4 - y_4)\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_4 - y_4\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	C
$\mathbf{B}_{18}$	=	$-x_4\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(-x_4 + \frac{1}{2}y_4\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_4 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	C
$\mathbf{B}_{19}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}(x_5 + y_5)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(-x_5 + y_5)\,a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O I
$\mathbf{B}_{20}$	=	$-y_5\,\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_5 - y_5\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O I
$\mathbf{B}_{21}$	=	$(-x_5 + y_5)\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(-x_5 + \frac{1}{2}y_5\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_5 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O I
$\mathbf{B}_{22}$	=	$y_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}(x_5 + y_5)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(x_5 - y_5)\,a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O I
$\mathbf{B}_{23}$	=	$(x_5 - y_5)\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_5 - y_5\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O I
$\mathbf{B}_{24}$	=	$-x_5\,\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(-x_5 + \frac{1}{2}y_5\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_5 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O I
$\mathbf{B}_{25}$	=	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\frac{1}{2}(x_6 + y_6)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}(-x_6 + y_6)\,a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(12i)$	O II

$$\mathbf{B}_{26} = -y_6\,\mathbf{a}_1 + (x_6 - y_6)\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 = \left(\tfrac{1}{2}x_6 - y_6\right)a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}} \quad (12i) \quad \text{O II}$$

$$\mathbf{B}_{27} = (-x_6 + y_6)\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 = \left(-x_6 + \tfrac{1}{2}y_6\right)a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}y_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}} \quad (12i) \quad \text{O II}$$

$$\mathbf{B}_{28} = x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_3 = \tfrac{1}{2}(x_6 + y_6)\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}(-x_6 + y_6)\,a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_6\right)c\,\hat{\mathbf{z}} \quad (12i) \quad \text{O II}$$

$$\mathbf{B}_{29} = -y_6\,\mathbf{a}_1 + (x_6 - y_6)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}x_6 - y_6\right)a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_6\right)c\,\hat{\mathbf{z}} \quad (12i) \quad \text{O II}$$

$$\mathbf{B}_{30} = (-x_6 + y_6)\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_3 = \left(-x_6 + \tfrac{1}{2}y_6\right)a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}y_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_6\right)c\,\hat{\mathbf{z}} \quad (12i) \quad \text{O II}$$

$$\mathbf{B}_{31} = y_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3 = \tfrac{1}{2}(x_6 + y_6)\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}(x_6 - y_6)\,a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}} \quad (12i) \quad \text{O II}$$

$$\mathbf{B}_{32} = (x_6 - y_6)\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3 = \left(\tfrac{1}{2}x_6 - y_6\right)a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}} \quad (12i) \quad \text{O II}$$

$$\mathbf{B}_{33} = -x_6\,\mathbf{a}_1 + (-x_6 + y_6)\,\mathbf{a}_2 - z_6\,\mathbf{a}_3 = \left(-x_6 + \tfrac{1}{2}y_6\right)a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}y_6 a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}} \quad (12i) \quad \text{O II}$$

$$\mathbf{B}_{34} = y_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 = \tfrac{1}{2}(x_6 + y_6)\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}(x_6 - y_6)\,a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right)c\,\hat{\mathbf{z}} \quad (12i) \quad \text{O II}$$

$$\mathbf{B}_{35} = (x_6 - y_6)\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}x_6 - y_6\right)a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right)c\,\hat{\mathbf{z}} \quad (12i) \quad \text{O II}$$

$$\mathbf{B}_{36} = -x_6\,\mathbf{a}_1 + (-x_6 + y_6)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 = \left(-x_6 + \tfrac{1}{2}y_6\right)a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}y_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right)c\,\hat{\mathbf{z}} \quad (12i) \quad \text{O II}$$

References:

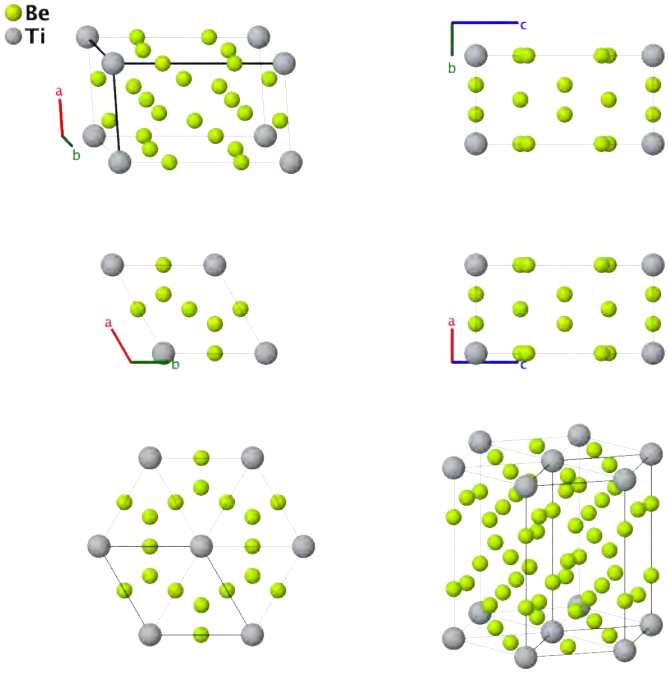
- Y. Ni, J. M. Hughes, and A. N. Mariano, *The atomic arrangement of bastnäsite-(Ce), Ce(CO_3)F, and structural elements of synchysite-(Ce), röntgenite-(Ce), and parisite-(Ce)*, Am. Mineral. **78**, 415–418 (1993).

Geometry files:

- CIF: pp. 1760
- POSCAR: pp. 1761

$TiBe_{12}$ (approximate, $D2_a$) Structure: A12B_hP13_191_cdei_a

http://aflow.org/prototype-encyclopedia/A12B_hP13_191_cdei_a

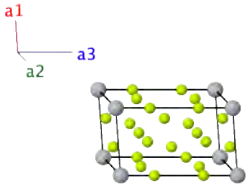


Prototype	:	$Be_{12}Ti$
AFLOW prototype label	:	A12B_hP13_191_cdei_a
***Strukturbericht* designation**	:	$D2_a$
Pearson symbol	:	hP13
Space group number	:	191
Space group symbol	:	$P6/mmm$
AFLOW prototype command	:	`aflow --proto=A12B_hP13_191_cdei_a --params=`$a, c/a, z_4, z_5$

- The structure of $TiBe_{12}$ is not settled. (Raeuchle, 1952) described the structure as a somewhat disordered supercell containing 48 atoms with lattice constants $a = 29.44$ Å and $c = 7.33$ Å, but they stated that a 'pseudo-cell' existed with dimensions $a = 4.23$ Å and $c = 7.33$ Å. This pseudo-cell is described here, and was assigned the $D2_a$ *Strukturbericht* type by Smithells (Brandes, 1992). Raeuchle and Rundle suggested that the larger primitive cell was constructed from the multiple pseudo-cells, with the titanium atom alternating between the ($1a$) and ($1b$) $(0, 0, 1/2)$ Wyckoff positions.
- Other experimenters have suggested that the actual structure of $TiBe_{12}$ is tetragonal. (Jackson, 2016) presents first-principles calculations which suggest that the actual structure is of the tetragonal $ThMn_{12}$ ($D2_b$) type.

Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(1*a*)	Ti
$\mathbf{B}_2$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}$	(2*c*)	Be I
$\mathbf{B}_3$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}$	(2*c*)	Be I
$\mathbf{B}_4$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(2*d*)	Be II
$\mathbf{B}_5$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(2*d*)	Be II
$\mathbf{B}_6$	=	$z_4\,\mathbf{a}_3$	=	$z_4 c\,\hat{\mathbf{z}}$	(2*e*)	Be III
$\mathbf{B}_7$	=	$-z_4\,\mathbf{a}_3$	=	$-z_4 c\,\hat{\mathbf{z}}$	(2*e*)	Be III
$\mathbf{B}_8$	=	$\frac{1}{2}\,\mathbf{a}_1 + z_5\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{4}a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(6*i*)	Be IV
$\mathbf{B}_9$	=	$\frac{1}{2}\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{4}a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(6*i*)	Be IV
$\mathbf{B}_{10}$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + z_5 c\,\hat{\mathbf{z}}$	(6*i*)	Be IV
$\mathbf{B}_{11}$	=	$\frac{1}{2}\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{4}a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(6*i*)	Be IV
$\mathbf{B}_{12}$	=	$\frac{1}{2}\,\mathbf{a}_1 - z_5\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{4}a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(6*i*)	Be IV
$\mathbf{B}_{13}$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - z_5 c\,\hat{\mathbf{z}}$	(6*i*)	Be IV

References:

- R. F. Raeuchle and R. E. Rundle, *The Structure of $TiBe_{12}$*, Acta Cryst. **5**, 85–93 (1952), doi:10.1107/S0365110X52000186.
- E. A. Brandes and G. B. Brook, eds., *Smithells Metals Reference Book* (Butterworth Heinemann, Oxford, Auckland, Boston, Johannesburg, Melbourne, New Delhi, 1992), seventh edn. Strukturbericht designations start on page 6-36 (163 in PDF), see table 6.3 on page 6-63 (190).

Found in:

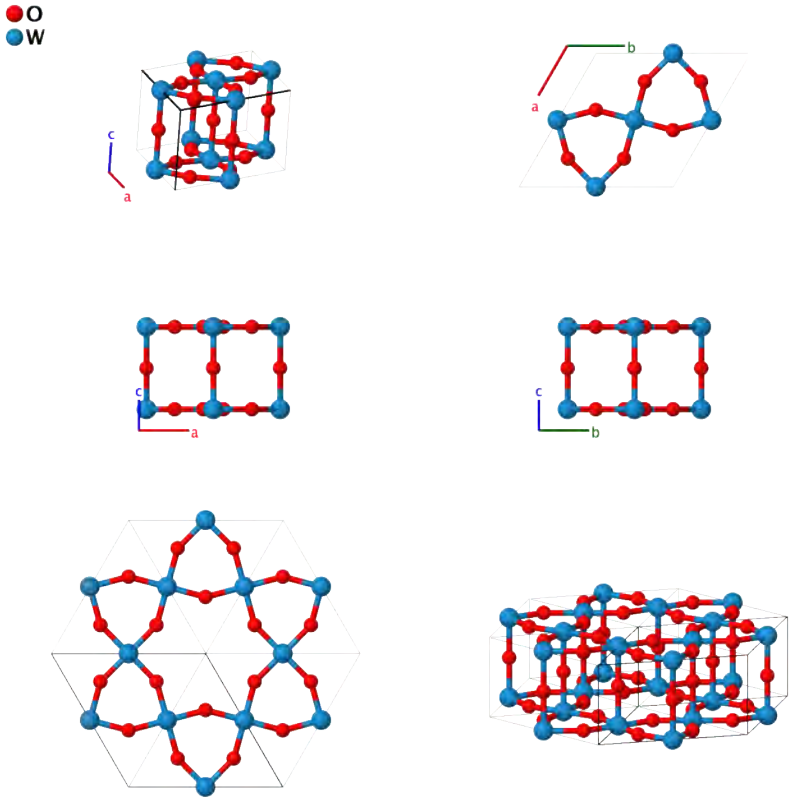
- M. L. Jackson, P. A. Burr, and R. W. Grimes, *Resolving the structure of $TiBe_{12}$*, Acta Crystallogr. Sect. B Struct. Sci. **72**, 277–280 (2016), doi:10.1107/S205252061600322X.

Geometry files:

- CIF: pp. 1761
- POSCAR: pp. 1761

Hexagonal WO_3 Structure: A3B_hP12_191_gl_f

http://aflow.org/prototype-encyclopedia/A3B_hP12_191_gl_f



Prototype	:	O_3W
AFLOW prototype label	:	A3B_hP12_191_gl_f
***Strukturbericht* designation**	:	None
Pearson symbol	:	hP12
Space group number	:	191
Space group symbol	:	$P6/mmm$
AFLOW prototype command	:	`aflow --proto=A3B_hP12_191_gl_f --params=`$a, c/a, x_3$

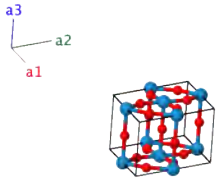
- All stable phases of WO_3 are distortions of the cubic α-ReO_3 ($D0_9$) phase. Based on (Woodward, 1997 and Vogt, 1999), the known stable phases and their approximate temperature ranges are:
  - α-WO_3 (1010-1170 K) (Vogt, 1999)
  - β-WO_3 (600-1170 K) (Vogt, 1999)
  - γ-WO_3 (290-600 K) (Vogt, 1999)
  - δ-WO_3 (230-290 K) (Diehl, 1978)
  - ϵ-WO_3 (below 23 K) (Woodward, 1997)

  Woodward's caption says "The transition temperatures display large hysteresis effects and universal agreement is not found in the literature."

- In addition, several other structures have been proposed and/or found:
  - The original $D0_{10}$ structure (Bräkken, 1931), (Hermann, 1937) superseded by δ-WO_3
  - Original β-WO_3 (Salje, 1977)
  - Hexagonal WO_3 (Gerand, 1979) (metastable), this structure

Hexagonal primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}\\ \mathbf{a}_2 &= \tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{2}\mathbf{a}_1$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{4}a\,\hat{\mathbf{y}}$	$(3f)$	W
$\mathbf{B}_2$	=	$\frac{1}{2}\mathbf{a}_2$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{4}a\,\hat{\mathbf{y}}$	$(3f)$	W
$\mathbf{B}_3$	=	$\frac{1}{2}\mathbf{a}_1 + \frac{1}{2}\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	$(3f)$	W
$\mathbf{B}_4$	=	$\frac{1}{2}\mathbf{a}_1 + \frac{1}{2}\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{4}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(3g)$	O I
$\mathbf{B}_5$	=	$\frac{1}{2}\mathbf{a}_2 + \frac{1}{2}\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{4}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(3g)$	O I
$\mathbf{B}_6$	=	$\frac{1}{2}\mathbf{a}_1 + \frac{1}{2}\mathbf{a}_2 + \frac{1}{2}\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(3g)$	O I
$\mathbf{B}_7$	=	$x_3\,\mathbf{a}_1 + 2x_3\,\mathbf{a}_2$	=	$\frac{3}{2}x_3a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}}$	$(6l)$	O II
$\mathbf{B}_8$	=	$-2x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2$	=	$-\frac{3}{2}x_3a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}}$	$(6l)$	O II
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2$	=	$-\sqrt{3}x_3a\,\hat{\mathbf{y}}$	$(6l)$	O II
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 - 2x_3\,\mathbf{a}_2$	=	$-\frac{3}{2}x_3a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}}$	$(6l)$	O II
$\mathbf{B}_{11}$	=	$2x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2$	=	$\frac{3}{2}x_3a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}}$	$(6l)$	O II
$\mathbf{B}_{12}$	=	$-x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2$	=	$\sqrt{3}x_3a\,\hat{\mathbf{y}}$	$(6l)$	O II

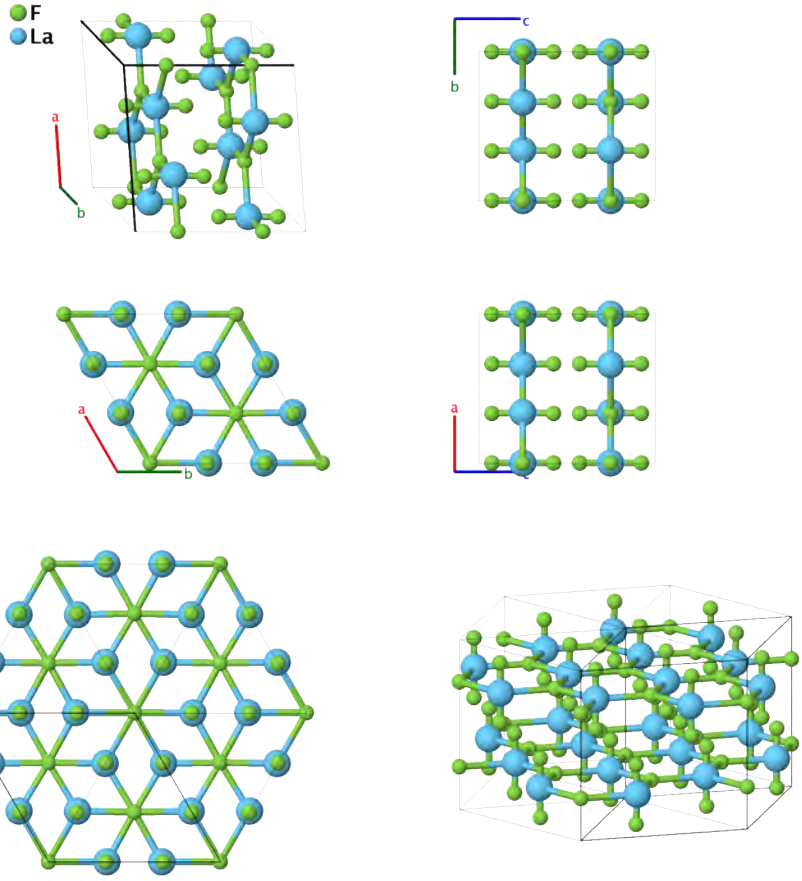
References:

- P. M. Woodward, A. W. Sleight, and T. Vogt, *Ferroelectric Tungsten Trioxide*, J. Solid State Chem. **131**, 9–17 (1997), doi:10.1006/jssc.1997.7268.
- T. Vogt, P. M. Woodward, and B. A. Hunter, *The High-Temperature Phases of WO_3*, J. Solid State Chem. **144**, 209–215 (1999), doi:10.1006/jssc.1999.8173.
- R. Diehl, G. Brandt, and E. Salje, *The Crystal Structure of Triclinic WO_3*, Acta Crystallogr. Sect. B Struct. Sci. **34**, 1105–1111 (1978), doi:10.1107/S0567740878005014.
- H. Bräkken, *Die Kristallstrukturen der Trioxyde von Chrom, Molybdän und Wolfram*, Zeitschrift für Kristallographie - Crystalline Materials **78**, 484–488 (1931), doi:10.1524/zkri.1931.78.1.484.
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).
- E. Salje, *The Orthorhombic Phase of WO_3*, Acta Crystallogr. Sect. B Struct. Sci. **33**, 574–577 (1977), doi:10.1107/S0567740877004130.
- B. Gerand, G. Nowogrocki, J. Guenot, and M. Figlarz, *Structural study of a new hexagonal form of tungsten trioxide*, J. Solid State Chem. **29**, 429–434 (1979), doi:10.1016/0022-4596(79)90199-3.

Geometry files:
- CIF: pp. 1762
- POSCAR: pp. 1762

$D0_6$ (Tysonite, LaF_3) (*obsolete*) Structure: A3B_hP24_193_ack_g

http://aflow.org/prototype-encyclopedia/A3B_hP24_193_ack_g

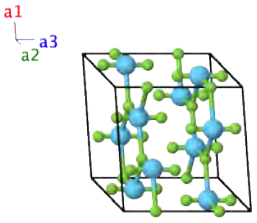


Prototype	:	F_3La
AFLOW prototype label	:	A3B_hP24_193_ack_g
***Strukturbericht* designation**	:	$D0_6$
Pearson symbol	:	hP24
Space group number	:	193
Space group symbol	:	$P6_3/mcm$
AFLOW prototype command	:	`aflow --proto=A3B_hP24_193_ack_g --params=`$a, c/a, x_3, x_4, z_4$

- This stucture was one of several considered by (Hermann, 1937) as a candidate structure for tysonite, and was chosen because it gave the best positioning of the fluorine atoms. Later, (Zalkin, 1985) showed that the structure was trigonal rather than hexagonal, and isostructural with Cu_3P ($D0_{21}$). We keep the original structure as part of the historical record.

Hexagonal primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\,\hat{\mathbf{z}}$	(2*a*)	F I
$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\,\hat{\mathbf{z}}$	(2*a*)	F I
$\mathbf{B}_3$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*c*)	F II
$\mathbf{B}_4$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(4*c*)	F II
$\mathbf{B}_5$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(4*c*)	F II
$\mathbf{B}_6$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(4*c*)	F II
$\mathbf{B}_7$	=	$x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}x_3a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*g*)	La
$\mathbf{B}_8$	=	$x_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}x_3a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*g*)	La
$\mathbf{B}_9$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*g*)	La
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_3$	=	$-\frac{1}{2}x_3a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*g*)	La
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-\frac{1}{2}x_3a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*g*)	La
$\mathbf{B}_{12}$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*g*)	La
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + z_4\,\mathbf{a}_3$	=	$\frac{1}{2}x_4a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	(12*k*)	F III
$\mathbf{B}_{14}$	=	$x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\frac{1}{2}x_4a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4a\,\hat{\mathbf{y}} + z_4c\,\hat{\mathbf{z}}$	(12*k*)	F III
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + z_4c\,\hat{\mathbf{z}}$	(12*k*)	F III
$\mathbf{B}_{16}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$-\frac{1}{2}x_4a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(12*k*)	F III
$\mathbf{B}_{17}$	=	$-x_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$-\frac{1}{2}x_4a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(12*k*)	F III
$\mathbf{B}_{18}$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(12*k*)	F III
$\mathbf{B}_{19}$	=	$x_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\frac{1}{2}x_4a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(12*k*)	F III
$\mathbf{B}_{20}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\frac{1}{2}x_4a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(12*k*)	F III
$\mathbf{B}_{21}$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(12*k*)	F III
$\mathbf{B}_{22}$	=	$-x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-\frac{1}{2}x_4a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	(12*k*)	F III
$\mathbf{B}_{23}$	=	$-x_4\,\mathbf{a}_1 - z_4\,\mathbf{a}_3$	=	$-\frac{1}{2}x_4a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4a\,\hat{\mathbf{y}} - z_4c\,\hat{\mathbf{z}}$	(12*k*)	F III
$\mathbf{B}_{24}$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} - z_4c\,\hat{\mathbf{z}}$	(12*k*)	F III

References:

- I. Oftedal, *Zur Kristallstruktur von Tysonit (Ce, La, ...)F_3*, Z. Phys. B: Condens. Matter **13**, 190–200 (1931), doi:10.1515/zpch-1931-1315.

- A. Zalkin and D. H. Templeton, *Refinement of the trigonal crystal structure of lanthanum trifluoride with neutron diffraction data*, Acta Crystallogr. Sect. B Struct. Sci. **41**, 91–93 (1985), doi:10.1107/S0108768185001689.

Found in:

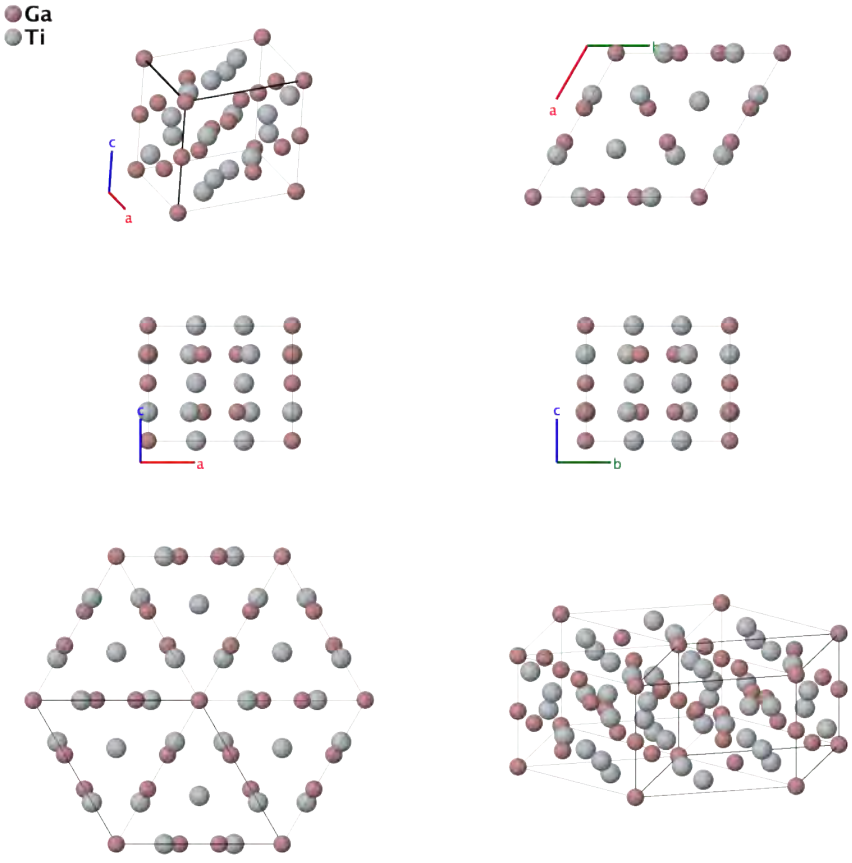
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1762
- POSCAR: pp. 1763

Ti_5Ga_4 Structure: A4B5_hP18_193_bg_dg

http://aflow.org/prototype-encyclopedia/A4B5_hP18_193_bg_dg



Prototype	:	Ga_4Ti_5
AFLOW prototype label	:	A4B5_hP18_193_bg_dg
***Strukturbericht* designation**	:	None
Pearson symbol	:	hP18
Space group number	:	193
Space group symbol	:	$P6_3/mcm$
AFLOW prototype command	:	`aflow --proto=A4B5_hP18_193_bg_dg --params=`$a, c/a, x_3, x_4$

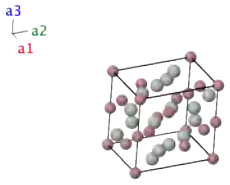
Other compounds with this structure

- Zr_5Ga_4, Hf_5Sn_4, Ba_3TlTe_5, and NbIrO

- This is the $D8_8$ structure with the addition of two atoms on the ($2b$) site.

Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(2*b*)	Ga I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	(2*b*)	Ga I
$\mathbf{B}_3$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}$	(4*d*)	Ti I
$\mathbf{B}_4$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*d*)	Ti I
$\mathbf{B}_5$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}}$	(4*d*)	Ti I
$\mathbf{B}_6$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4*d*)	Ti I
$\mathbf{B}_7$	=	$x_3\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}x_3a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*g*)	Ga II
$\mathbf{B}_8$	=	$x_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}x_3a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*g*)	Ga II
$\mathbf{B}_9$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*g*)	Ga II
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_3$	=	$-\frac{1}{2}x_3a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*g*)	Ga II
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-\frac{1}{2}x_3a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*g*)	Ga II
$\mathbf{B}_{12}$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*g*)	Ga II
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}x_4a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*g*)	Ti II
$\mathbf{B}_{14}$	=	$x_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}x_4a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*g*)	Ti II
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*g*)	Ti II
$\mathbf{B}_{16}$	=	$-x_4\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_3$	=	$-\frac{1}{2}x_4a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*g*)	Ti II
$\mathbf{B}_{17}$	=	$-x_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-\frac{1}{2}x_4a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*g*)	Ti II
$\mathbf{B}_{18}$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*g*)	Ti II

References:

- K. Schubert, H. G. Meissner, M. Pötzschke, W. Rossteutscher, and E. Stolz, *Einige Strukturdaten metallischer Phasen (7)*, Naturwissenschaften **49**, 57 (1962), doi:10.1007/BF00595382.

Found in:

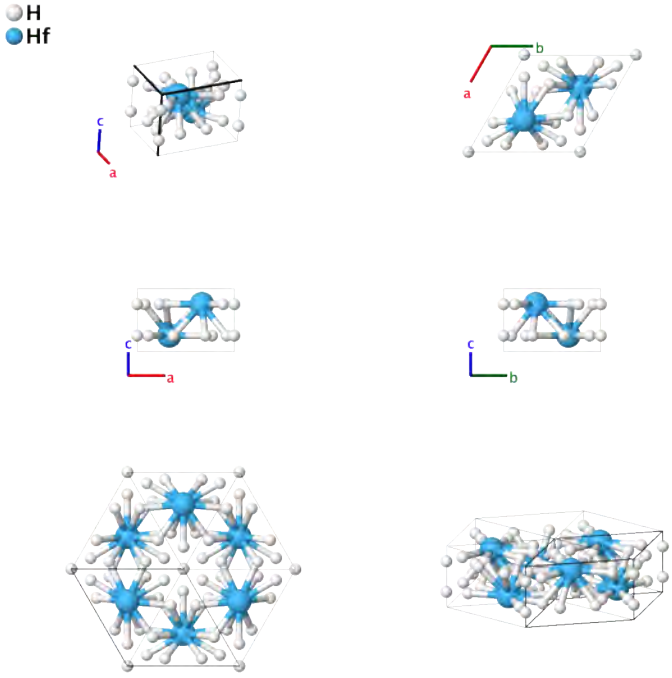
- P. Villars (Chief Editor), *Ti_5Ga_4 Crystal Structure*, http://materials.springer.com/isp/crystallographic/docs/sd_0528839 (2016). PAULING FILE in: Inorganic Solid Phases, SpringerMaterials (online database), Springer, Heidelberg (ed.) SpringerMaterials.

Geometry files:

- CIF: pp. 1763
- POSCAR: pp. 1763

Proposed 300 GPa HfH_{10} Structure: A10B_hP22_194_bhj_c

http://aflow.org/prototype-encyclopedia/A10B_hP22_194_bhj_c



Prototype	:	$H_{10}Hf$
AFLOW prototype label	:	A10B_hP22_194_bhj_c
***Strukturbericht* designation**	:	None
Pearson symbol	:	hP22
Space group number	:	194
Space group symbol	:	$P6_3/mmc$
AFLOW prototype command	:	`aflow --proto=A10B_hP22_194_bhj_c --params=`$a, c/a, x_3, x_4, y_4$

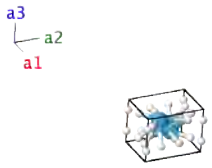
Other compounds with this structure

- LuH_{10}, ScH_{10}, and ZrH_{10}

- This structure has only been determined computationally. At 300 GPa it is predicted to have a superconducting transitions temperature T_c above 200 K.

Hexagonal primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\,\hat{\mathbf{z}}$	(2*b*)	H I
$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\,\hat{\mathbf{z}}$	(2*b*)	H I
$\mathbf{B}_3$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(2*c*)	Hf
$\mathbf{B}_4$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(2*c*)	Hf
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + 2x_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{3}{2}x_3a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	H II
$\mathbf{B}_6$	=	$-2x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-\frac{3}{2}x_3a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	H II
$\mathbf{B}_7$	=	$x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-\sqrt{3}x_3a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	H II
$\mathbf{B}_8$	=	$-x_3\,\mathbf{a}_1 - 2x_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-\frac{3}{2}x_3a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	H II
$\mathbf{B}_9$	=	$2x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{2}x_3a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	H II
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\sqrt{3}x_3a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	H II
$\mathbf{B}_{11}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_4 + y_4)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_4 + y_4)\,a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(12*j*)	H III
$\mathbf{B}_{12}$	=	$-y_4\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_4 - y_4\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(12*j*)	H III
$\mathbf{B}_{13}$	=	$(-x_4 + y_4)\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(-x_4 + \frac{1}{2}y_4\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_4a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(12*j*)	H III
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-\frac{1}{2}\,(x_4 + y_4)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(x_4 - y_4)\,a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(12*j*)	H III
$\mathbf{B}_{15}$	=	$y_4\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_4 + y_4\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(12*j*)	H III
$\mathbf{B}_{16}$	=	$(x_4 - y_4)\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(x_4 - \frac{1}{2}y_4\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_4a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(12*j*)	H III
$\mathbf{B}_{17}$	=	$y_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}\,(x_4 + y_4)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(x_4 - y_4)\,a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(12*j*)	H III
$\mathbf{B}_{18}$	=	$(x_4 - y_4)\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}x_4 - y_4\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(12*j*)	H III
$\mathbf{B}_{19}$	=	$-x_4\,\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(-x_4 + \frac{1}{2}y_4\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}y_4a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(12*j*)	H III
$\mathbf{B}_{20}$	=	$-y_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-\frac{1}{2}\,(x_4 + y_4)\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,(-x_4 + y_4)\,a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(12*j*)	H III
$\mathbf{B}_{21}$	=	$(-x_4 + y_4)\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(-\frac{1}{2}x_4 + y_4\right)a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(12*j*)	H III
$\mathbf{B}_{22}$	=	$x_4\,\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(x_4 - \frac{1}{2}y_4\right)a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}y_4a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(12*j*)	H III

References:

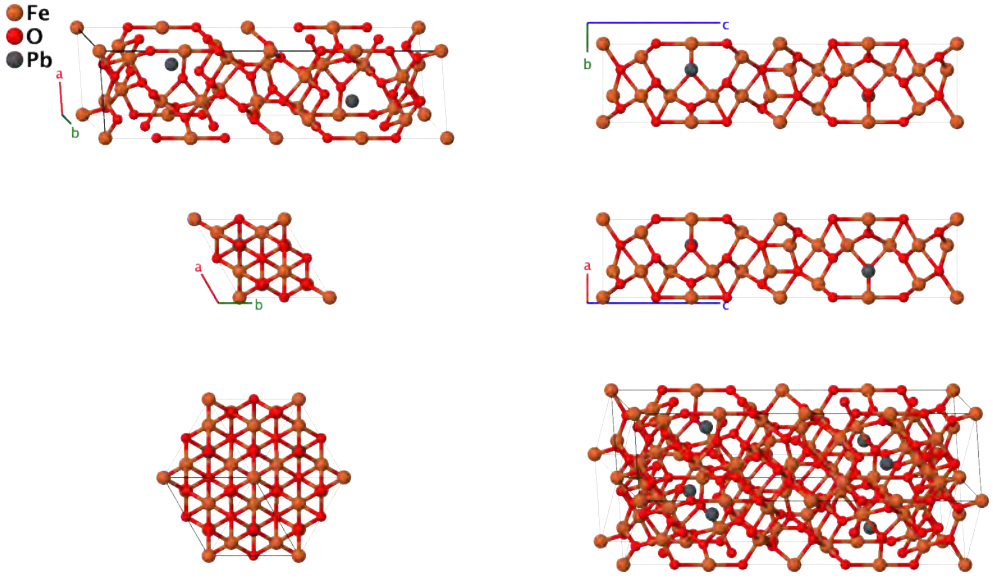
- H. Xie, Y. Yao, X. Feng, D. Duan, H. Song, Z. Zhang, S. Jiang, S. A. T. Redfern, V. Z. Kresin, C. J. Pickard, and T. Cui, *Hydrogen pentagraphenelike structure stabilized by hafnium: a high-temperature conventional superconductor*, Phys. Rev. Lett. **125**, 217001 (2020), doi:10.1103/PhysRevLett.125.217001.

Geometry files:

- CIF: pp. 1763
- POSCAR: pp. 1764

Magnetoplumbite ($PbFe_{12}O_{19}$) Structure: A12B19C_hP64_194_ab2fk_efh2k_d

http://aflow.org/prototype-encyclopedia/A12B19C_hP64_194_ab2fk_efh2k_d



Prototype	:	$Fe_{12}O_{19}Pb$
AFLOW prototype label	:	A12B19C_hP64_194_ab2fk_efh2k_d
***Strukturbericht* designation**	:	None
Pearson symbol	:	hP64
Space group number	:	194
Space group symbol	:	$P6_3/mmc$
AFLOW prototype command	:	`aflow --proto=A12B19C_hP64_194_ab2fk_efh2k_d --params=`$a, c/a, z_4, z_5, z_6, z_7, x_8, x_9, z_9, x_{10}, z_{10}, x_{11}, z_{11}$

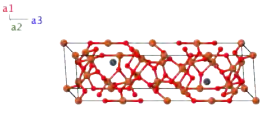
Other compounds with this structure

- $PbAl_{12}O_{19}$, $PbGa_{12}O_{19}$, $PbMn_{12}O_{19}$, $Pb(Co,Ti)_{12}O_{19}$, $BaAl_{12}O_{19}$, $BaGa_{12}O_{19}$, $BaMn_{12}O_{19}$, $Ba(Co,Ti)_{12}O_{19}$, $CaAl_{12}O_{19}$, $CaGa_{12}O_{19}$, $CaMn_{12}O_{19}$, $Ca(Co,Ti)_{12}O_{19}$, $SrAl_{12}O_{19}$, $SrGa_{12}O_{19}$, $SrMn_{12}O_{19}$, and $Sr(Co,Ti)_{12}O_{19}$

- In addition to the listed compounds, the lead and iron sites may be alloyed with a wide variety of metals and semi-metals resulting in high-entropy phases (Vinnik, 2019).

Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(2*a*)	Fe I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	(2*a*)	Fe I
$\mathbf{B}_3$	=	$\frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\,\hat{\mathbf{z}}$	(2*b*)	Fe II
$\mathbf{B}_4$	=	$\frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\,\hat{\mathbf{z}}$	(2*b*)	Fe II
$\mathbf{B}_5$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(2*d*)	Pb
$\mathbf{B}_6$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(2*d*)	Pb
$\mathbf{B}_7$	=	$z_4\,\mathbf{a}_3$	=	$z_4 c\,\hat{\mathbf{z}}$	(4*e*)	O I
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + z_4\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	(4*e*)	O I
$\mathbf{B}_9$	=	$-z_4\,\mathbf{a}_3$	=	$-z_4 c\,\hat{\mathbf{z}}$	(4*e*)	O I
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - z_4\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	(4*e*)	O I
$\mathbf{B}_{11}$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	(4*f*)	Fe III
$\mathbf{B}_{12}$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	(4*f*)	Fe III
$\mathbf{B}_{13}$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	(4*f*)	Fe III
$\mathbf{B}_{14}$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	(4*f*)	Fe III
$\mathbf{B}_{15}$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	(4*f*)	Fe IV
$\mathbf{B}_{16}$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\,\hat{\mathbf{z}}$	(4*f*)	Fe IV
$\mathbf{B}_{17}$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	(4*f*)	Fe IV
$\mathbf{B}_{18}$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\,\hat{\mathbf{z}}$	(4*f*)	Fe IV
$\mathbf{B}_{19}$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	(4*f*)	O II
$\mathbf{B}_{20}$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right)c\,\hat{\mathbf{z}}$	(4*f*)	O II
$\mathbf{B}_{21}$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	(4*f*)	O II
$\mathbf{B}_{22}$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right)c\,\hat{\mathbf{z}}$	(4*f*)	O II
$\mathbf{B}_{23}$	=	$x_8\,\mathbf{a}_1 + 2x_8\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{3}{2}x_8 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_8 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	O III
$\mathbf{B}_{24}$	=	$-2x_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-\frac{3}{2}x_8 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_8 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	O III
$\mathbf{B}_{25}$	=	$x_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-\sqrt{3}x_8 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	(6*h*)	O III
$\mathbf{B}_{26}$	=	$-x_8\,\mathbf{a}_1 - 2x_8\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-\frac{3}{2}x_8 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_8 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	O III
$\mathbf{B}_{27}$	=	$2x_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{2}x_8 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_8 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	O III
$\mathbf{B}_{28}$	=	$-x_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\sqrt{3}x_8 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	(6*h*)	O III
$\mathbf{B}_{29}$	=	$x_9\,\mathbf{a}_1 + 2x_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$\frac{3}{2}x_9 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_9 a\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(12*k*)	Fe V
$\mathbf{B}_{30}$	=	$-2x_9\,\mathbf{a}_1 - x_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$-\frac{3}{2}x_9 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_9 a\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(12*k*)	Fe V
$\mathbf{B}_{31}$	=	$x_9\,\mathbf{a}_1 - x_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$-\sqrt{3}x_9 a\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	(12*k*)	Fe V
$\mathbf{B}_{32}$	=	$-x_9\,\mathbf{a}_1 - 2x_9\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\,\mathbf{a}_3$	=	$-\frac{3}{2}x_9 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_9 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right)c\,\hat{\mathbf{z}}$	(12*k*)	Fe V
$\mathbf{B}_{33}$	=	$2x_9\,\mathbf{a}_1 + x_9\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\,\mathbf{a}_3$	=	$\frac{3}{2}x_9 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_9 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right)c\,\hat{\mathbf{z}}$	(12*k*)	Fe V
$\mathbf{B}_{34}$	=	$-x_9\,\mathbf{a}_1 + x_9\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\,\mathbf{a}_3$	=	$\sqrt{3}x_9 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right)c\,\hat{\mathbf{z}}$	(12*k*)	Fe V
$\mathbf{B}_{35}$	=	$2x_9\,\mathbf{a}_1 + x_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$\frac{3}{2}x_9 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_9 a\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}}$	(12*k*)	Fe V
$\mathbf{B}_{36}$	=	$-x_9\,\mathbf{a}_1 - 2x_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$-\frac{3}{2}x_9 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_9 a\,\hat{\mathbf{y}} - z_9 c\,\hat{\mathbf{z}}$	(12*k*)	Fe V

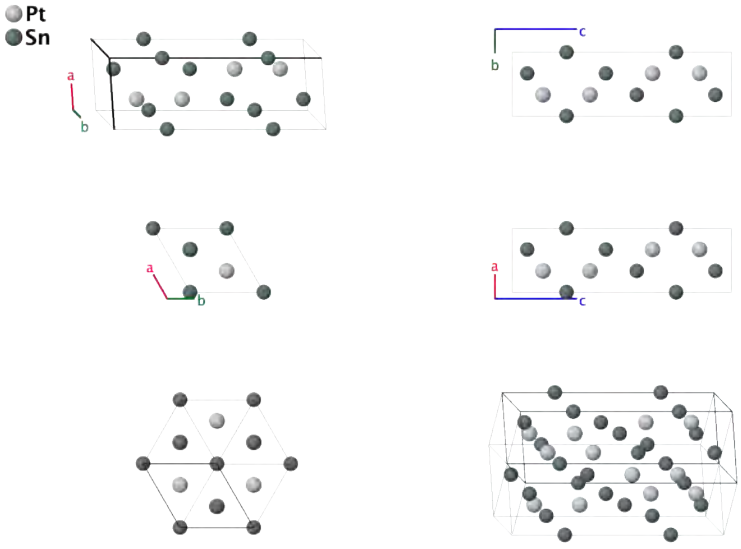
$\mathbf{B}_{37}$	=	$-x_9\,\mathbf{a}_1 + x_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$\sqrt{3}x_9a\,\hat{\mathbf{y}} - z_9c\,\hat{\mathbf{z}}$	$(12k)$	Fe V
$\mathbf{B}_{38}$	=	$-2x_9\,\mathbf{a}_1 - x_9\,\mathbf{a}_2 + \left(\frac{1}{2} - z_9\right)\mathbf{a}_3$	=	$-\frac{3}{2}x_9a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_9a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_9\right)c\,\hat{\mathbf{z}}$	$(12k)$	Fe V
$\mathbf{B}_{39}$	=	$x_9\,\mathbf{a}_1 + 2x_9\,\mathbf{a}_2 + \left(\frac{1}{2} - z_9\right)\mathbf{a}_3$	=	$\frac{3}{2}x_9a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_9a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_9\right)c\,\hat{\mathbf{z}}$	$(12k)$	Fe V
$\mathbf{B}_{40}$	=	$x_9\,\mathbf{a}_1 - x_9\,\mathbf{a}_2 + \left(\frac{1}{2} - z_9\right)\mathbf{a}_3$	=	$-\sqrt{3}x_9a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_9\right)c\,\hat{\mathbf{z}}$	$(12k)$	Fe V
$\mathbf{B}_{41}$	=	$x_{10}\,\mathbf{a}_1 + 2x_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$\frac{3}{2}x_{10}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_{10}a\,\hat{\mathbf{y}} + z_{10}c\,\hat{\mathbf{z}}$	$(12k)$	O IV
$\mathbf{B}_{42}$	=	$-2x_{10}\,\mathbf{a}_1 - x_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$-\frac{3}{2}x_{10}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_{10}a\,\hat{\mathbf{y}} + z_{10}c\,\hat{\mathbf{z}}$	$(12k)$	O IV
$\mathbf{B}_{43}$	=	$x_{10}\,\mathbf{a}_1 - x_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$-\sqrt{3}x_{10}a\,\hat{\mathbf{y}} + z_{10}c\,\hat{\mathbf{z}}$	$(12k)$	O IV
$\mathbf{B}_{44}$	=	$-x_{10}\,\mathbf{a}_1 - 2x_{10}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{10}\right)\mathbf{a}_3$	=	$-\frac{3}{2}x_{10}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_{10}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{10}\right)c\,\hat{\mathbf{z}}$	$(12k)$	O IV
$\mathbf{B}_{45}$	=	$2x_{10}\,\mathbf{a}_1 + x_{10}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{10}\right)\mathbf{a}_3$	=	$\frac{3}{2}x_{10}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_{10}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{10}\right)c\,\hat{\mathbf{z}}$	$(12k)$	O IV
$\mathbf{B}_{46}$	=	$-x_{10}\,\mathbf{a}_1 + x_{10}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{10}\right)\mathbf{a}_3$	=	$\sqrt{3}x_{10}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{10}\right)c\,\hat{\mathbf{z}}$	$(12k)$	O IV
$\mathbf{B}_{47}$	=	$2x_{10}\,\mathbf{a}_1 + x_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3$	=	$\frac{3}{2}x_{10}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_{10}a\,\hat{\mathbf{y}} - z_{10}c\,\hat{\mathbf{z}}$	$(12k)$	O IV
$\mathbf{B}_{48}$	=	$-x_{10}\,\mathbf{a}_1 - 2x_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3$	=	$-\frac{3}{2}x_{10}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_{10}a\,\hat{\mathbf{y}} - z_{10}c\,\hat{\mathbf{z}}$	$(12k)$	O IV
$\mathbf{B}_{49}$	=	$-x_{10}\,\mathbf{a}_1 + x_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3$	=	$\sqrt{3}x_{10}a\,\hat{\mathbf{y}} - z_{10}c\,\hat{\mathbf{z}}$	$(12k)$	O IV
$\mathbf{B}_{50}$	=	$-2x_{10}\,\mathbf{a}_1 - x_{10}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{10}\right)\mathbf{a}_3$	=	$-\frac{3}{2}x_{10}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_{10}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{10}\right)c\,\hat{\mathbf{z}}$	$(12k)$	O IV
$\mathbf{B}_{51}$	=	$x_{10}\,\mathbf{a}_1 + 2x_{10}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{10}\right)\mathbf{a}_3$	=	$\frac{3}{2}x_{10}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_{10}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{10}\right)c\,\hat{\mathbf{z}}$	$(12k)$	O IV
$\mathbf{B}_{52}$	=	$x_{10}\,\mathbf{a}_1 - x_{10}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{10}\right)\mathbf{a}_3$	=	$-\sqrt{3}x_{10}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{10}\right)c\,\hat{\mathbf{z}}$	$(12k)$	O IV
$\mathbf{B}_{53}$	=	$x_{11}\,\mathbf{a}_1 + 2x_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$\frac{3}{2}x_{11}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_{11}a\,\hat{\mathbf{y}} + z_{11}c\,\hat{\mathbf{z}}$	$(12k)$	O V
$\mathbf{B}_{54}$	=	$-2x_{11}\,\mathbf{a}_1 - x_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$-\frac{3}{2}x_{11}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_{11}a\,\hat{\mathbf{y}} + z_{11}c\,\hat{\mathbf{z}}$	$(12k)$	O V
$\mathbf{B}_{55}$	=	$x_{11}\,\mathbf{a}_1 - x_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$-\sqrt{3}x_{11}a\,\hat{\mathbf{y}} + z_{11}c\,\hat{\mathbf{z}}$	$(12k)$	O V
$\mathbf{B}_{56}$	=	$-x_{11}\,\mathbf{a}_1 - 2x_{11}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{11}\right)\mathbf{a}_3$	=	$-\frac{3}{2}x_{11}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_{11}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{11}\right)c\,\hat{\mathbf{z}}$	$(12k)$	O V
$\mathbf{B}_{57}$	=	$2x_{11}\,\mathbf{a}_1 + x_{11}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{11}\right)\mathbf{a}_3$	=	$\frac{3}{2}x_{11}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_{11}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{11}\right)c\,\hat{\mathbf{z}}$	$(12k)$	O V
$\mathbf{B}_{58}$	=	$-x_{11}\,\mathbf{a}_1 + x_{11}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{11}\right)\mathbf{a}_3$	=	$\sqrt{3}x_{11}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{11}\right)c\,\hat{\mathbf{z}}$	$(12k)$	O V
$\mathbf{B}_{59}$	=	$2x_{11}\,\mathbf{a}_1 + x_{11}\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3$	=	$\frac{3}{2}x_{11}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_{11}a\,\hat{\mathbf{y}} - z_{11}c\,\hat{\mathbf{z}}$	$(12k)$	O V
$\mathbf{B}_{60}$	=	$-x_{11}\,\mathbf{a}_1 - 2x_{11}\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3$	=	$-\frac{3}{2}x_{11}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_{11}a\,\hat{\mathbf{y}} - z_{11}c\,\hat{\mathbf{z}}$	$(12k)$	O V
$\mathbf{B}_{61}$	=	$-x_{11}\,\mathbf{a}_1 + x_{11}\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3$	=	$\sqrt{3}x_{11}a\,\hat{\mathbf{y}} - z_{11}c\,\hat{\mathbf{z}}$	$(12k)$	O V
$\mathbf{B}_{62}$	=	$-2x_{11}\,\mathbf{a}_1 - x_{11}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{11}\right)\mathbf{a}_3$	=	$-\frac{3}{2}x_{11}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_{11}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{11}\right)c\,\hat{\mathbf{z}}$	$(12k)$	O V
$\mathbf{B}_{63}$	=	$x_{11}\,\mathbf{a}_1 + 2x_{11}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{11}\right)\mathbf{a}_3$	=	$\frac{3}{2}x_{11}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_{11}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{11}\right)c\,\hat{\mathbf{z}}$	$(12k)$	O V
$\mathbf{B}_{64}$	=	$x_{11}\,\mathbf{a}_1 - x_{11}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{11}\right)\mathbf{a}_3$	=	$-\sqrt{3}x_{11}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{11}\right)c\,\hat{\mathbf{z}}$	$(12k)$	O V

References:
- R. Gerber, Z. Šimša, and L. Jenšovský, *A note on the magnetoplumbite crystal structure*, Czech. J. Phys. **44**, 937–940 (1994), doi:10.1007/BF01715487.
- D. A. Vinnik, E. A. Trofimov, V. E. Zhivulin, O. V. Zaitseva, S. A. Gudkova, A. Y. Starikov, D. A. Zherebtsov, A. A. Kirsanova, M. Häßner, and R. Niewa, *High-entropy oxide phases with magnetoplumbite structure*, Ceram. Int. **45**, 12942–12948 (2019), doi:10.1016/j.ceramint.2019.03.221.

Geometry files:
- CIF: pp. 1764
- POSCAR: pp. 1764

Pt_2Sn_3 ($D5_b$) Structure: A2B3_hP10_194_f_bf

http://aflow.org/prototype-encyclopedia/A2B3_hP10_194_f_bf



Prototype	:	Pt_2Sn_3
AFLOW prototype label	:	A2B3_hP10_194_f_bf
***Strukturbericht* designation**	:	$D5_b$
Pearson symbol	:	hP10
Space group number	:	194
Space group symbol	:	$P6_3/mmc$
AFLOW prototype command	:	`aflow --proto=A2B3_hP10_194_f_bf --params=`$a, c/a, z_2, z_3$

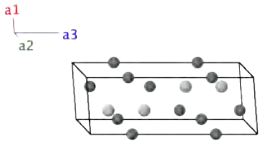
Other compounds with this structure

- Pt_2Si_3, $InAu_2Ga_2$, and $Au_4In_3Sn_3$

- We were unable to obtain a copy of (Schubert, 1949), and so we use the structural parameters given by (Pearson, 1958), using (Wood, 1947) to convert the kX units used by Pearson into Ångström units.

Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{4}\mathbf{a}_3$	=	$\frac{1}{4}c\,\hat{\mathbf{z}}$	($2b$)	Sn I
$\mathbf{B}_2$	=	$\frac{3}{4}\mathbf{a}_3$	=	$\frac{3}{4}c\,\hat{\mathbf{z}}$	($2b$)	Sn I

$\mathbf{B}_3$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(4f)$	Pt
$\mathbf{B}_4$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	$(4f)$	Pt
$\mathbf{B}_5$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	$(4f)$	Pt
$\mathbf{B}_6$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	$(4f)$	Pt
$\mathbf{B}_7$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(4f)$	Sn II
$\mathbf{B}_8$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	$(4f)$	Sn II
$\mathbf{B}_9$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(4f)$	Sn II
$\mathbf{B}_{10}$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	$(4f)$	Sn II

References:

- K. Schubert and H. Pfisterer, *Kristallstruktur von Pt_2Sn_3*, Z. Metallkd. **40**, 405 (1949).
- E. A. Wood, *The Conversion Factor for kX Units to Angström Units*, J. Appl. Phys. **18**, 929 (1947), doi:10.1063/1.1697570.

Found in:

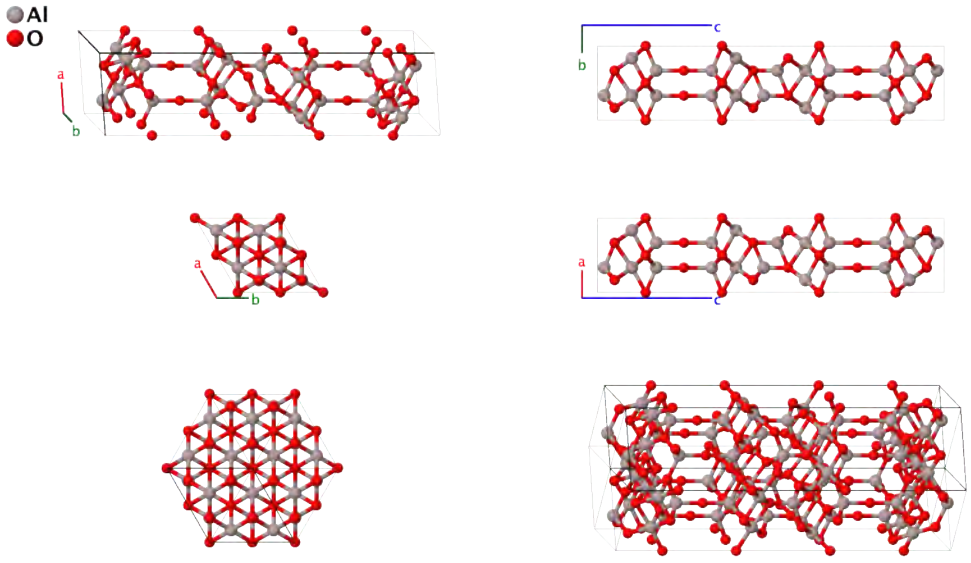
- W. B. Pearson, *A Handbook of Lattice Spacings and Structures of Metals and Alloys*, *International Series of Monographs on Metal Physics and Physical Metallurgy*, vol. 4 (Pergamon Press, Oxford, London, Edinburgh, New York, Paris, Frankfort, 1958), 1964 reprint with corrections edn. N. R. C. No. 4303.

Geometry files:

- CIF: pp. 1765
- POSCAR: pp. 1765

β-Alumina (Al_2O_3, $D5_6$) Structure: A2B3_hP60_194_3fk_cdef2k

http://aflow.org/prototype-encyclopedia/A2B3_hP60_194_3fk_cdef2k

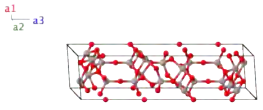


Prototype	:	Al_2O_3
AFLOW prototype label	:	A2B3_hP60_194_3fk_cdef2k
***Strukturbericht* designation**	:	$D5_6$
Pearson symbol	:	hP60
Space group number	:	194
Space group symbol	:	$P6_3/mmc$
AFLOW prototype command	:	`aflow --proto=A2B3_hP60_194_3fk_cdef2k --params=`$a, c/a, z_3, z_4, z_5, z_6, z_7, x_8, z_8, x_9, z_9, x_{10}, z_{10}$

- (Hermann, 1937) assigned this the *Strukturbericht* designation $D5_6$, calling it β-corundum, and subtitles the section "with small Na_2O impurities." As noted by (Gottfried, 1937) and (Le Cars, 1975), the Na impurities replace Al atoms on the ($4f$) Wyckoff positions, along with some unspecified O sites. Charge neutrality requires that some of the displaced atoms remain in the lattice, and they are moved to ($2a$) Wyckoff positions at (0, 0, 0) and (0, 0, 1/2). Here we only list the "ideal" structure.

Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$ =	$\frac{1}{3}\mathbf{a}_1 + \frac{2}{3}\mathbf{a}_2 + \frac{1}{4}\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($2c$)	O I

$\mathbf{B}_2$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(2c)$	O I
$\mathbf{B}_3$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(2d)$	O II
$\mathbf{B}_4$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(2d)$	O II
$\mathbf{B}_5$	=	$z_3\,\mathbf{a}_3$	=	$z_3 c\,\hat{\mathbf{z}}$	$(4e)$	O III
$\mathbf{B}_6$	=	$\left(\frac{1}{2} + z_3\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	$(4e)$	O III
$\mathbf{B}_7$	=	$-z_3\,\mathbf{a}_3$	=	$-z_3 c\,\hat{\mathbf{z}}$	$(4e)$	O III
$\mathbf{B}_8$	=	$\left(\frac{1}{2} - z_3\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	$(4e)$	O III
$\mathbf{B}_9$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(4f)$	Al I
$\mathbf{B}_{10}$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	$(4f)$	Al I
$\mathbf{B}_{11}$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	$(4f)$	Al I
$\mathbf{B}_{12}$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	$(4f)$	Al I
$\mathbf{B}_{13}$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(4f)$	Al II
$\mathbf{B}_{14}$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	$(4f)$	Al II
$\mathbf{B}_{15}$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	$(4f)$	Al II
$\mathbf{B}_{16}$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	$(4f)$	Al II
$\mathbf{B}_{17}$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(4f)$	Al III
$\mathbf{B}_{18}$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\,\hat{\mathbf{z}}$	$(4f)$	Al III
$\mathbf{B}_{19}$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_6 c\,\hat{\mathbf{z}}$	$(4f)$	Al III
$\mathbf{B}_{20}$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\,\hat{\mathbf{z}}$	$(4f)$	Al III
$\mathbf{B}_{21}$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_7\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_7 c\,\hat{\mathbf{z}}$	$(4f)$	O IV
$\mathbf{B}_{22}$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_7\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_7\right)c\,\hat{\mathbf{z}}$	$(4f)$	O IV
$\mathbf{B}_{23}$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_7 c\,\hat{\mathbf{z}}$	$(4f)$	O IV
$\mathbf{B}_{24}$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_7\right)\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_7\right)c\,\hat{\mathbf{z}}$	$(4f)$	O IV
$\mathbf{B}_{25}$	=	$x_8\,\mathbf{a}_1 + 2x_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$\frac{3}{2}x_8 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_8 a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	$(12k)$	Al IV
$\mathbf{B}_{26}$	=	$-2x_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$-\frac{3}{2}x_8 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_8 a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	$(12k)$	Al IV
$\mathbf{B}_{27}$	=	$x_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$-\sqrt{3}x_8 a\,\hat{\mathbf{y}} + z_8 c\,\hat{\mathbf{z}}$	$(12k)$	Al IV
$\mathbf{B}_{28}$	=	$-x_8\,\mathbf{a}_1 - 2x_8\,\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\,\mathbf{a}_3$	=	$-\frac{3}{2}x_8 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_8 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right)c\,\hat{\mathbf{z}}$	$(12k)$	Al IV
$\mathbf{B}_{29}$	=	$2x_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\,\mathbf{a}_3$	=	$\frac{3}{2}x_8 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_8 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right)c\,\hat{\mathbf{z}}$	$(12k)$	Al IV
$\mathbf{B}_{30}$	=	$-x_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\,\mathbf{a}_3$	=	$\sqrt{3}x_8 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right)c\,\hat{\mathbf{z}}$	$(12k)$	Al IV
$\mathbf{B}_{31}$	=	$2x_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$\frac{3}{2}x_8 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_8 a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	$(12k)$	Al IV
$\mathbf{B}_{32}$	=	$-x_8\,\mathbf{a}_1 - 2x_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$-\frac{3}{2}x_8 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_8 a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	$(12k)$	Al IV
$\mathbf{B}_{33}$	=	$-x_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$\sqrt{3}x_8 a\,\hat{\mathbf{y}} - z_8 c\,\hat{\mathbf{z}}$	$(12k)$	Al IV
$\mathbf{B}_{34}$	=	$-2x_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\,\mathbf{a}_3$	=	$-\frac{3}{2}x_8 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_8 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right)c\,\hat{\mathbf{z}}$	$(12k)$	Al IV
$\mathbf{B}_{35}$	=	$x_8\,\mathbf{a}_1 + 2x_8\,\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\,\mathbf{a}_3$	=	$\frac{3}{2}x_8 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_8 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right)c\,\hat{\mathbf{z}}$	$(12k)$	Al IV
$\mathbf{B}_{36}$	=	$x_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 + \left(\frac{1}{2} - z_8\right)\,\mathbf{a}_3$	=	$-\sqrt{3}x_8 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_8\right)c\,\hat{\mathbf{z}}$	$(12k)$	Al IV
$\mathbf{B}_{37}$	=	$x_9\,\mathbf{a}_1 + 2x_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$\frac{3}{2}x_9 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_9 a\,\hat{\mathbf{y}} + z_9 c\,\hat{\mathbf{z}}$	$(12k)$	O V

$\mathbf{B}_{38}$	=	$-2x_9\,\mathbf{a}_1 - x_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$-\frac{3}{2}x_9a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_9a\,\hat{\mathbf{y}} + z_9c\,\hat{\mathbf{z}}$	(12*k*)	O V
$\mathbf{B}_{39}$	=	$x_9\,\mathbf{a}_1 - x_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$-\sqrt{3}x_9a\,\hat{\mathbf{y}} + z_9c\,\hat{\mathbf{z}}$	(12*k*)	O V
$\mathbf{B}_{40}$	=	$-x_9\,\mathbf{a}_1 - 2x_9\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$-\frac{3}{2}x_9a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_9a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right)c\,\hat{\mathbf{z}}$	(12*k*)	O V
$\mathbf{B}_{41}$	=	$2x_9\,\mathbf{a}_1 + x_9\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$\frac{3}{2}x_9a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_9a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right)c\,\hat{\mathbf{z}}$	(12*k*)	O V
$\mathbf{B}_{42}$	=	$-x_9\,\mathbf{a}_1 + x_9\,\mathbf{a}_2 + \left(\frac{1}{2} + z_9\right)\mathbf{a}_3$	=	$\sqrt{3}x_9a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_9\right)c\,\hat{\mathbf{z}}$	(12*k*)	O V
$\mathbf{B}_{43}$	=	$2x_9\,\mathbf{a}_1 + x_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$\frac{3}{2}x_9a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_9a\,\hat{\mathbf{y}} - z_9c\,\hat{\mathbf{z}}$	(12*k*)	O V
$\mathbf{B}_{44}$	=	$-x_9\,\mathbf{a}_1 - 2x_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$-\frac{3}{2}x_9a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_9a\,\hat{\mathbf{y}} - z_9c\,\hat{\mathbf{z}}$	(12*k*)	O V
$\mathbf{B}_{45}$	=	$-x_9\,\mathbf{a}_1 + x_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3$	=	$\sqrt{3}x_9a\,\hat{\mathbf{y}} - z_9c\,\hat{\mathbf{z}}$	(12*k*)	O V
$\mathbf{B}_{46}$	=	$-2x_9\,\mathbf{a}_1 - x_9\,\mathbf{a}_2 + \left(\frac{1}{2} - z_9\right)\mathbf{a}_3$	=	$-\frac{3}{2}x_9a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_9a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_9\right)c\,\hat{\mathbf{z}}$	(12*k*)	O V
$\mathbf{B}_{47}$	=	$x_9\,\mathbf{a}_1 + 2x_9\,\mathbf{a}_2 + \left(\frac{1}{2} - z_9\right)\mathbf{a}_3$	=	$\frac{3}{2}x_9a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_9a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_9\right)c\,\hat{\mathbf{z}}$	(12*k*)	O V
$\mathbf{B}_{48}$	=	$x_9\,\mathbf{a}_1 - x_9\,\mathbf{a}_2 + \left(\frac{1}{2} - z_9\right)\mathbf{a}_3$	=	$-\sqrt{3}x_9a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_9\right)c\,\hat{\mathbf{z}}$	(12*k*)	O V
$\mathbf{B}_{49}$	=	$x_{10}\,\mathbf{a}_1 + 2x_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$\frac{3}{2}x_{10}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_{10}a\,\hat{\mathbf{y}} + z_{10}c\,\hat{\mathbf{z}}$	(12*k*)	O VI
$\mathbf{B}_{50}$	=	$-2x_{10}\,\mathbf{a}_1 - x_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$-\frac{3}{2}x_{10}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_{10}a\,\hat{\mathbf{y}} + z_{10}c\,\hat{\mathbf{z}}$	(12*k*)	O VI
$\mathbf{B}_{51}$	=	$x_{10}\,\mathbf{a}_1 - x_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$-\sqrt{3}x_{10}a\,\hat{\mathbf{y}} + z_{10}c\,\hat{\mathbf{z}}$	(12*k*)	O VI
$\mathbf{B}_{52}$	=	$-x_{10}\,\mathbf{a}_1 - 2x_{10}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{10}\right)\mathbf{a}_3$	=	$-\frac{3}{2}x_{10}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_{10}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{10}\right)c\,\hat{\mathbf{z}}$	(12*k*)	O VI
$\mathbf{B}_{53}$	=	$2x_{10}\,\mathbf{a}_1 + x_{10}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{10}\right)\mathbf{a}_3$	=	$\frac{3}{2}x_{10}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_{10}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{10}\right)c\,\hat{\mathbf{z}}$	(12*k*)	O VI
$\mathbf{B}_{54}$	=	$-x_{10}\,\mathbf{a}_1 + x_{10}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{10}\right)\mathbf{a}_3$	=	$\sqrt{3}x_{10}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{10}\right)c\,\hat{\mathbf{z}}$	(12*k*)	O VI
$\mathbf{B}_{55}$	=	$2x_{10}\,\mathbf{a}_1 + x_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3$	=	$\frac{3}{2}x_{10}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_{10}a\,\hat{\mathbf{y}} - z_{10}c\,\hat{\mathbf{z}}$	(12*k*)	O VI
$\mathbf{B}_{56}$	=	$-x_{10}\,\mathbf{a}_1 - 2x_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3$	=	$-\frac{3}{2}x_{10}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_{10}a\,\hat{\mathbf{y}} - z_{10}c\,\hat{\mathbf{z}}$	(12*k*)	O VI
$\mathbf{B}_{57}$	=	$-x_{10}\,\mathbf{a}_1 + x_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3$	=	$\sqrt{3}x_{10}a\,\hat{\mathbf{y}} - z_{10}c\,\hat{\mathbf{z}}$	(12*k*)	O VI
$\mathbf{B}_{58}$	=	$-2x_{10}\,\mathbf{a}_1 - x_{10}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{10}\right)\mathbf{a}_3$	=	$-\frac{3}{2}x_{10}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_{10}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{10}\right)c\,\hat{\mathbf{z}}$	(12*k*)	O VI
$\mathbf{B}_{59}$	=	$x_{10}\,\mathbf{a}_1 + 2x_{10}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{10}\right)\mathbf{a}_3$	=	$\frac{3}{2}x_{10}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_{10}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{10}\right)c\,\hat{\mathbf{z}}$	(12*k*)	O VI
$\mathbf{B}_{60}$	=	$x_{10}\,\mathbf{a}_1 - x_{10}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{10}\right)\mathbf{a}_3$	=	$-\sqrt{3}x_{10}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{10}\right)c\,\hat{\mathbf{z}}$	(12*k*)	O VI

References:

- W. L. Bragg, C. Gottfried, and J. West, *The Structure of β Alumina*, Zeitschrift für Kristallographie - Crystalline Materials **77**, 255–274 (1931), doi:10.1524/zkri.1931.77.1.255.
- Y. Le Cars, D. Gratias, R. Portier, and J. Théry, *Planar defects in β-alumina*, J. Solid State Chem. **15**, 218–222 (1975), doi:10.1016/0022-4596(75)90205-4.

Found in:

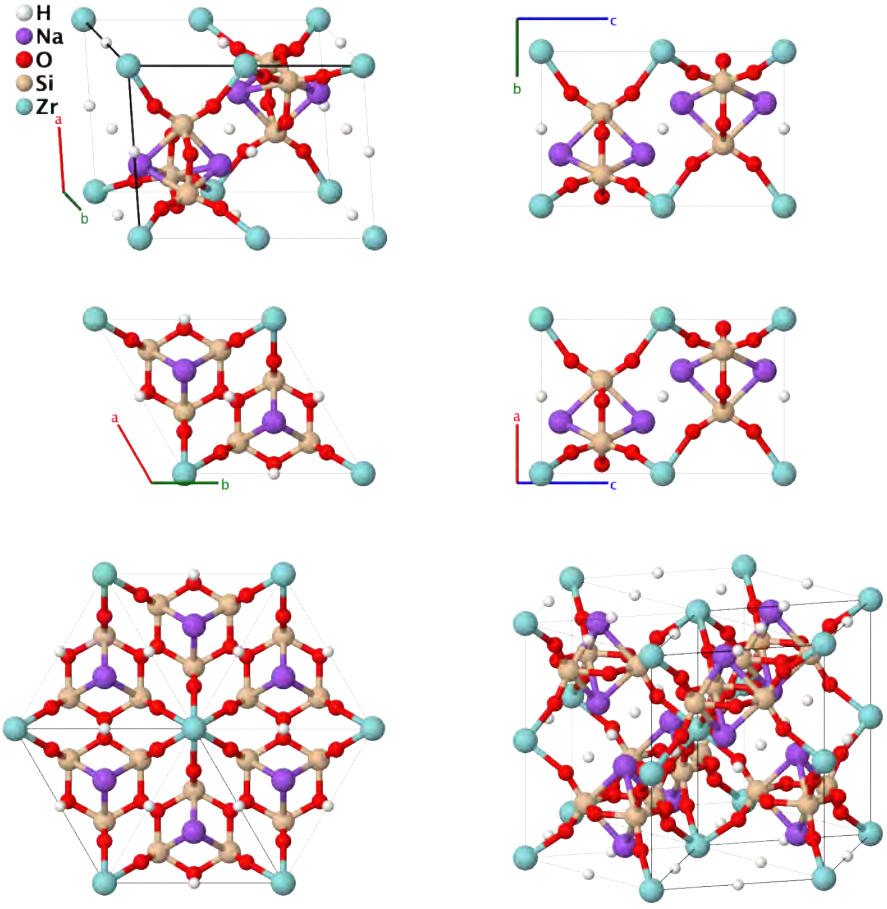
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1765
- POSCAR: pp. 1766

$S\,3_4$ (II) (Catapleiite, $Na_2Zr(SiO_3)_3 \cdot H_2O$) (*obsolete*) Structure: A3B2C9D3E_hP36_194_g_f_hk_h_a

http://aflow.org/prototype-encyclopedia/A3B2C9D3E_hP36_194_g_f_hk_h_a

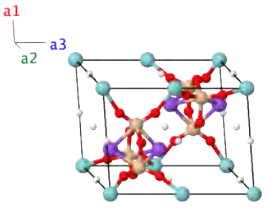


Prototype	:	$(H_2O)Na_2O_9Si_3Zr$
AFLOW prototype label	:	A3B2C9D3E_hP36_194_g_f_hk_h_a
***Strukturbericht* designation**	:	$S\,3_4$ (II)
Pearson symbol	:	hP36
Space group number	:	194
Space group symbol	:	$P6_3/mmc$
AFLOW prototype command	:	`aflow --proto=A3B2C9D3E_hP36_194_g_f_hk_h_a --params=`$a, c/a, z_2, x_4, x_5, x_6, z_6$

- This hexagonal structure has been superseded by the monoclinic structure of (Ilyushin, 1981). We present it here for historical interest.
- We were unable to procure the original reference, so we use the data provided by (Gottfried, 1937).
- Four of the (6*g*) sites are randomly occupied by water molecules.
- (Gottfried, 1937) originally gave the *Strukturbericht* designation $S\,3_4$ to chabazite, but (Gottfried, 1940) gave it to this structure of catapleiite. We resolve this dilemma by using $S\,3_4$(I) for chabazite and $S\,3_4$(II) for catapleiite.

Hexagonal primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	Zr
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	$(2a)$	Zr
$\mathbf{B}_3$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(4f)$	Na
$\mathbf{B}_4$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	$(4f)$	Na
$\mathbf{B}_5$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	$(4f)$	Na
$\mathbf{B}_6$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	$(4f)$	Na
$\mathbf{B}_7$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{4}a\,\hat{\mathbf{y}}$	$(6g)$	H_2O
$\mathbf{B}_8$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{4}a\,\hat{\mathbf{y}}$	$(6g)$	H_2O
$\mathbf{B}_9$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	$(6g)$	H_2O
$\mathbf{B}_{10}$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{4}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(6g)$	H_2O
$\mathbf{B}_{11}$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{4}a\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(6g)$	H_2O
$\mathbf{B}_{12}$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	$(6g)$	H_2O
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + 2x_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{3}{2}x_4 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O I
$\mathbf{B}_{14}$	=	$-2x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-\frac{3}{2}x_4 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O I
$\mathbf{B}_{15}$	=	$x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-\sqrt{3}x_4 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O I
$\mathbf{B}_{16}$	=	$-x_4\,\mathbf{a}_1 - 2x_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-\frac{3}{2}x_4 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O I
$\mathbf{B}_{17}$	=	$2x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{2}x_4 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O I
$\mathbf{B}_{18}$	=	$-x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\sqrt{3}x_4 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	O I
$\mathbf{B}_{19}$	=	$x_5\,\mathbf{a}_1 + 2x_5\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{3}{2}x_5 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	Si
$\mathbf{B}_{20}$	=	$-2x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-\frac{3}{2}x_5 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	Si
$\mathbf{B}_{21}$	=	$x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-\sqrt{3}x_5 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	Si
$\mathbf{B}_{22}$	=	$-x_5\,\mathbf{a}_1 - 2x_5\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-\frac{3}{2}x_5 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	Si
$\mathbf{B}_{23}$	=	$2x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{2}x_5 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	Si
$\mathbf{B}_{24}$	=	$-x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\sqrt{3}x_5 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	Si
$\mathbf{B}_{25}$	=	$x_6\,\mathbf{a}_1 + 2x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$\frac{3}{2}x_6 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(12k)$	O II
$\mathbf{B}_{26}$	=	$-2x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$-\frac{3}{2}x_6 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(12k)$	O II
$\mathbf{B}_{27}$	=	$x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	=	$-\sqrt{3}x_6 a\,\hat{\mathbf{y}} + z_6 c\,\hat{\mathbf{z}}$	$(12k)$	O II
$\mathbf{B}_{28}$	=	$-x_6\,\mathbf{a}_1 - 2x_6\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$-\frac{3}{2}x_6 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_6 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\,\hat{\mathbf{z}}$	$(12k)$	O II

$\mathbf{B}_{29}$	=	$2x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\frac{3}{2}x_6a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_6a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\,\hat{\mathbf{z}}$	(12*k*)	O II
$\mathbf{B}_{30}$	=	$-x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	=	$\sqrt{3}x_6a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)c\,\hat{\mathbf{z}}$	(12*k*)	O II
$\mathbf{B}_{31}$	=	$2x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$\frac{3}{2}x_6a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_6a\,\hat{\mathbf{y}} - z_6c\,\hat{\mathbf{z}}$	(12*k*)	O II
$\mathbf{B}_{32}$	=	$-x_6\,\mathbf{a}_1 - 2x_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$-\frac{3}{2}x_6a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_6a\,\hat{\mathbf{y}} - z_6c\,\hat{\mathbf{z}}$	(12*k*)	O II
$\mathbf{B}_{33}$	=	$-x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3$	=	$\sqrt{3}x_6a\,\hat{\mathbf{y}} - z_6c\,\hat{\mathbf{z}}$	(12*k*)	O II
$\mathbf{B}_{34}$	=	$-2x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$-\frac{3}{2}x_6a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_6a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\,\hat{\mathbf{z}}$	(12*k*)	O II
$\mathbf{B}_{35}$	=	$x_6\,\mathbf{a}_1 + 2x_6\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$\frac{3}{2}x_6a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_6a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\,\hat{\mathbf{z}}$	(12*k*)	O II
$\mathbf{B}_{36}$	=	$x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	=	$-\sqrt{3}x_6a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)c\,\hat{\mathbf{z}}$	(12*k*)	O II

References:

- B. Brunowsky, *Die Struktur des Katapleits*, Acta Physicochim. USSR **5**, 863–892 (1936).
- C. Gottfried and F. Schossberger, eds., *Strukturbericht Band III 1933-1935* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Found in:

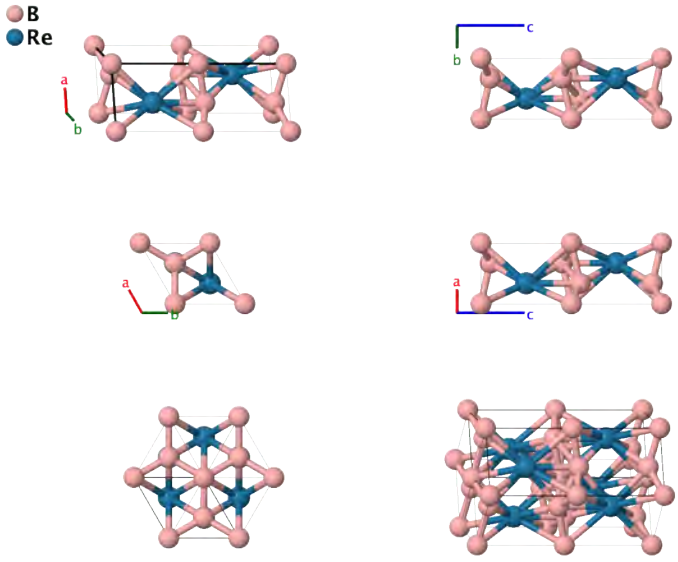
- C. Gottfried, ed., *Strukturbericht Band V 1937* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1940).

Geometry files:

- CIF: pp. 1766
- POSCAR: pp. 1767

ReB_3 Structure: A3B_hP8_194_af_c

http://aflow.org/prototype-encyclopedia/A3B_hP8_194_af_c



Prototype	:	B_3Re
AFLOW prototype label	:	A3B_hP8_194_af_c
***Strukturbericht* designation**	:	None
Pearson symbol	:	hP8
Space group number	:	194
Space group symbol	:	$P6_3/mmc$
AFLOW prototype command	:	`aflow --proto=A3B_hP8_194_af_c --params=`$a, c/a, z_3$

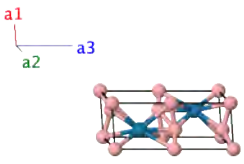
Other compounds with this structure

- TcB_3, $CaNi_2Si$, and $GdPt_2Sn$

- The lattice constants a and c were inferred from the nearest-neighbor distances in (Aronsson, 1960).

Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	B I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	$(2a)$	B I

$\mathbf{B}_3$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(2c)$	Re
$\mathbf{B}_4$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(2c)$	Re
$\mathbf{B}_5$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(4f)$	B II
$\mathbf{B}_6$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	$(4f)$	B II
$\mathbf{B}_7$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(4f)$	B II
$\mathbf{B}_8$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	$(4f)$	B II

References:

- B. Aronsson, E. Stenberg, and J. Åselius, *Borides of Rhenium and the Platinum Metals. The Crystal Structure of Re_7B_3, ReB_3, Rh_7B_3, $RhB_{\sim 1.1}$, $IrB_{\sim 1.1}$ and PtB*, Acta Chem. Scand. **14**, 733–741 (1960), doi:10.3891/acta.chem.scand.14-0733.

Found in:

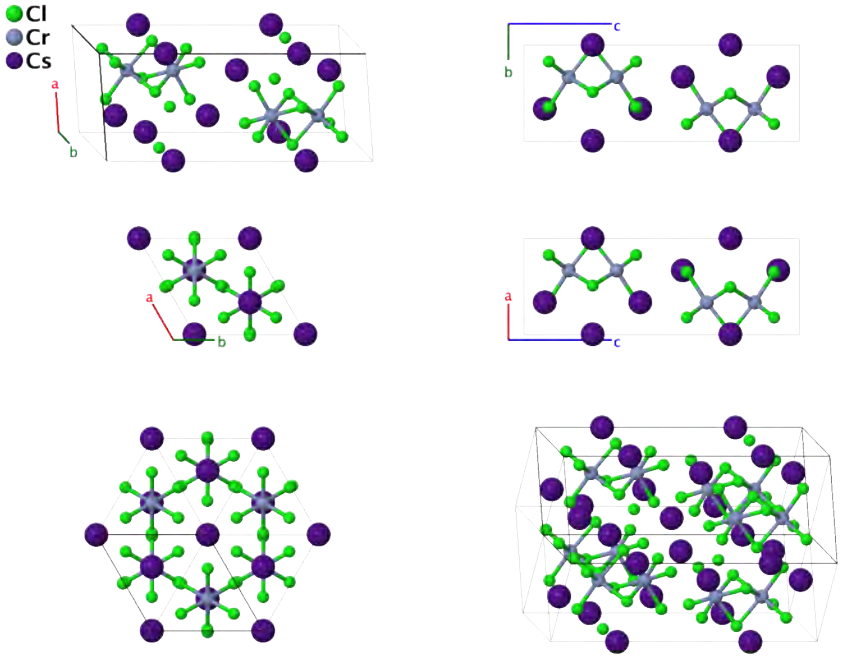
- P. Villars and L. Calvert, *Pearson's Handbook of Crystallographic Data for Intermetallic Phases* (ASM International, Materials Park, OK, 1991), vol. I, p. 612.

Geometry files:

- CIF: pp. 1767
- POSCAR: pp. 1767

$Cs_3Cr_2Cl_9$ Structure: A9B2C3_hP28_194_hk_f_bf

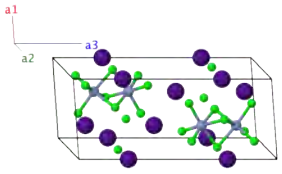
http://aflow.org/prototype-encyclopedia/A9B2C3_hP28_194_hk_f_bf



Prototype	:	$Cl_9Cr_2Cs_3$
AFLOW prototype label	:	A9B2C3_hP28_194_hk_f_bf
***Strukturbericht* designation**	:	None
Pearson symbol	:	hP28
Space group number	:	194
Space group symbol	:	$P6_3/mmc$
AFLOW prototype command	:	`aflow --proto=A9B2C3_hP28_194_hk_f_bf --params=`$a, c/a, z_2, z_3, x_4, x_5, z_5$

Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{4}\mathbf{a}_3$	=	$\frac{1}{4}c\,\hat{\mathbf{z}}$	$(2b)$	Cs I
$\mathbf{B}_2$	=	$\frac{3}{4}\mathbf{a}_3$	=	$\frac{3}{4}c\,\hat{\mathbf{z}}$	$(2b)$	Cs I
$\mathbf{B}_3$	=	$\frac{1}{3}\mathbf{a}_1 + \frac{2}{3}\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(4f)$	Cr
$\mathbf{B}_4$	=	$\frac{2}{3}\mathbf{a}_1 + \frac{1}{3}\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	$(4f)$	Cr

$\mathbf{B}_5$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	$(4f)$	Cr
$\mathbf{B}_6$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	$(4f)$	Cr
$\mathbf{B}_7$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(4f)$	Cs II
$\mathbf{B}_8$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	$(4f)$	Cs II
$\mathbf{B}_9$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(4f)$	Cs II
$\mathbf{B}_{10}$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	$(4f)$	Cs II
$\mathbf{B}_{11}$	=	$x_4\,\mathbf{a}_1 + 2x_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{3}{2}x_4 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	Cl I
$\mathbf{B}_{12}$	=	$-2x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-\frac{3}{2}x_4 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	Cl I
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-\sqrt{3}x_4 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	Cl I
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1 - 2x_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-\frac{3}{2}x_4 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	Cl I
$\mathbf{B}_{15}$	=	$2x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{2}x_4 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	Cl I
$\mathbf{B}_{16}$	=	$-x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\sqrt{3}x_4 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	Cl I
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + 2x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\frac{3}{2}x_5 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(12k)$	Cl II
$\mathbf{B}_{18}$	=	$-2x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$-\frac{3}{2}x_5 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(12k)$	Cl II
$\mathbf{B}_{19}$	=	$x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$-\sqrt{3}x_5 a\,\hat{\mathbf{y}} + z_5 c\,\hat{\mathbf{z}}$	$(12k)$	Cl II
$\mathbf{B}_{20}$	=	$-x_5\,\mathbf{a}_1 - 2x_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$-\frac{3}{2}x_5 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	$(12k)$	Cl II
$\mathbf{B}_{21}$	=	$2x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\frac{3}{2}x_5 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	$(12k)$	Cl II
$\mathbf{B}_{22}$	=	$-x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \left(\frac{1}{2} + z_5\right)\mathbf{a}_3$	=	$\sqrt{3}x_5 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_5\right)c\,\hat{\mathbf{z}}$	$(12k)$	Cl II
$\mathbf{B}_{23}$	=	$2x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\frac{3}{2}x_5 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	$(12k)$	Cl II
$\mathbf{B}_{24}$	=	$-x_5\,\mathbf{a}_1 - 2x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-\frac{3}{2}x_5 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	$(12k)$	Cl II
$\mathbf{B}_{25}$	=	$-x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\sqrt{3}x_5 a\,\hat{\mathbf{y}} - z_5 c\,\hat{\mathbf{z}}$	$(12k)$	Cl II
$\mathbf{B}_{26}$	=	$-2x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$-\frac{3}{2}x_5 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	$(12k)$	Cl II
$\mathbf{B}_{27}$	=	$x_5\,\mathbf{a}_1 + 2x_5\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\frac{3}{2}x_5 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_5 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	$(12k)$	Cl II
$\mathbf{B}_{28}$	=	$x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$-\sqrt{3}x_5 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)c\,\hat{\mathbf{z}}$	$(12k)$	Cl II

References:

- G. G. Wessel and D. J. W. IJdo, *The Crystal Structure of $Cs_3Cr_2Cl_9$*, Acta Cryst. **10**, 466–468 (1957), doi:10.1107/S0365110X57001577.

Found in:

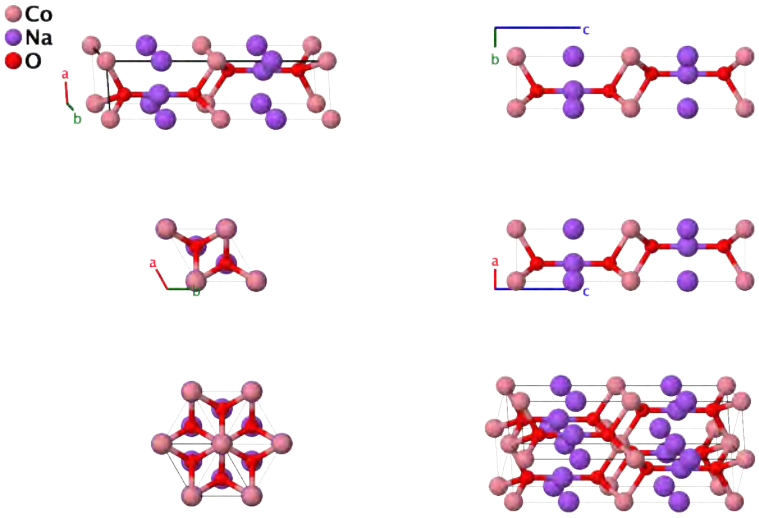
- A. Dönni, A. Furrer, and H. U. Güdel, *Structure of the dimer compounds $Cs_3R_2Br_9$ (R = Tb, Dy, Ho, Er, Yb) at 8 and 295 K studied by neutron diffraction*, J. Solid State Chem. **81**, 278–284 (1989), doi:10.1016/0022-4596(89)90015-7.

Geometry files:

- CIF: pp. 1767
- POSCAR: pp. 1768

$Na_{0.74}CoO_2$ Structure: AB2C2_hP10_194_a_bc_f

http://aflow.org/prototype-encyclopedia/AB2C2_hP10_194_a_bc_f

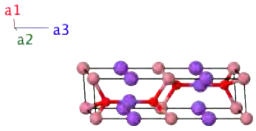


Prototype	:	$CoNa_{0.74}O_2$
AFLOW prototype label	:	AB2C2_hP10_194_a_bc_f
***Strukturbericht* designation**	:	None
Pearson symbol	:	hP10
Space group number	:	194
Space group symbol	:	$P6_3/mmc$
AFLOW prototype command	:	`aflow --proto=AB2C2_hP10_194_a_bc_f --params=`$a, c/a, z_4$

- $Na_{0.74}CoO_2$ is a high figure-of-merit thermoelectric (Sk, 2019). The sodium sites are only partial filled, with the Na-I ($2b$) site having 21% occupancy while the Na-II ($2c$) site is at 51%.

Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($2a$)	Co
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	($2a$)	Co
$\mathbf{B}_3$	=	$\frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\,\hat{\mathbf{z}}$	($2b$)	Na I
$\mathbf{B}_4$	=	$\frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\,\hat{\mathbf{z}}$	($2b$)	Na I
$\mathbf{B}_5$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($2c$)	Na II

$\mathbf{B}_6$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(2c)$	Na II
$\mathbf{B}_7$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(4f)$	O
$\mathbf{B}_8$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	$(4f)$	O
$\mathbf{B}_9$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	$(4f)$	O
$\mathbf{B}_{10}$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	$(4f)$	O

References:

- R. J. Balsys and R. L. Davis, *Refinement of the structure of $Na_{0.74}CoO_2$ using neutron powder diffraction*, Solid State Ion. **93**, 279–282 (1997), doi:10.1016/S0167-2738(96)00557-7.

Found in:

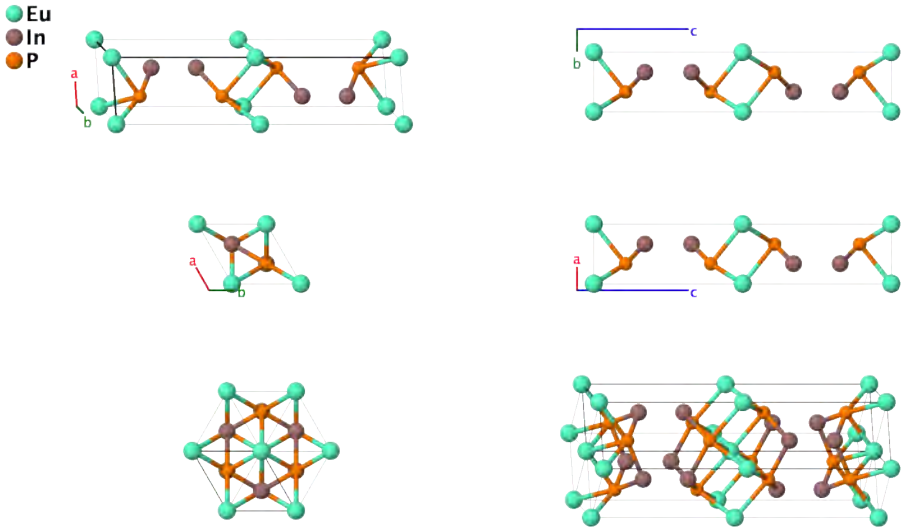
- S. Sk, J. Pati, R. S. Dhaka, and S. K. Pandey, *Exploring the possibility of enhancing the figure-of-merit (> 2) of $Na_{0.74}CoO_2$: A combined experimental and theoretical study*, Eur. Phys. J. B **93**, 155 (2020), doi:10.1140/epjb/e2020-10227-x.

Geometry files:

- CIF: pp. 1768
- POSCAR: pp. 1768

$EuIn_2P_2$ Structure: AB2C2_hP10_194_a_f_f

http://aflow.org/prototype-encyclopedia/AB2C2_hP10_194_a_f_f



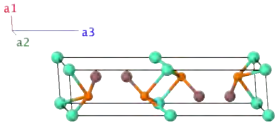
Prototype	:	$EuIn_2P_2$
AFLOW prototype label	:	AB2C2_hP10_194_a_f_f
***Strukturbericht* designation**	:	None
Pearson symbol	:	hP10
Space group number	:	194
Space group symbol	:	$P6_3/mmc$
AFLOW prototype command	:	`aflow --proto=AB2C2_hP10_194_a_f_f --params=`$a, c/a, z_2, z_3$

Other compounds with this structure

- $EuIn_2As_2$

Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	Eu
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	$(2a)$	Eu
$\mathbf{B}_3$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(4f)$	In
$\mathbf{B}_4$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	$(4f)$	In
$\mathbf{B}_5$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	$(4f)$	In

$\mathbf{B}_6$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	$(4f)$	In
$\mathbf{B}_7$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(4f)$	P
$\mathbf{B}_8$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	$(4f)$	P
$\mathbf{B}_9$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(4f)$	P
$\mathbf{B}_{10}$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	$(4f)$	P

References:

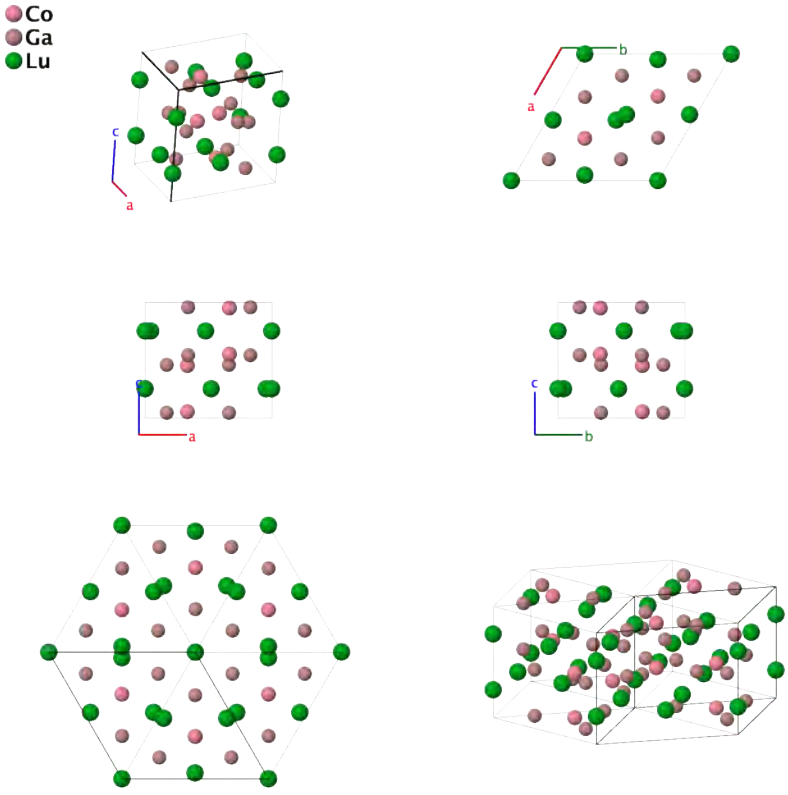
- J. Jiang and S. M. Kauzlarich, *Colossal Magnetoresistance in a Rare Earth Zintl Compound with a New Structure Type: $EuIn_2P_2$*, Chem. Mater. **18**, 435–441 (2006), doi:10.1021/cm0520362.

Geometry files:

- CIF: pp. 1769
- POSCAR: pp. 1769

Lu_2CoGa_3 Structure: AB3C2_hP24_194_f_k_bh

http://aflow.org/prototype-encyclopedia/AB3C2_hP24_194_f_k_bh



Prototype	:	$CoGa_3Lu_2$
AFLOW prototype label	:	AB3C2_hP24_194_f_k_bh
***Strukturbericht* designation**	:	None
Pearson symbol	:	hP24
Space group number	:	194
Space group symbol	:	$P6_3/mmc$
AFLOW prototype command	:	`aflow --proto=AB3C2_hP24_194_f_k_bh --params=`$a, c/a, z_2, x_3, x_4, z_4$

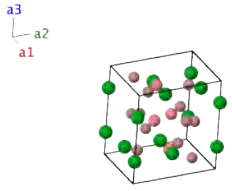
Other compounds with this structure

- Dy_2CoGa_3, Er_2CoGa_3, Er_2RhSi_3, Ho_2CoGa_3, Nd_2PdSi_3, Tb_2CoGa_3, Tm_2CoGa_3, Y_2CoGa_3, and Yb_2CoGa_3

- Although (Gladyshevskii, 1992) refers to these compounds as "new members of the AlB_2 structure family" and this structure is related to AlB_2, the two are not isostructural.

Hexagonal primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= c\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}c\,\hat{\mathbf{z}}$	$(2b)$	Lu I
$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}c\,\hat{\mathbf{z}}$	$(2b)$	Lu I
$\mathbf{B}_3$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_2 c\,\hat{\mathbf{z}}$	$(4f)$	Co
$\mathbf{B}_4$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_2\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_2\right)c\,\hat{\mathbf{z}}$	$(4f)$	Co
$\mathbf{B}_5$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_2 c\,\hat{\mathbf{z}}$	$(4f)$	Co
$\mathbf{B}_6$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_2\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)c\,\hat{\mathbf{z}}$	$(4f)$	Co
$\mathbf{B}_7$	=	$x_3\,\mathbf{a}_1 + 2x_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{3}{2}x_3 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	Lu II
$\mathbf{B}_8$	=	$-2x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-\frac{3}{2}x_3 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	Lu II
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$-\sqrt{3}x_3 a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(6h)$	Lu II
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 - 2x_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-\frac{3}{2}x_3 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	Lu II
$\mathbf{B}_{11}$	=	$2x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{2}x_3 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_3 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	Lu II
$\mathbf{B}_{12}$	=	$-x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\sqrt{3}x_3 a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(6h)$	Lu II
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + 2x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\frac{3}{2}x_4 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(12k)$	Ga
$\mathbf{B}_{14}$	=	$-2x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$-\frac{3}{2}x_4 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(12k)$	Ga
$\mathbf{B}_{15}$	=	$x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$-\sqrt{3}x_4 a\,\hat{\mathbf{y}} + z_4 c\,\hat{\mathbf{z}}$	$(12k)$	Ga
$\mathbf{B}_{16}$	=	$-x_4\,\mathbf{a}_1 - 2x_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$-\frac{3}{2}x_4 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	$(12k)$	Ga
$\mathbf{B}_{17}$	=	$2x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\frac{3}{2}x_4 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	$(12k)$	Ga
$\mathbf{B}_{18}$	=	$-x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\sqrt{3}x_4 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)c\,\hat{\mathbf{z}}$	$(12k)$	Ga
$\mathbf{B}_{19}$	=	$2x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\frac{3}{2}x_4 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	$(12k)$	Ga
$\mathbf{B}_{20}$	=	$-x_4\,\mathbf{a}_1 - 2x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-\frac{3}{2}x_4 a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	$(12k)$	Ga
$\mathbf{B}_{21}$	=	$-x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\sqrt{3}x_4 a\,\hat{\mathbf{y}} - z_4 c\,\hat{\mathbf{z}}$	$(12k)$	Ga
$\mathbf{B}_{22}$	=	$-2x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$-\frac{3}{2}x_4 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	$(12k)$	Ga
$\mathbf{B}_{23}$	=	$x_4\,\mathbf{a}_1 + 2x_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\frac{3}{2}x_4 a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}x_4 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	$(12k)$	Ga
$\mathbf{B}_{24}$	=	$x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$-\sqrt{3}x_4 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)c\,\hat{\mathbf{z}}$	$(12k)$	Ga

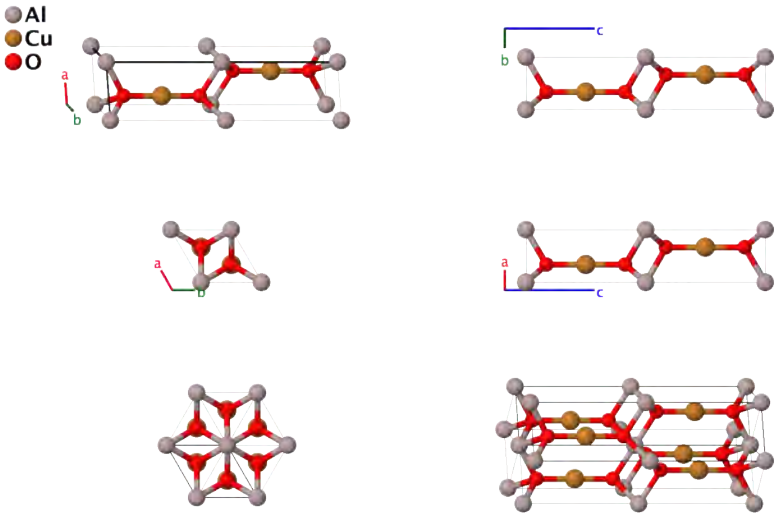
References:

- R. E. Gladyshevskii, K. Cenzual, and E. Parthé, *Er_2RhSi_3 and R_2CoGa_3 (R = Y, Tb, Dy, Ho, Er, Tm, Yb) with Lu_2CoGa_3 type structure: new members of the AlB_2 structure family*, J. Alloys Compd. **189**, 221–228 (1992), doi:10.1016/0925-8388(92)90711-H.

Geometry files:
- CIF: pp. 1769
- POSCAR: pp. 1769

Hexagonal Delafossite ($CuAlO_2$) Structure: ABC2_hP8_194_a_c_f

http://aflow.org/prototype-encyclopedia/ABC2_hP8_194_a_c_f



Prototype	:	$AlCuO_2$
AFLOW prototype label	:	ABC2_hP8_194_a_c_f
***Strukturbericht* designation**	:	None
Pearson symbol	:	hP8
Space group number	:	194
Space group symbol	:	$P6_3/mmc$
AFLOW prototype command	:	`aflow --proto=ABC2_hP8_194_a_c_f --params=`$a, c/a, z_3$

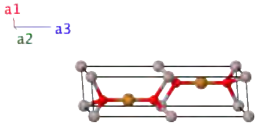
Other compounds with this structure

- $CuGaO_2$, $CuScO_2$, and $CuYO_2$

- Delafossite appears in two forms which differ in the stacking of the layers: rhombohedral, prototype $CuFeO_2$, and hexagonal, shown here. Most of the structures found in the hexagonal phase can also be found in the rhombohedral structure (Marquardt, 2006).

Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	Al
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	$(2a)$	Al
$\mathbf{B}_3$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(2c)$	Cu
$\mathbf{B}_4$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(2c)$	Cu
$\mathbf{B}_5$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + z_3 c\,\hat{\mathbf{z}}$	$(4f)$	O
$\mathbf{B}_6$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)c\,\hat{\mathbf{z}}$	$(4f)$	O
$\mathbf{B}_7$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} - z_3 c\,\hat{\mathbf{z}}$	$(4f)$	O
$\mathbf{B}_8$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)c\,\hat{\mathbf{z}}$	$(4f)$	O

References:

- B. U. Köhler and M. Jansen, *Darstellung und Strukturdaten von, "Delafossiten" $CuMO_2$ (M = Al, Ga, Sc, Y)*, Z. Anorg. Allg. Chem. **543**, 73–80 (1986), doi:10.1002/zaac.19865431209.

Found in:

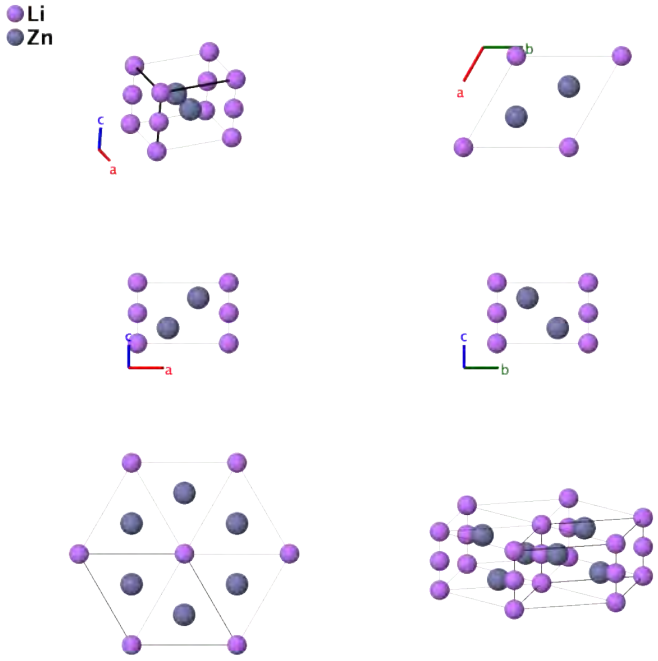
- M. A. Marquardt, N. A. Ashmore, and D. P. Cann, *Crystal chemistry and electrical properties of the delafossite structure*, Thin Solid Films **496**, 146–156 (2006), doi:10.1016/j.tsf.2005.08.316.

Geometry files:

- CIF: pp. 1770
- POSCAR: pp. 1770

$LiZn_2$ (C_k) Structure: AB_hP4_194_a_c

http://aflow.org/prototype-encyclopedia/AB_hP4_194_a_c

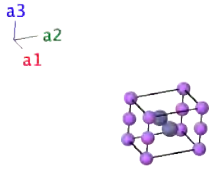


Prototype	:	$LiZn_2$
AFLOW prototype label	:	AB_hP4_194_a_c
***Strukturbericht* designation**	:	C_k
Pearson symbol	:	hP4
Space group number	:	194
Space group symbol	:	$P6_3/mmc$
AFLOW prototype command	:	`aflow --proto=AB_hP4_194_a_c --params=`$a, c/a$

- The actual stoichiometry of this compounds is $Li_{0.8}Zn_2$, so each ($2a$) site is only filled 40% of the time. This means that the short (1.26 Å) distance between the lithium atoms shown here is illusionary.

Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}\,a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($2a$)	Li

$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	($2a$)	Li
$\mathbf{B}_3$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	($2c$)	Zn
$\mathbf{B}_4$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	($2c$)	Zn

References:

- E. Zintl and A. Schneider, *Röntgenanalyse der Lithium-Zink-Legierungen (15. Mitteilung über Metalle und Legierungen)*, Z. Elektrochem. **41**, 764–767 (1935), doi:10.1002/bbpc.19350411103.

Found in:

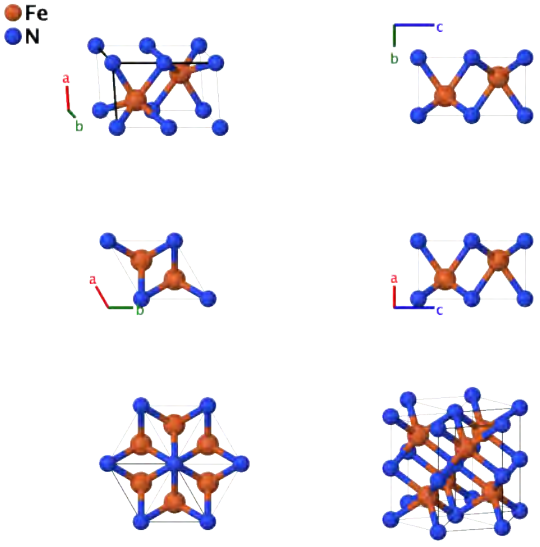
- W. B. Pearson, *A Handbook of Lattice Spacings and Structures of Metals and Alloys*, *International Series of Monographs on Metal Physics and Physical Metallurgy*, vol. 4 (Pergamon Press, Oxford, London, Edinburgh, New York, Paris, Frankfort, 1958), 1964 reprint with corrections edn. N. R. C. No. 4303.

Geometry files:

- CIF: pp. 1770
- POSCAR: pp. 1771

Fe_2N (approximate, $L'3_0$) Structure: AB_hP4_194_c_a

http://aflow.org/prototype-encyclopedia/AB_hP4_194_c_a.Fe2N



Prototype	:	Fe_2N
AFLOW prototype label	:	AB_hP4_194_c_a
***Strukturbericht* designation**	:	$L'3_0$
Pearson symbol	:	hP4
Space group number	:	194
Space group symbol	:	$P6_3/mmc$
AFLOW prototype command	:	`aflow --proto=AB_hP4_194_c_a --params=`$a, c/a$

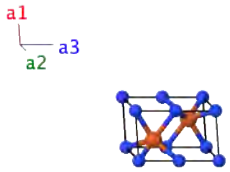
Other compounds with this structure

- β-CTa_2, CV_2, CW_2, Mn_2N, Nb_2N, and $Ta_{\approx 2}N$

- The $L'3_0$ designation is only found in (Smithells, 1955) and (Pearson, 1958), as well as their following volumes. Smithells gives two possible positions of the nitrogen atoms. We use the one that matches Pearson, which (Parthé, 1993) notes is equivalent to the NiAs $B8_1$ structure, but here the ($2a$) nitrogen site is only half-filled to maintain stoichiometry.
- Neither of the defining references gives a source for the structure. It is likely an approximation to ϵ-Fe_2N, which has the β-V_2N $L'3_2$ structure. We take the lattice constants for this structure from the Fe-Fe distances in ϵ-Fe_2N provided by (Hendricks, 1930).
- The similar compounds list is taken from (Pearson, 1967). As it is quite similar to the list for β-V_2N, it is likely that this structure is only an approximation to the correct structure for these compounds as well.
- This structure has the same AFLOW designation, AB_hP4_194_c_a.Fe2N, as $B8_1$. The two structures differ by their lattice constants and the half-occupancy of the ($2a$) site in $L'3_0$.

Hexagonal primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = c\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(2a)$	N
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}c\,\hat{\mathbf{z}}$	$(2a)$	N
$\mathbf{B}_3$	=	$\frac{1}{3}\,\mathbf{a}_1 + \frac{2}{3}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{1}{4}c\,\hat{\mathbf{z}}$	$(2c)$	Fe
$\mathbf{B}_4$	=	$\frac{2}{3}\,\mathbf{a}_1 + \frac{1}{3}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2\sqrt{3}}a\,\hat{\mathbf{y}} + \frac{3}{4}c\,\hat{\mathbf{z}}$	$(2c)$	Fe

References:

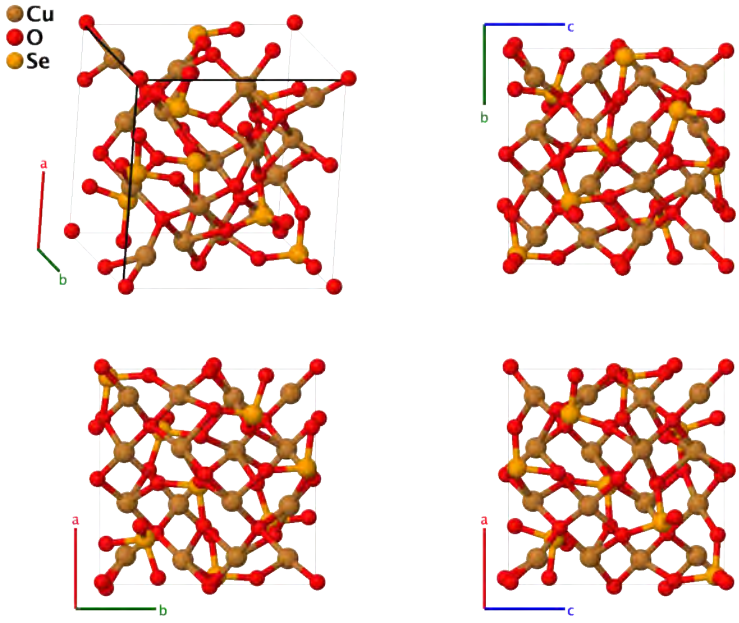
- C. J. Smithells, *Metals Reference Book* (Butterworths Scientific, London, 1955), second edn.
- W. B. Pearson, *A Handbook of Lattice Spacings and Structures of Metals and Alloys*, *International Series of Monographs on Metal Physics and Physical Metallurgy*, vol. 4 (Pergamon Press, Oxford, London, Edinburgh, New York, Paris, Frankfort, 1958), 1964 reprint with corrections edn. N. R. C. No. 4303.
- E. Parthé, L. Gelato, B. Chabot, M. Penzo, K. Cenzual, and R. Gladyshevskii, in *Standardized Data and Crystal Chemical Characterization of Inorganic Structure Types* (Springer-Verlag, Berlin, Heidelberg, 1993), *Gmelin Handbook of Inorganic and Organometallic Chemistry*, vol. 2, chap. Crystal Chemical Characterization of Inorganic Structure Types, 8 edn., doi:10.1007/978-3-662-02909-1_3.
- S. B. Hendricks and P. R. Kosting, *The Crystal Structure of Fe_2P, Fe_2N, Fe_3N and FeB*, Zeitschrift für Kristallographie - Crystalline Materials **74**, 511–533 (1930), doi:10.1524/zkri.1930.74.1.511.
- W. B. Pearson, *A Handbook of Lattice Spacings and Structures of Metals and Alloys, Volume 2*, *International Series of Monographs on Metal Physics and Physical Metallurgy*, vol. 8 (Pergamon Press, Oxford, London, Edinburgh, New York, Toronto, Sydney, Paris, Braunschweig, 1967). N. R. C. No. 8752.

Geometry files:

- CIF: pp. 1771
- POSCAR: pp. 1771

Cubic Cu_2OSeO_3 Structure: A2B4C_cP56_198_ab_2a2b_2a

http://aflow.org/prototype-encyclopedia/A2B4C_cP56_198_ab_2a2b_2a

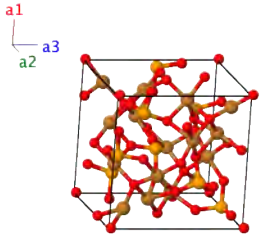


Prototype	:	Cu_2O_4Se
AFLOW prototype label	:	A2B4C_cP56_198_ab_2a2b_2a
***Strukturbericht* designation**	:	None
Pearson symbol	:	cP56
Space group number	:	198
Space group symbol	:	$P2_13$
AFLOW prototype command	:	`aflow --proto=A2B4C_cP56_198_ab_2a2b_2a --params=`$a, x_1, x_2, x_3, x_4, x_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8$

- This is the cubic phase of Cu_2OSeO_3. There is also a monoclinic phase.

Simple Cubic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = a\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$ =	$x_1\,\mathbf{a}_1 + x_1\,\mathbf{a}_2 + x_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + x_1 a\,\hat{\mathbf{y}} + x_1 a\,\hat{\mathbf{z}}$	($4a$)	Cu I

$$
\begin{aligned}
\mathbf{B}_{2} &= \left(\tfrac{1}{2}-x_1\right)\mathbf{a}_1 - x_1\,\mathbf{a}_2 + \left(\tfrac{1}{2}+x_1\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_1\right)a\,\hat{\mathbf{x}} - x_1 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_1\right)a\,\hat{\mathbf{z}} && (4a) && \text{Cu I}\\
\mathbf{B}_{3} &= -x_1\,\mathbf{a}_1 + \left(\tfrac{1}{2}+x_1\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-x_1\right)\mathbf{a}_3 &&= -x_1 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_1\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-x_1\right)a\,\hat{\mathbf{z}} && (4a) && \text{Cu I}\\
\mathbf{B}_{4} &= \left(\tfrac{1}{2}+x_1\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-x_1\right)\mathbf{a}_2 - x_1\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_1\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_1\right)a\,\hat{\mathbf{y}} - x_1 a\,\hat{\mathbf{z}} && (4a) && \text{Cu I}\\
\mathbf{B}_{5} &= x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + x_2\,\mathbf{a}_3 &&= x_2 a\,\hat{\mathbf{x}} + x_2 a\,\hat{\mathbf{y}} + x_2 a\,\hat{\mathbf{z}} && (4a) && \text{O I}\\
\mathbf{B}_{6} &= \left(\tfrac{1}{2}-x_2\right)\mathbf{a}_1 - x_2\,\mathbf{a}_2 + \left(\tfrac{1}{2}+x_2\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_2\right)a\,\hat{\mathbf{x}} - x_2 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_2\right)a\,\hat{\mathbf{z}} && (4a) && \text{O I}\\
\mathbf{B}_{7} &= -x_2\,\mathbf{a}_1 + \left(\tfrac{1}{2}+x_2\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-x_2\right)\mathbf{a}_3 &&= -x_2 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_2\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-x_2\right)a\,\hat{\mathbf{z}} && (4a) && \text{O I}\\
\mathbf{B}_{8} &= \left(\tfrac{1}{2}+x_2\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-x_2\right)\mathbf{a}_2 - x_2\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_2\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_2\right)a\,\hat{\mathbf{y}} - x_2 a\,\hat{\mathbf{z}} && (4a) && \text{O I}\\
\mathbf{B}_{9} &= x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3 &&= x_3 a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}} + x_3 a\,\hat{\mathbf{z}} && (4a) && \text{O II}\\
\mathbf{B}_{10} &= \left(\tfrac{1}{2}-x_3\right)\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\tfrac{1}{2}+x_3\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_3\right)a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_3\right)a\,\hat{\mathbf{z}} && (4a) && \text{O II}\\
\mathbf{B}_{11} &= -x_3\,\mathbf{a}_1 + \left(\tfrac{1}{2}+x_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-x_3\right)\mathbf{a}_3 &&= -x_3 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_3\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-x_3\right)a\,\hat{\mathbf{z}} && (4a) && \text{O II}\\
\mathbf{B}_{12} &= \left(\tfrac{1}{2}+x_3\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-x_3\right)\mathbf{a}_2 - x_3\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_3\right)a\,\hat{\mathbf{y}} - x_3 a\,\hat{\mathbf{z}} && (4a) && \text{O II}\\
\mathbf{B}_{13} &= x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3 &&= x_4 a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}} + x_4 a\,\hat{\mathbf{z}} && (4a) && \text{Se I}\\
\mathbf{B}_{14} &= \left(\tfrac{1}{2}-x_4\right)\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \left(\tfrac{1}{2}+x_4\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_4\right)a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_4\right)a\,\hat{\mathbf{z}} && (4a) && \text{Se I}\\
\mathbf{B}_{15} &= -x_4\,\mathbf{a}_1 + \left(\tfrac{1}{2}+x_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-x_4\right)\mathbf{a}_3 &&= -x_4 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-x_4\right)a\,\hat{\mathbf{z}} && (4a) && \text{Se I}\\
\mathbf{B}_{16} &= \left(\tfrac{1}{2}+x_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-x_4\right)\mathbf{a}_2 - x_4\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_4\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_4\right)a\,\hat{\mathbf{y}} - x_4 a\,\hat{\mathbf{z}} && (4a) && \text{Se I}\\
\mathbf{B}_{17} &= x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3 &&= x_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}} && (4a) && \text{Se II}\\
\mathbf{B}_{18} &= \left(\tfrac{1}{2}-x_5\right)\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \left(\tfrac{1}{2}+x_5\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_5\right)a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_5\right)a\,\hat{\mathbf{z}} && (4a) && \text{Se II}\\
\mathbf{B}_{19} &= -x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2}+x_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-x_5\right)\mathbf{a}_3 &&= -x_5 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-x_5\right)a\,\hat{\mathbf{z}} && (4a) && \text{Se II}\\
\mathbf{B}_{20} &= \left(\tfrac{1}{2}+x_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-x_5\right)\mathbf{a}_2 - x_5\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_5\right)a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}} && (4a) && \text{Se II}\\
\mathbf{B}_{21} &= x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 &&= x_6 a\,\hat{\mathbf{x}} + y_6 a\,\hat{\mathbf{y}} + z_6 a\,\hat{\mathbf{z}} && (12b) && \text{Cu II}\\
\mathbf{B}_{22} &= \left(\tfrac{1}{2}-x_6\right)\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{x}} - y_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_6\right)a\,\hat{\mathbf{z}} && (12b) && \text{Cu II}\\
\mathbf{B}_{23} &= -x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_6\right)\mathbf{a}_3 &&= -x_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_6\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_6\right)a\,\hat{\mathbf{z}} && (12b) && \text{Cu II}\\
\mathbf{B}_{24} &= \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_6\right)a\,\hat{\mathbf{y}} - z_6 a\,\hat{\mathbf{z}} && (12b) && \text{Cu II}\\
\mathbf{B}_{25} &= z_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + y_6\,\mathbf{a}_3 &&= z_6 a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} + y_6 a\,\hat{\mathbf{z}} && (12b) && \text{Cu II}\\
\mathbf{B}_{26} &= \left(\tfrac{1}{2}+z_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-x_6\right)\mathbf{a}_2 - y_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+z_6\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{y}} - y_6 a\,\hat{\mathbf{z}} && (12b) && \text{Cu II}\\
\mathbf{B}_{27} &= \left(\tfrac{1}{2}-z_6\right)\mathbf{a}_1 - x_6\,\mathbf{a}_2 + \left(\tfrac{1}{2}+y_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-z_6\right)a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+y_6\right)a\,\hat{\mathbf{z}} && (12b) && \text{Cu II}\\
\mathbf{B}_{28} &= -z_6\,\mathbf{a}_1 + \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-y_6\right)\mathbf{a}_3 &&= -z_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-y_6\right)a\,\hat{\mathbf{z}} && (12b) && \text{Cu II}\\
\mathbf{B}_{29} &= y_6\,\mathbf{a}_1 + z_6\,\mathbf{a}_2 + x_6\,\mathbf{a}_3 &&= y_6 a\,\hat{\mathbf{x}} + z_6 a\,\hat{\mathbf{y}} + x_6 a\,\hat{\mathbf{z}} && (12b) && \text{Cu II}\\
\mathbf{B}_{30} &= -y_6\,\mathbf{a}_1 + \left(\tfrac{1}{2}+z_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-x_6\right)\mathbf{a}_3 &&= -y_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+z_6\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{z}} && (12b) && \text{Cu II}\\
\mathbf{B}_{31} &= \left(\tfrac{1}{2}+y_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-z_6\right)\mathbf{a}_2 - x_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+y_6\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-z_6\right)a\,\hat{\mathbf{y}} - x_6 a\,\hat{\mathbf{z}} && (12b) && \text{Cu II}\\
\mathbf{B}_{32} &= \left(\tfrac{1}{2}-y_6\right)\mathbf{a}_1 - z_6\,\mathbf{a}_2 + \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-y_6\right)a\,\hat{\mathbf{x}} - z_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{z}} && (12b) && \text{Cu II}\\
\mathbf{B}_{33} &= x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 &&= x_7 a\,\hat{\mathbf{x}} + y_7 a\,\hat{\mathbf{y}} + z_7 a\,\hat{\mathbf{z}} && (12b) && \text{O III}\\
\mathbf{B}_{34} &= \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{x}} - y_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{z}} && (12b) && \text{O III}\\
\mathbf{B}_{35} &= -x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_3 &&= -x_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_7\right)a\,\hat{\mathbf{z}} && (12b) && \text{O III}\\
\mathbf{B}_{36} &= \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_7\right)\mathbf{a}_2 - z_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_7\right)a\,\hat{\mathbf{y}} - z_7 a\,\hat{\mathbf{z}} && (12b) && \text{O III}\\
\mathbf{B}_{37} &= z_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + y_7\,\mathbf{a}_3 &&= z_7 a\,\hat{\mathbf{x}} + x_7 a\,\hat{\mathbf{y}} + y_7 a\,\hat{\mathbf{z}} && (12b) && \text{O III}
\end{aligned}
$$

$$\mathbf{B}_{38} = \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}-x_7\right)\mathbf{a}_2-y_7\,\mathbf{a}_3 = \left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{y}}-y_7a\,\hat{\mathbf{z}} \quad (12b) \quad \text{O III}$$

$$\mathbf{B}_{39} = \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_1-x_7\,\mathbf{a}_2+\left(\tfrac{1}{2}+y_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-z_7\right)a\,\hat{\mathbf{x}}-x_7a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+y_7\right)a\,\hat{\mathbf{z}} \quad (12b) \quad \text{O III}$$

$$\mathbf{B}_{40} = -z_7\,\mathbf{a}_1+\left(\tfrac{1}{2}+x_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}-y_7\right)\mathbf{a}_3 = -z_7a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-y_7\right)a\,\hat{\mathbf{z}} \quad (12b) \quad \text{O III}$$

$$\mathbf{B}_{41} = y_7\,\mathbf{a}_1+z_7\,\mathbf{a}_2+x_7\,\mathbf{a}_3 = y_7a\,\hat{\mathbf{x}}+z_7a\,\hat{\mathbf{y}}+x_7a\,\hat{\mathbf{z}} \quad (12b) \quad \text{O III}$$

$$\mathbf{B}_{42} = -y_7\,\mathbf{a}_1+\left(\tfrac{1}{2}+z_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}-x_7\right)\mathbf{a}_3 = -y_7a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{z}} \quad (12b) \quad \text{O III}$$

$$\mathbf{B}_{43} = \left(\tfrac{1}{2}+y_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}-z_7\right)\mathbf{a}_2-x_7\,\mathbf{a}_3 = \left(\tfrac{1}{2}+y_7\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-z_7\right)a\,\hat{\mathbf{y}}-x_7a\,\hat{\mathbf{z}} \quad (12b) \quad \text{O III}$$

$$\mathbf{B}_{44} = \left(\tfrac{1}{2}-y_7\right)\mathbf{a}_1-z_7\,\mathbf{a}_2+\left(\tfrac{1}{2}+x_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-y_7\right)a\,\hat{\mathbf{x}}-z_7a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{z}} \quad (12b) \quad \text{O III}$$

$$\mathbf{B}_{45} = x_8\,\mathbf{a}_1+y_8\,\mathbf{a}_2+z_8\,\mathbf{a}_3 = x_8a\,\hat{\mathbf{x}}+y_8a\,\hat{\mathbf{y}}+z_8a\,\hat{\mathbf{z}} \quad (12b) \quad \text{O IV}$$

$$\mathbf{B}_{46} = \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_1-y_8\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{x}}-y_8a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_8\right)a\,\hat{\mathbf{z}} \quad (12b) \quad \text{O IV}$$

$$\mathbf{B}_{47} = -x_8\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_8\right)\mathbf{a}_3 = -x_8a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_8\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_8\right)a\,\hat{\mathbf{z}} \quad (12b) \quad \text{O IV}$$

$$\mathbf{B}_{48} = \left(\tfrac{1}{2}+x_8\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_8\right)\mathbf{a}_2-z_8\,\mathbf{a}_3 = \left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_8\right)a\,\hat{\mathbf{y}}-z_8a\,\hat{\mathbf{z}} \quad (12b) \quad \text{O IV}$$

$$\mathbf{B}_{49} = z_8\,\mathbf{a}_1+x_8\,\mathbf{a}_2+y_8\,\mathbf{a}_3 = z_8a\,\hat{\mathbf{x}}+x_8a\,\hat{\mathbf{y}}+y_8a\,\hat{\mathbf{z}} \quad (12b) \quad \text{O IV}$$

$$\mathbf{B}_{50} = \left(\tfrac{1}{2}+z_8\right)\mathbf{a}_1+\left(\tfrac{1}{2}-x_8\right)\mathbf{a}_2-y_8\,\mathbf{a}_3 = \left(\tfrac{1}{2}+z_8\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{y}}-y_8a\,\hat{\mathbf{z}} \quad (12b) \quad \text{O IV}$$

$$\mathbf{B}_{51} = \left(\tfrac{1}{2}-z_8\right)\mathbf{a}_1-x_8\,\mathbf{a}_2+\left(\tfrac{1}{2}+y_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-z_8\right)a\,\hat{\mathbf{x}}-x_8a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+y_8\right)a\,\hat{\mathbf{z}} \quad (12b) \quad \text{O IV}$$

$$\mathbf{B}_{52} = -z_8\,\mathbf{a}_1+\left(\tfrac{1}{2}+x_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}-y_8\right)\mathbf{a}_3 = -z_8a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-y_8\right)a\,\hat{\mathbf{z}} \quad (12b) \quad \text{O IV}$$

$$\mathbf{B}_{53} = y_8\,\mathbf{a}_1+z_8\,\mathbf{a}_2+x_8\,\mathbf{a}_3 = y_8a\,\hat{\mathbf{x}}+z_8a\,\hat{\mathbf{y}}+x_8a\,\hat{\mathbf{z}} \quad (12b) \quad \text{O IV}$$

$$\mathbf{B}_{54} = -y_8\,\mathbf{a}_1+\left(\tfrac{1}{2}+z_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}-x_8\right)\mathbf{a}_3 = -y_8a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+z_8\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{z}} \quad (12b) \quad \text{O IV}$$

$$\mathbf{B}_{55} = \left(\tfrac{1}{2}+y_8\right)\mathbf{a}_1+\left(\tfrac{1}{2}-z_8\right)\mathbf{a}_2-x_8\,\mathbf{a}_3 = \left(\tfrac{1}{2}+y_8\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-z_8\right)a\,\hat{\mathbf{y}}-x_8a\,\hat{\mathbf{z}} \quad (12b) \quad \text{O IV}$$

$$\mathbf{B}_{56} = \left(\tfrac{1}{2}-y_8\right)\mathbf{a}_1-z_8\,\mathbf{a}_2+\left(\tfrac{1}{2}+x_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-y_8\right)a\,\hat{\mathbf{x}}-z_8a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{z}} \quad (12b) \quad \text{O IV}$$

References:

- H. Effenberger and F. Pertlik, *Die Kristallstrukturen der Kupfer(II)-oxo-selenite $Cu_2O(SeO_3)$ (kubisch und monoklin) und $Cu_4O(SeO_3)_3$ (monoklin und triklin)*, Monatshefte für Chemie - Chemical Monthly **117**, 887–896 (1986), doi:10.1007/BF00811258.

Found in:

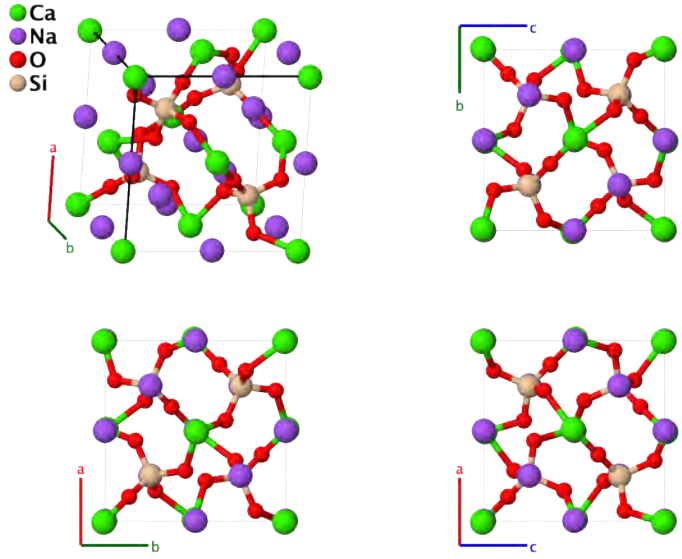
- P. Y. Portnichenko, J. Romhányi, Y. A. Onykiienko, A. Henschel, M. Schmidt, A. S. Cameron, M. A. Surmach, J. A. Lim, J. T. Park, A. Schneidewind, D. L. Abernathy, H. Rosner, J. van den Brink, and D. S. Inosov, *Magnon spectrum of the helimagnetic insulator Cu_2OSeO_3*, Nat. Commun. **7**, 10725 (2016), doi:10.1038/ncomms10725.

Geometry files:

- CIF: pp. 1771
- POSCAR: pp. 1771

Na_2CaSiO_4 ($S6_6$) Structure: AB2C4D_cP32_198_a_2a_ab_a

http://aflow.org/prototype-encyclopedia/AB2C4D_cP32_198_a_2a_ab_a



Prototype	:	$CaNa_2O_4Si$
AFLOW prototype label	:	AB2C4D_cP32_198_a_2a_ab_a
***Strukturbericht* designation**	:	$S6_6$
Pearson symbol	:	cP32
Space group number	:	198
Space group symbol	:	$P2_13$
AFLOW prototype command	:	`aflow --proto=AB2C4D_cP32_198_a_2a_ab_a --params=`$a, x_1, x_2, x_3, x_4, x_5, x_6, y_6, z_6$

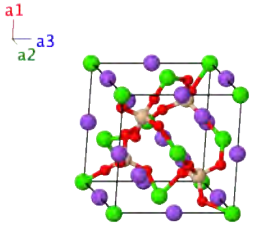
Other compounds with this structure

- Li_2SrSi_4

- Substitution of small fractions of rare earth elements onto the Ca site can produce photoluminescence.
- This is the crystal structure found by (Barth, 1932). The coordinates on pp. 159 of (Hermann, 1937) are incorrect, as they do not correctly transform from Barth and Posnjak's coordinates in degrees to fractional coordinates. See (Hermann, 1937) pp. 557-8 for the correct coordinates.
- (Dollase, 1991) found that the SiO_4 tetrahedra were orientationally disordered, and that the resulting $Na_2Ca(SiO_4)$ structure took on the Heusler ($L1_2$) structure.

Simple Cubic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= a\,\hat{\mathbf{x}}\\ \mathbf{a}_2 &= a\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= a\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$x_1\,\mathbf{a}_1 + x_1\,\mathbf{a}_2 + x_1\,\mathbf{a}_3$	$x_1 a\,\hat{\mathbf{x}} + x_1 a\,\hat{\mathbf{y}} + x_1 a\,\hat{\mathbf{z}}$	(4*a*)	Ca
$\mathbf{B}_2$	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 - x_1\,\mathbf{a}_2 + \left(\frac{1}{2} + x_1\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_1\right)a\,\hat{\mathbf{x}} - x_1 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_1\right)a\,\hat{\mathbf{z}}$	(4*a*)	Ca
$\mathbf{B}_3$	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2} + x_1\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_1\right)\mathbf{a}_3$	$-x_1 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_1\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_1\right)a\,\hat{\mathbf{z}}$	(4*a*)	Ca
$\mathbf{B}_4$	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_1\right)\mathbf{a}_2 - x_1\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_1\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_1\right)a\,\hat{\mathbf{y}} - x_1 a\,\hat{\mathbf{z}}$	(4*a*)	Ca
$\mathbf{B}_5$	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	$x_2 a\,\hat{\mathbf{x}} + x_2 a\,\hat{\mathbf{y}} + x_2 a\,\hat{\mathbf{z}}$	(4*a*)	Na I
$\mathbf{B}_6$	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 - x_2\,\mathbf{a}_2 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} - x_2 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{z}}$	(4*a*)	Na I
$\mathbf{B}_7$	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_3$	$-x_2 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{z}}$	(4*a*)	Na I
$\mathbf{B}_8$	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_2 - x_2\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{y}} - x_2 a\,\hat{\mathbf{z}}$	(4*a*)	Na I
$\mathbf{B}_9$	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	$x_3 a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}} + x_3 a\,\hat{\mathbf{z}}$	(4*a*)	Na II
$\mathbf{B}_{10}$	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{z}}$	(4*a*)	Na II
$\mathbf{B}_{11}$	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_3$	$-x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{z}}$	(4*a*)	Na II
$\mathbf{B}_{12}$	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 - x_3\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} - x_3 a\,\hat{\mathbf{z}}$	(4*a*)	Na II
$\mathbf{B}_{13}$	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	$x_4 a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}} + x_4 a\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_{14}$	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \left(\frac{1}{2} + x_4\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_{15}$	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + x_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_3$	$-x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_{16}$	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_2 - x_4\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{y}} - x_4 a\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_{17}$	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	$x_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}}$	(4*a*)	Si
$\mathbf{B}_{18}$	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \left(\frac{1}{2} + x_5\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{z}}$	(4*a*)	Si
$\mathbf{B}_{19}$	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2} + x_5\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_5\right)\mathbf{a}_3$	$-x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{z}}$	(4*a*)	Si
$\mathbf{B}_{20}$	$\left(\frac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_5\right)\mathbf{a}_2 - x_5\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}}$	(4*a*)	Si
$\mathbf{B}_{21}$	$x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3$	$x_6 a\,\hat{\mathbf{x}} + y_6 a\,\hat{\mathbf{y}} + z_6 a\,\hat{\mathbf{z}}$	(12*b*)	O II
$\mathbf{B}_{22}$	$\left(\frac{1}{2} - x_6\right)\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\frac{1}{2} + z_6\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{x}} - y_6 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_6\right)a\,\hat{\mathbf{z}}$	(12*b*)	O II
$\mathbf{B}_{23}$	$-x_6\,\mathbf{a}_1 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_6\right)\mathbf{a}_3$	$-x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_6\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_6\right)a\,\hat{\mathbf{z}}$	(12*b*)	O II
$\mathbf{B}_{24}$	$\left(\frac{1}{2} + x_6\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_6\right)a\,\hat{\mathbf{y}} - z_6 a\,\hat{\mathbf{z}}$	(12*b*)	O II
$\mathbf{B}_{25}$	$z_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + y_6\,\mathbf{a}_3$	$z_6 a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} + y_6 a\,\hat{\mathbf{z}}$	(12*b*)	O II
$\mathbf{B}_{26}$	$\left(\frac{1}{2} + z_6\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_6\right)\mathbf{a}_2 - y_6\,\mathbf{a}_3$	$\left(\frac{1}{2} + z_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{y}} - y_6 a\,\hat{\mathbf{z}}$	(12*b*)	O II
$\mathbf{B}_{27}$	$\left(\frac{1}{2} - z_6\right)\mathbf{a}_1 - x_6\,\mathbf{a}_2 + \left(\frac{1}{2} + y_6\right)\mathbf{a}_3$	$\left(\frac{1}{2} - z_6\right)a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + y_6\right)a\,\hat{\mathbf{z}}$	(12*b*)	O II

$$\mathbf{B}_{28} = -z_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - y_6\right)\mathbf{a}_3 = -z_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_6\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - y_6\right)a\,\hat{\mathbf{z}} \quad (12b) \quad \text{O II}$$

$$\mathbf{B}_{29} = y_6\,\mathbf{a}_1 + z_6\,\mathbf{a}_2 + x_6\,\mathbf{a}_3 = y_6 a\,\hat{\mathbf{x}} + z_6 a\,\hat{\mathbf{y}} + x_6 a\,\hat{\mathbf{z}} \quad (12b) \quad \text{O II}$$

$$\mathbf{B}_{30} = -y_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_3 = -y_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + z_6\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_6\right)a\,\hat{\mathbf{z}} \quad (12b) \quad \text{O II}$$

$$\mathbf{B}_{31} = \left(\tfrac{1}{2} + y_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_2 - x_6\,\mathbf{a}_3 = \left(\tfrac{1}{2} + y_6\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - z_6\right)a\,\hat{\mathbf{y}} - x_6 a\,\hat{\mathbf{z}} \quad (12b) \quad \text{O II}$$

$$\mathbf{B}_{32} = \left(\tfrac{1}{2} - y_6\right)\mathbf{a}_1 - z_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - y_6\right)a\,\hat{\mathbf{x}} - z_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_6\right)a\,\hat{\mathbf{z}} \quad (12b) \quad \text{O II}$$

References:

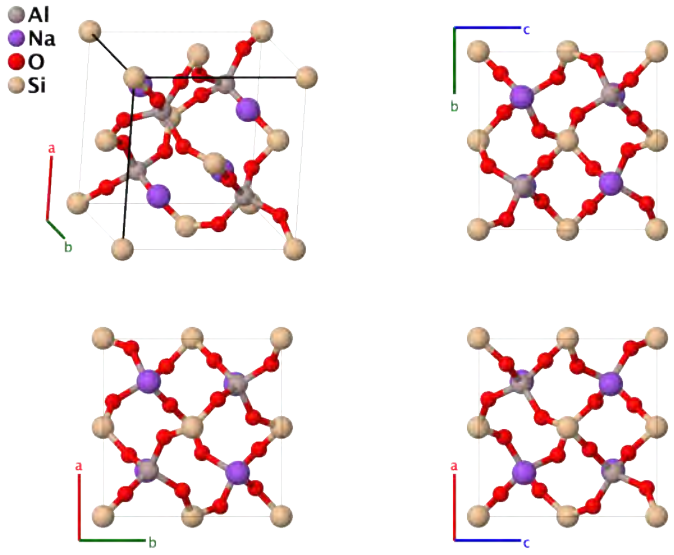
- T. F. W. Barth and E. Posnjak, *Silicate structures of the cristobalite type: II. The crystal structure of Na_2CaSiO_4*, Zeitschrift für Kristallographie - Crystalline Materials **81**, 370–375 (1932), doi:10.1524/zkri.1932.81.1.370.
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).
- W. A. Dollase and C. R. Ross II, *Crystal structure of orientationally disordered $Na_2(Ca,Sr)SiO_4$*, Zeitschrift für Kristallographie - Crystalline Materials **197**, 13–26 (1991), doi:10.1524/zkri.1991.197.1-2.13.

Geometry files:

- CIF: pp. 1772
- POSCAR: pp. 1772

α-Carnegieite ($NaAlSiO_4$, $S6_5$) Structure: ABC4D_cP28_198_a_a_ab_a

http://aflow.org/prototype-encyclopedia/ABC4D_cP28_198_a_a_ab_a

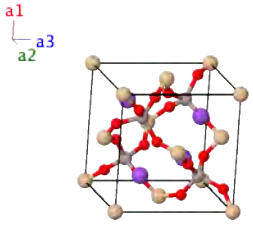


Prototype	:	$AlNaO_4Si$
AFLOW prototype label	:	ABC4D_cP28_198_a_a_ab_a
***Strukturbericht* designation**	:	$S6_5$
Pearson symbol	:	cP28
Space group number	:	198
Space group symbol	:	$P2_13$
AFLOW prototype command	:	`aflow --proto=ABC4D_cP28_198_a_a_ab_a --params=`$a, x_1, x_2, x_3, x_4, x_5, y_5, z_5$

- This high-temperature form of carnegieite is stable above 970 K. To our knowledge, the atomic positions of the low temperature β-carnegieite structure have not been determined.

Simple Cubic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = a\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$ =	$x_1\,\mathbf{a}_1 + x_1\,\mathbf{a}_2 + x_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + x_1 a\,\hat{\mathbf{y}} + x_1 a\,\hat{\mathbf{z}}$	($4a$)	Al

$\mathbf{B}_2$	=	$\left(\frac{1}{2}-x_1\right)\mathbf{a}_1 - x_1\,\mathbf{a}_2 + \left(\frac{1}{2}+x_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_1\right)a\,\hat{\mathbf{x}} - x_1 a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+x_1\right)a\,\hat{\mathbf{z}}$	(4*a*)	Al
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2}+x_1\right)\mathbf{a}_2 + \left(\frac{1}{2}-x_1\right)\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_1\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-x_1\right)a\,\hat{\mathbf{z}}$	(4*a*)	Al
$\mathbf{B}_4$	=	$\left(\frac{1}{2}+x_1\right)\mathbf{a}_1 + \left(\frac{1}{2}-x_1\right)\mathbf{a}_2 - x_1\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_1\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_1\right)a\,\hat{\mathbf{y}} - x_1 a\,\hat{\mathbf{z}}$	(4*a*)	Al
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + x_2 a\,\hat{\mathbf{y}} + x_2 a\,\hat{\mathbf{z}}$	(4*a*)	Na
$\mathbf{B}_6$	=	$\left(\frac{1}{2}-x_2\right)\mathbf{a}_1 - x_2\,\mathbf{a}_2 + \left(\frac{1}{2}+x_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_2\right)a\,\hat{\mathbf{x}} - x_2 a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+x_2\right)a\,\hat{\mathbf{z}}$	(4*a*)	Na
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2}+x_2\right)\mathbf{a}_2 + \left(\frac{1}{2}-x_2\right)\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_2\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-x_2\right)a\,\hat{\mathbf{z}}$	(4*a*)	Na
$\mathbf{B}_8$	=	$\left(\frac{1}{2}+x_2\right)\mathbf{a}_1 + \left(\frac{1}{2}-x_2\right)\mathbf{a}_2 - x_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_2\right)a\,\hat{\mathbf{y}} - x_2 a\,\hat{\mathbf{z}}$	(4*a*)	Na
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}} + x_3 a\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2}-x_3\right)\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\frac{1}{2}+x_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2}+x_3\right)\mathbf{a}_2 + \left(\frac{1}{2}-x_3\right)\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2}+x_3\right)\mathbf{a}_1 + \left(\frac{1}{2}-x_3\right)\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{y}} - x_3 a\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}} + x_4 a\,\hat{\mathbf{z}}$	(4*a*)	Si
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2}-x_4\right)\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \left(\frac{1}{2}+x_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{z}}$	(4*a*)	Si
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2}+x_4\right)\mathbf{a}_2 + \left(\frac{1}{2}-x_4\right)\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{z}}$	(4*a*)	Si
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2}+x_4\right)\mathbf{a}_1 + \left(\frac{1}{2}-x_4\right)\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{y}} - x_4 a\,\hat{\mathbf{z}}$	(4*a*)	Si
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + y_5 a\,\hat{\mathbf{y}} + z_5 a\,\hat{\mathbf{z}}$	(12*b*)	O II
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2}-x_5\right)\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{x}} - y_5 a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_5\right)a\,\hat{\mathbf{z}}$	(12*b*)	O II
$\mathbf{B}_{19}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{1}{2}+y_5\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_5\right)\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_5\right)a\,\hat{\mathbf{z}}$	(12*b*)	O II
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2}+x_5\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_5\right)a\,\hat{\mathbf{y}} - z_5 a\,\hat{\mathbf{z}}$	(12*b*)	O II
$\mathbf{B}_{21}$	=	$z_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + y_5\,\mathbf{a}_3$	=	$z_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + y_5 a\,\hat{\mathbf{z}}$	(12*b*)	O II
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2}+z_5\right)\mathbf{a}_1 + \left(\frac{1}{2}-x_5\right)\mathbf{a}_2 - y_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+z_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{y}} - y_5 a\,\hat{\mathbf{z}}$	(12*b*)	O II
$\mathbf{B}_{23}$	=	$\left(\frac{1}{2}-z_5\right)\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \left(\frac{1}{2}+y_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-z_5\right)a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+y_5\right)a\,\hat{\mathbf{z}}$	(12*b*)	O II
$\mathbf{B}_{24}$	=	$-z_5\,\mathbf{a}_1 + \left(\frac{1}{2}+x_5\right)\mathbf{a}_2 + \left(\frac{1}{2}-y_5\right)\mathbf{a}_3$	=	$-z_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-y_5\right)a\,\hat{\mathbf{z}}$	(12*b*)	O II
$\mathbf{B}_{25}$	=	$y_5\,\mathbf{a}_1 + z_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$y_5 a\,\hat{\mathbf{x}} + z_5 a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}}$	(12*b*)	O II
$\mathbf{B}_{26}$	=	$-y_5\,\mathbf{a}_1 + \left(\frac{1}{2}+z_5\right)\mathbf{a}_2 + \left(\frac{1}{2}-x_5\right)\mathbf{a}_3$	=	$-y_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+z_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{z}}$	(12*b*)	O II
$\mathbf{B}_{27}$	=	$\left(\frac{1}{2}+y_5\right)\mathbf{a}_1 + \left(\frac{1}{2}-z_5\right)\mathbf{a}_2 - x_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-z_5\right)a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}}$	(12*b*)	O II
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2}-y_5\right)\mathbf{a}_1 - z_5\,\mathbf{a}_2 + \left(\frac{1}{2}+x_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_5\right)a\,\hat{\mathbf{x}} - z_5 a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{z}}$	(12*b*)	O II

References:

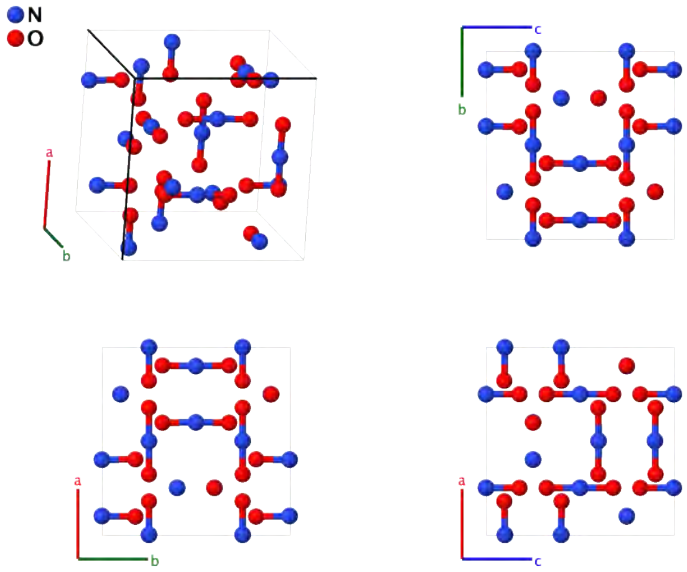
- T. F. W. Barth and E. Posnjak, *Silicate structures of the cristobalite type: I. The crystal structure of α-carnegieite (NaAlSiO$_4$)*, Zeitschrift für Kristallographie - Crystalline Materials **81**, 135–141 (1932), doi:10.1524/zkri.1932.81.1.135.

Geometry files:

- CIF: pp. 1772
- POSCAR: pp. 1773

$C26_a$ (NO_2) (*obsolete*) Structure: AB2_cI36_199_b_c

http://aflow.org/prototype-encyclopedia/AB2_cI36_199_b_c

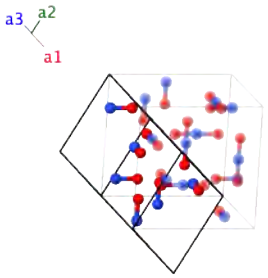


Prototype	:	NO_2
AFLOW prototype label	:	AB2_cI36_199_b_c
***Strukturbericht* designation**	:	$C26_a$
Pearson symbol	:	cI36
Space group number	:	199
Space group symbol	:	$I2_13$
AFLOW prototype command	:	`aflow --proto=AB2_cI36_199_b_c --params=`a, x_1, x_2, y_2, z_2

- (Hermann, 1937) listed two possible structures for the low temperature solid cubic phase of NO_2, which were given *Strukturbericht* designations $C26_a$ and $C26_b$, the only structures with Roman subscripts in the original series.
- $C26_a$ (AB2_cI36_199_b_c) was set in space group $I2_13$ #199. Hermann noted that this structure has a very short distance (1.88 Å) between oxygen atoms on different NO_2 molecules, and that this structure does not form the $(NO_2)_2$ aggregate molecule found in the $C26_b$ structure, making "making this proposed structure very unlikely."
- Recognizing this, (Hendricks, 1931) suggested that NO_2 was actually in space group $I23$ #197. (Hermann, 1997) gave this structure the $C26_b$ designation, but noted that based on Hendricks's atomic positions the space group was actually $Im\bar{3}$ #204.
- The modern accepted structure for NO_2 (AB2_cI36_204_d_g) is set in space group $Im\bar{3}$ #204, confirming Hendricks. We follow (Villars, 2005) and give this the $C26$ designation.

Body-centered Cubic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= -\tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{y}} + \tfrac{1}{2}\,a\,\hat{\mathbf{z}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2}\,a\,\hat{\mathbf{y}} + \tfrac{1}{2}\,a\,\hat{\mathbf{z}}\\ \mathbf{a}_3 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{y}} - \tfrac{1}{2}\,a\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{4}\,\mathbf{a}_1 + \left(\frac{1}{4} + x_1\right)\,\mathbf{a}_2 + x_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(12*b*)	N
$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_1 + \left(\frac{1}{4} - x_1\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - x_1\right)\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(12*b*)	N
$\mathbf{B}_3$	=	$x_1\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{4} + x_1\right)\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + x_1 a\,\hat{\mathbf{y}}$	(12*b*)	N
$\mathbf{B}_4$	=	$\left(\frac{1}{2} - x_1\right)\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{4} - x_1\right)\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} - x_1 a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(12*b*)	N
$\mathbf{B}_5$	=	$\left(\frac{1}{4} + x_1\right)\,\mathbf{a}_1 + x_1\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{y}} + x_1 a\,\hat{\mathbf{z}}$	(12*b*)	N
$\mathbf{B}_6$	=	$\left(\frac{1}{4} - x_1\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - x_1\right)\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} - x_1 a\,\hat{\mathbf{z}}$	(12*b*)	N
$\mathbf{B}_7$	=	$(y_2 + z_2)\,\mathbf{a}_1 + (x_2 + z_2)\,\mathbf{a}_2 + (x_2 + y_2)\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + y_2 a\,\hat{\mathbf{y}} + z_2 a\,\hat{\mathbf{z}}$	(24*c*)	O
$\mathbf{B}_8$	=	$\left(\frac{1}{2} - y_2 + z_2\right)\,\mathbf{a}_1 + (-x_2 + z_2)\,\mathbf{a}_2 + \left(\frac{1}{2} - x_2 - y_2\right)\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_2\right)a\,\hat{\mathbf{y}} + z_2 a\,\hat{\mathbf{z}}$	(24*c*)	O
$\mathbf{B}_9$	=	$(y_2 - z_2)\,\mathbf{a}_1 + \left(\frac{1}{2} - x_2 - z_2\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - x_2 + y_2\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + y_2 a\,\hat{\mathbf{y}} - z_2 a\,\hat{\mathbf{z}}$	(24*c*)	O
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - y_2 - z_2\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + x_2 - z_2\right)\,\mathbf{a}_2 + (x_2 - y_2)\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} - y_2 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_2\right)a\,\hat{\mathbf{z}}$	(24*c*)	O
$\mathbf{B}_{11}$	=	$(x_2 + y_2)\,\mathbf{a}_1 + (y_2 + z_2)\,\mathbf{a}_2 + (x_2 + z_2)\,\mathbf{a}_3$	=	$z_2 a\,\hat{\mathbf{x}} + x_2 a\,\hat{\mathbf{y}} + y_2 a\,\hat{\mathbf{z}}$	(24*c*)	O
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} - x_2 - y_2\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - y_2 + z_2\right)\,\mathbf{a}_2 + (-x_2 + z_2)\,\mathbf{a}_3$	=	$z_2 a\,\hat{\mathbf{x}} - x_2 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - y_2\right)a\,\hat{\mathbf{z}}$	(24*c*)	O
$\mathbf{B}_{13}$	=	$\left(\frac{1}{2} - x_2 + y_2\right)\,\mathbf{a}_1 + (y_2 - z_2)\,\mathbf{a}_2 + \left(\frac{1}{2} - x_2 - z_2\right)\,\mathbf{a}_3$	=	$-z_2 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{y}} + y_2 a\,\hat{\mathbf{z}}$	(24*c*)	O
$\mathbf{B}_{14}$	=	$(x_2 - y_2)\,\mathbf{a}_1 + \left(\frac{1}{2} - y_2 - z_2\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + x_2 - z_2\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - z_2\right)a\,\hat{\mathbf{x}} + x_2 a\,\hat{\mathbf{y}} - y_2 a\,\hat{\mathbf{z}}$	(24*c*)	O
$\mathbf{B}_{15}$	=	$(x_2 + z_2)\,\mathbf{a}_1 + (x_2 + y_2)\,\mathbf{a}_2 + (y_2 + z_2)\,\mathbf{a}_3$	=	$y_2 a\,\hat{\mathbf{x}} + z_2 a\,\hat{\mathbf{y}} + x_2 a\,\hat{\mathbf{z}}$	(24*c*)	O
$\mathbf{B}_{16}$	=	$(-x_2 + z_2)\,\mathbf{a}_1 + \left(\frac{1}{2} - x_2 - y_2\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - y_2 + z_2\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_2\right)a\,\hat{\mathbf{x}} + z_2 a\,\hat{\mathbf{y}} - x_2 a\,\hat{\mathbf{z}}$	(24*c*)	O
$\mathbf{B}_{17}$	=	$\left(\frac{1}{2} - x_2 - z_2\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - x_2 + y_2\right)\,\mathbf{a}_2 + (y_2 - z_2)\,\mathbf{a}_3$	=	$y_2 a\,\hat{\mathbf{x}} - z_2 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{z}}$	(24*c*)	O
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} + x_2 - z_2\right)\,\mathbf{a}_1 + (x_2 - y_2)\,\mathbf{a}_2 + \left(\frac{1}{2} - y_2 - z_2\right)\,\mathbf{a}_3$	=	$-y_2 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - z_2\right)a\,\hat{\mathbf{y}} + x_2 a\,\hat{\mathbf{z}}$	(24*c*)	O

References:

- L. Vegard, *Die Struktur von festem N_2O_4 bei der Temperatur von flüssiger Luft*, Z. Phys. **68**, 184–203 (1931),

doi:10.1007/BF01390966.
- P. Villars and K. Cenzual, eds., *Crystal Structure Data of Inorganic Compounds* (Springer-Verlag, Berlin, Heidelberg, 2005). Landolt-Bornstein Volume III 43A2.
- S. B. Hendricks, *Die Kristallstruktur von N_2O_4*, Z. Phys. **70**, 699–700 (1931), doi:10.1007/BF01340758.

Found in:

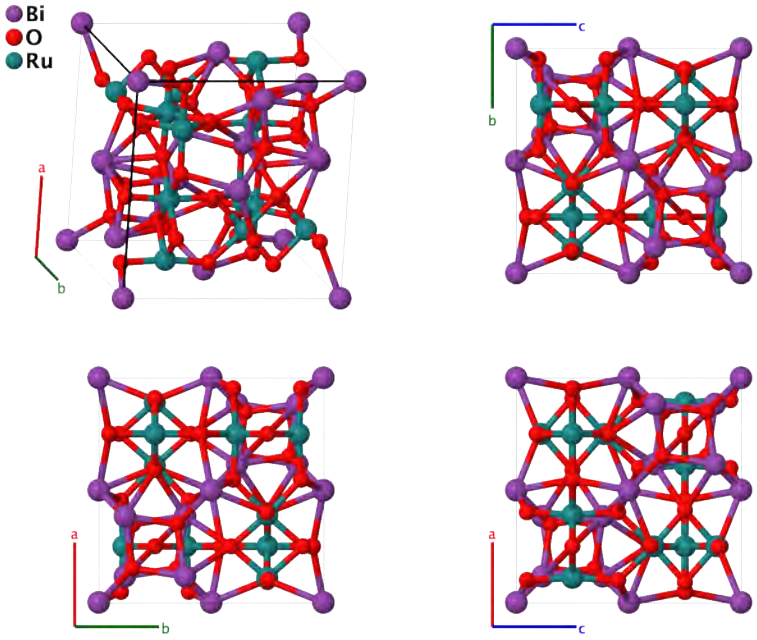
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1773
- POSCAR: pp. 1773

$Bi_3Ru_3O_{11}$ Structure: A3B11C3_cP68_201_be_efh_g

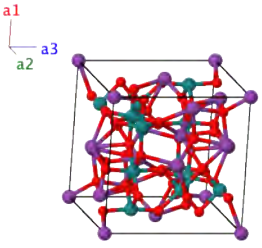
http://aflow.org/prototype-encyclopedia/A3B11C3_cP68_201_be_efh_g



Prototype	:	$Bi_3O_{11}Ru_3$
AFLOW prototype label	:	A3B11C3_cP68_201_be_efh_g
***Strukturbericht* designation**	:	None
Pearson symbol	:	cP68
Space group number	:	201
Space group symbol	:	$Pn\bar{3}$
AFLOW prototype command	:	`aflow --proto=A3B11C3_cP68_201_be_efh_g --params=`$a, x_2, x_3, x_4, x_5, x_6, y_6, z_6$

Simple Cubic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = a\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(4*b*)	Bi I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$	(4*b*)	Bi I
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(4*b*)	Bi I

$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(4*b*)	Bi I
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$x_2a\,\hat{\mathbf{x}} + x_2a\,\hat{\mathbf{y}} + x_2a\,\hat{\mathbf{z}}$	(8*e*)	Bi II
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{y}} + x_2a\,\hat{\mathbf{z}}$	(8*e*)	Bi II
$\mathbf{B}_7$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + x_2\,\mathbf{a}_2 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + x_2a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{z}}$	(8*e*)	Bi II
$\mathbf{B}_8$	=	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_3$	=	$x_2a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{z}}$	(8*e*)	Bi II
$\mathbf{B}_9$	=	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 - x_2\,\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}} - x_2a\,\hat{\mathbf{y}} - x_2a\,\hat{\mathbf{z}}$	(8*e*)	Bi II
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_2 - x_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{y}} - x_2a\,\hat{\mathbf{z}}$	(8*e*)	Bi II
$\mathbf{B}_{11}$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 - x_2\,\mathbf{a}_2 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} - x_2a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{z}}$	(8*e*)	Bi II
$\mathbf{B}_{12}$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{z}}$	(8*e*)	Bi II
$\mathbf{B}_{13}$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{15}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{16}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{17}$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{19}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{20}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{21}$	=	$x_4\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(12*f*)	O II
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(12*f*)	O II
$\mathbf{B}_{23}$	=	$\frac{1}{4}\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(12*f*)	O II
$\mathbf{B}_{24}$	=	$\frac{1}{4}\,\mathbf{a}_1 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(12*f*)	O II
$\mathbf{B}_{25}$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	(12*f*)	O II
$\mathbf{B}_{26}$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{z}}$	(12*f*)	O II
$\mathbf{B}_{27}$	=	$-x_4\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(12*f*)	O II
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(12*f*)	O II
$\mathbf{B}_{29}$	=	$\frac{3}{4}\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(12*f*)	O II
$\mathbf{B}_{30}$	=	$\frac{3}{4}\,\mathbf{a}_1 + \left(\frac{1}{2} + x_4\right)\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(12*f*)	O II
$\mathbf{B}_{31}$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} - x_4a\,\hat{\mathbf{z}}$	(12*f*)	O II
$\mathbf{B}_{32}$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} + x_4\right)\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{z}}$	(12*f*)	O II
$\mathbf{B}_{33}$	=	$x_5\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(12*g*)	Ru
$\mathbf{B}_{34}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(12*g*)	Ru
$\mathbf{B}_{35}$	=	$\frac{1}{4}\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + x_5a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(12*g*)	Ru
$\mathbf{B}_{36}$	=	$\frac{1}{4}\,\mathbf{a}_1 + \left(\frac{1}{2} - x_5\right)\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(12*g*)	Ru
$\mathbf{B}_{37}$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + x_5a\,\hat{\mathbf{z}}$	(12*g*)	Ru
$\mathbf{B}_{38}$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - x_5\right)\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{z}}$	(12*g*)	Ru
$\mathbf{B}_{39}$	=	$-x_5\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(12*g*)	Ru

$\mathbf{B}_{40}$	=	$\left(\frac{1}{2}+x_5\right)\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+\frac{3}{4}\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{x}}+\frac{1}{4}a\,\hat{\mathbf{y}}+\frac{3}{4}a\,\hat{\mathbf{z}}$	(12*g*)	Ru
$\mathbf{B}_{41}$	=	$\frac{3}{4}\mathbf{a}_1-x_5\,\mathbf{a}_2+\frac{1}{4}\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}}-x_5a\,\hat{\mathbf{y}}+\frac{1}{4}a\,\hat{\mathbf{z}}$	(12*g*)	Ru
$\mathbf{B}_{42}$	=	$\frac{3}{4}\mathbf{a}_1+\left(\frac{1}{2}+x_5\right)\mathbf{a}_2+\frac{1}{4}\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{y}}+\frac{1}{4}a\,\hat{\mathbf{z}}$	(12*g*)	Ru
$\mathbf{B}_{43}$	=	$\frac{1}{4}\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2-x_5\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}}+\frac{3}{4}a\,\hat{\mathbf{y}}-x_5a\,\hat{\mathbf{z}}$	(12*g*)	Ru
$\mathbf{B}_{44}$	=	$\frac{1}{4}\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2+\left(\frac{1}{2}+x_5\right)\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}}+\frac{3}{4}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{z}}$	(12*g*)	Ru
$\mathbf{B}_{45}$	=	$x_6\,\mathbf{a}_1+y_6\,\mathbf{a}_2+z_6\,\mathbf{a}_3$	=	$x_6a\,\hat{\mathbf{x}}+y_6a\,\hat{\mathbf{y}}+z_6a\,\hat{\mathbf{z}}$	(24*h*)	O III
$\mathbf{B}_{46}$	=	$\left(\frac{1}{2}-x_6\right)\mathbf{a}_1+\left(\frac{1}{2}-y_6\right)\mathbf{a}_2+z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_6\right)a\,\hat{\mathbf{y}}+z_6a\,\hat{\mathbf{z}}$	(24*h*)	O III
$\mathbf{B}_{47}$	=	$\left(\frac{1}{2}-x_6\right)\mathbf{a}_1+y_6\,\mathbf{a}_2+\left(\frac{1}{2}-z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{x}}+y_6a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_6\right)a\,\hat{\mathbf{z}}$	(24*h*)	O III
$\mathbf{B}_{48}$	=	$x_6\,\mathbf{a}_1+\left(\frac{1}{2}-y_6\right)\mathbf{a}_2+\left(\frac{1}{2}-z_6\right)\mathbf{a}_3$	=	$x_6a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_6\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_6\right)a\,\hat{\mathbf{z}}$	(24*h*)	O III
$\mathbf{B}_{49}$	=	$z_6\,\mathbf{a}_1+x_6\,\mathbf{a}_2+y_6\,\mathbf{a}_3$	=	$z_6a\,\hat{\mathbf{x}}+x_6a\,\hat{\mathbf{y}}+y_6a\,\hat{\mathbf{z}}$	(24*h*)	O III
$\mathbf{B}_{50}$	=	$z_6\,\mathbf{a}_1+\left(\frac{1}{2}-x_6\right)\mathbf{a}_2+\left(\frac{1}{2}-y_6\right)\mathbf{a}_3$	=	$z_6a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-y_6\right)a\,\hat{\mathbf{z}}$	(24*h*)	O III
$\mathbf{B}_{51}$	=	$\left(\frac{1}{2}-z_6\right)\mathbf{a}_1+\left(\frac{1}{2}-x_6\right)\mathbf{a}_2+y_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-z_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{y}}+y_6a\,\hat{\mathbf{z}}$	(24*h*)	O III
$\mathbf{B}_{52}$	=	$\left(\frac{1}{2}-z_6\right)\mathbf{a}_1+x_6\,\mathbf{a}_2+\left(\frac{1}{2}-y_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-z_6\right)a\,\hat{\mathbf{x}}+x_6a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-y_6\right)a\,\hat{\mathbf{z}}$	(24*h*)	O III
$\mathbf{B}_{53}$	=	$y_6\,\mathbf{a}_1+z_6\,\mathbf{a}_2+x_6\,\mathbf{a}_3$	=	$y_6a\,\hat{\mathbf{x}}+z_6a\,\hat{\mathbf{y}}+x_6a\,\hat{\mathbf{z}}$	(24*h*)	O III
$\mathbf{B}_{54}$	=	$\left(\frac{1}{2}-y_6\right)\mathbf{a}_1+z_6\,\mathbf{a}_2+\left(\frac{1}{2}-x_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_6\right)a\,\hat{\mathbf{x}}+z_6a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{z}}$	(24*h*)	O III
$\mathbf{B}_{55}$	=	$y_6\,\mathbf{a}_1+\left(\frac{1}{2}-z_6\right)\mathbf{a}_2+\left(\frac{1}{2}-x_6\right)\mathbf{a}_3$	=	$y_6a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-z_6\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{z}}$	(24*h*)	O III
$\mathbf{B}_{56}$	=	$\left(\frac{1}{2}-y_6\right)\mathbf{a}_1+\left(\frac{1}{2}-z_6\right)\mathbf{a}_2+x_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-z_6\right)a\,\hat{\mathbf{y}}+x_6a\,\hat{\mathbf{z}}$	(24*h*)	O III
$\mathbf{B}_{57}$	=	$-x_6\,\mathbf{a}_1-y_6\,\mathbf{a}_2-z_6\,\mathbf{a}_3$	=	$-x_6a\,\hat{\mathbf{x}}-y_6a\,\hat{\mathbf{y}}-z_6a\,\hat{\mathbf{z}}$	(24*h*)	O III
$\mathbf{B}_{58}$	=	$\left(\frac{1}{2}+x_6\right)\mathbf{a}_1+\left(\frac{1}{2}+y_6\right)\mathbf{a}_2-z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_6\right)a\,\hat{\mathbf{y}}-z_6a\,\hat{\mathbf{z}}$	(24*h*)	O III
$\mathbf{B}_{59}$	=	$\left(\frac{1}{2}+x_6\right)\mathbf{a}_1-y_6\,\mathbf{a}_2+\left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{x}}-y_6a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_6\right)a\,\hat{\mathbf{z}}$	(24*h*)	O III
$\mathbf{B}_{60}$	=	$-x_6\,\mathbf{a}_1+\left(\frac{1}{2}+y_6\right)\mathbf{a}_2+\left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	=	$-x_6a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_6\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_6\right)a\,\hat{\mathbf{z}}$	(24*h*)	O III
$\mathbf{B}_{61}$	=	$-z_6\,\mathbf{a}_1-x_6\,\mathbf{a}_2-y_6\,\mathbf{a}_3$	=	$-z_6a\,\hat{\mathbf{x}}-x_6a\,\hat{\mathbf{y}}-y_6a\,\hat{\mathbf{z}}$	(24*h*)	O III
$\mathbf{B}_{62}$	=	$-z_6\,\mathbf{a}_1+\left(\frac{1}{2}+x_6\right)\mathbf{a}_2+\left(\frac{1}{2}+y_6\right)\mathbf{a}_3$	=	$-z_6a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+y_6\right)a\,\hat{\mathbf{z}}$	(24*h*)	O III
$\mathbf{B}_{63}$	=	$\left(\frac{1}{2}+z_6\right)\mathbf{a}_1+\left(\frac{1}{2}+x_6\right)\mathbf{a}_2-y_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+z_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{y}}-y_6a\,\hat{\mathbf{z}}$	(24*h*)	O III
$\mathbf{B}_{64}$	=	$\left(\frac{1}{2}+z_6\right)\mathbf{a}_1-x_6\,\mathbf{a}_2+\left(\frac{1}{2}+y_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+z_6\right)a\,\hat{\mathbf{x}}-x_6a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+y_6\right)a\,\hat{\mathbf{z}}$	(24*h*)	O III
$\mathbf{B}_{65}$	=	$-y_6\,\mathbf{a}_1-z_6\,\mathbf{a}_2-x_6\,\mathbf{a}_3$	=	$-y_6a\,\hat{\mathbf{x}}-z_6a\,\hat{\mathbf{y}}-x_6a\,\hat{\mathbf{z}}$	(24*h*)	O III
$\mathbf{B}_{66}$	=	$\left(\frac{1}{2}+y_6\right)\mathbf{a}_1-z_6\,\mathbf{a}_2+\left(\frac{1}{2}+x_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_6\right)a\,\hat{\mathbf{x}}-z_6a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{z}}$	(24*h*)	O III
$\mathbf{B}_{67}$	=	$-y_6\,\mathbf{a}_1+\left(\frac{1}{2}+z_6\right)\mathbf{a}_2+\left(\frac{1}{2}+x_6\right)\mathbf{a}_3$	=	$-y_6a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+z_6\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{z}}$	(24*h*)	O III
$\mathbf{B}_{68}$	=	$\left(\frac{1}{2}+y_6\right)\mathbf{a}_1+\left(\frac{1}{2}+z_6\right)\mathbf{a}_2-x_6\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+z_6\right)a\,\hat{\mathbf{y}}-x_6a\,\hat{\mathbf{z}}$	(24*h*)	O III

References:

- F. Abraham, D. Thomas, and G. Nowogrocki, *Structure cristalline de $Bi_3Ru_3O_{11}$*, Bull. Soc. Fr. Mineral. Cristallogr. **98**, 25–29 (1975), doi:10.3406/bulmi.1975.6954.

Found in:

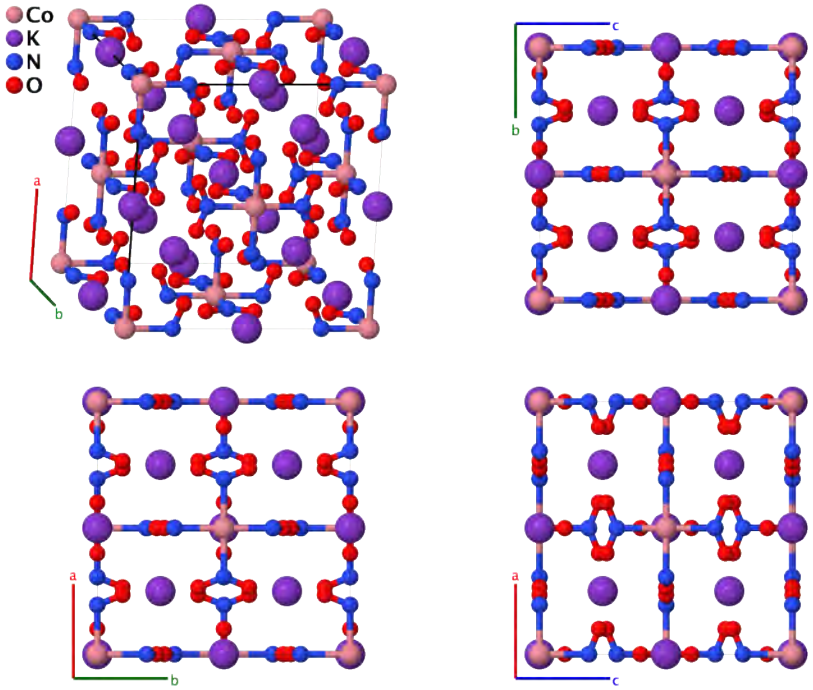
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1774
- POSCAR: pp. 1774

$K_3Co(NO_2)_6$ ($J2_4$) Structure: AB3C6D12_cF88_202_a_bc_e_h

http://aflow.org/prototype-encyclopedia/AB3C6D12_cF88_202_a_bc_e_h



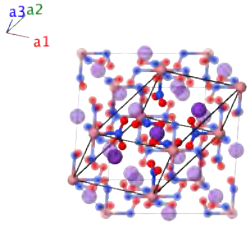
Prototype	:	$CoK_3N_6O_{12}$
AFLOW prototype label	:	AB3C6D12_cF88_202_a_bc_e_h
***Strukturbericht* designation**	:	$J2_4$
Pearson symbol	:	cF88
Space group number	:	202
Space group symbol	:	$Fm\bar{3}$
AFLOW prototype command	:	`aflow --proto=AB3C6D12_cF88_202_a_bc_e_h --params=`a, x_4, y_5, z_5

Other compounds with this structure

- $(NH_4)_2AgBi(NO_2)_6$, $(NH_4)_2LiBi(NO_2)_6$, $(NH_4)_2NaBi(NO_2)_6$, $(NH_4)_2NaCo(NO_2)_6$, $(NH_4)_2NaRh(NO_2)_6$, $(NH_4)_3Co(NO_2)_6$, $Cs_2AgBi(NO_2)_6$, $Cs_2LiBi(NO_2)_6$, $Cs_2NaBi(NO_2)_6$, $Cs_3Bi(NO_2)_6$, $K_2LiBi(NO_2)_6$, $K_2NaBi(NO_2)_6$, $K_2NaCo(NO_2)_6$, $K_2PbCu(NO_2)_6$, $K_3Ca(NO_2)_6$, $Rb_2AgBi(NO_2)_6$, $Rb_2NaBi(NO_2)_6$, $Tl_2AgBi(NO_2)_6$, $Tl_2LiBi(NO_2)_6$, and $Tl_2NaBi(NO_2)_6$

Face-centered Cubic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{y}} + \frac{1}{2}\,a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{1}{2}\,a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{1}{2}\,a\,\hat{\mathbf{y}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(4*a*)	Co
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(4*b*)	K I
$\mathbf{B}_3$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(8*c*)	K II
$\mathbf{B}_4$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(8*c*)	K II
$\mathbf{B}_5$	=	$-x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}}$	(24*e*)	N
$\mathbf{B}_6$	=	$x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}}$	(24*e*)	N
$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{y}}$	(24*e*)	N
$\mathbf{B}_8$	=	$-x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{y}}$	(24*e*)	N
$\mathbf{B}_9$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{z}}$	(24*e*)	N
$\mathbf{B}_{10}$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{z}}$	(24*e*)	N
$\mathbf{B}_{11}$	=	$(y_5 + z_5)\,\mathbf{a}_1 + (-y_5 + z_5)\,\mathbf{a}_2 + (y_5 - z_5)\,\mathbf{a}_3$	=	$y_5 a\,\hat{\mathbf{y}} + z_5 a\,\hat{\mathbf{z}}$	(48*h*)	O
$\mathbf{B}_{12}$	=	$(-y_5 + z_5)\,\mathbf{a}_1 + (y_5 + z_5)\,\mathbf{a}_2 + (-y_5 - z_5)\,\mathbf{a}_3$	=	$-y_5 a\,\hat{\mathbf{y}} + z_5 a\,\hat{\mathbf{z}}$	(48*h*)	O
$\mathbf{B}_{13}$	=	$(y_5 - z_5)\,\mathbf{a}_1 + (-y_5 - z_5)\,\mathbf{a}_2 + (y_5 + z_5)\,\mathbf{a}_3$	=	$y_5 a\,\hat{\mathbf{y}} - z_5 a\,\hat{\mathbf{z}}$	(48*h*)	O
$\mathbf{B}_{14}$	=	$(-y_5 - z_5)\,\mathbf{a}_1 + (y_5 - z_5)\,\mathbf{a}_2 + (-y_5 + z_5)\,\mathbf{a}_3$	=	$-y_5 a\,\hat{\mathbf{y}} - z_5 a\,\hat{\mathbf{z}}$	(48*h*)	O
$\mathbf{B}_{15}$	=	$(y_5 - z_5)\,\mathbf{a}_1 + (y_5 + z_5)\,\mathbf{a}_2 + (-y_5 + z_5)\,\mathbf{a}_3$	=	$z_5 a\,\hat{\mathbf{x}} + y_5 a\,\hat{\mathbf{z}}$	(48*h*)	O
$\mathbf{B}_{16}$	=	$(-y_5 - z_5)\,\mathbf{a}_1 + (-y_5 + z_5)\,\mathbf{a}_2 + (y_5 + z_5)\,\mathbf{a}_3$	=	$z_5 a\,\hat{\mathbf{x}} - y_5 a\,\hat{\mathbf{z}}$	(48*h*)	O
$\mathbf{B}_{17}$	=	$(y_5 + z_5)\,\mathbf{a}_1 + (y_5 - z_5)\,\mathbf{a}_2 + (-y_5 - z_5)\,\mathbf{a}_3$	=	$-z_5 a\,\hat{\mathbf{x}} + y_5 a\,\hat{\mathbf{z}}$	(48*h*)	O
$\mathbf{B}_{18}$	=	$(-y_5 + z_5)\,\mathbf{a}_1 + (-y_5 - z_5)\,\mathbf{a}_2 + (y_5 - z_5)\,\mathbf{a}_3$	=	$-z_5 a\,\hat{\mathbf{x}} - y_5 a\,\hat{\mathbf{z}}$	(48*h*)	O
$\mathbf{B}_{19}$	=	$(-y_5 + z_5)\,\mathbf{a}_1 + (y_5 - z_5)\,\mathbf{a}_2 + (y_5 + z_5)\,\mathbf{a}_3$	=	$y_5 a\,\hat{\mathbf{x}} + z_5 a\,\hat{\mathbf{y}}$	(48*h*)	O
$\mathbf{B}_{20}$	=	$(y_5 + z_5)\,\mathbf{a}_1 + (-y_5 - z_5)\,\mathbf{a}_2 + (-y_5 + z_5)\,\mathbf{a}_3$	=	$-y_5 a\,\hat{\mathbf{x}} + z_5 a\,\hat{\mathbf{y}}$	(48*h*)	O
$\mathbf{B}_{21}$	=	$(-y_5 - z_5)\,\mathbf{a}_1 + (y_5 + z_5)\,\mathbf{a}_2 + (y_5 - z_5)\,\mathbf{a}_3$	=	$y_5 a\,\hat{\mathbf{x}} - z_5 a\,\hat{\mathbf{y}}$	(48*h*)	O
$\mathbf{B}_{22}$	=	$(y_5 - z_5)\,\mathbf{a}_1 + (-y_5 + z_5)\,\mathbf{a}_2 + (-y_5 - z_5)\,\mathbf{a}_3$	=	$-y_5 a\,\hat{\mathbf{x}} - z_5 a\,\hat{\mathbf{y}}$	(48*h*)	O

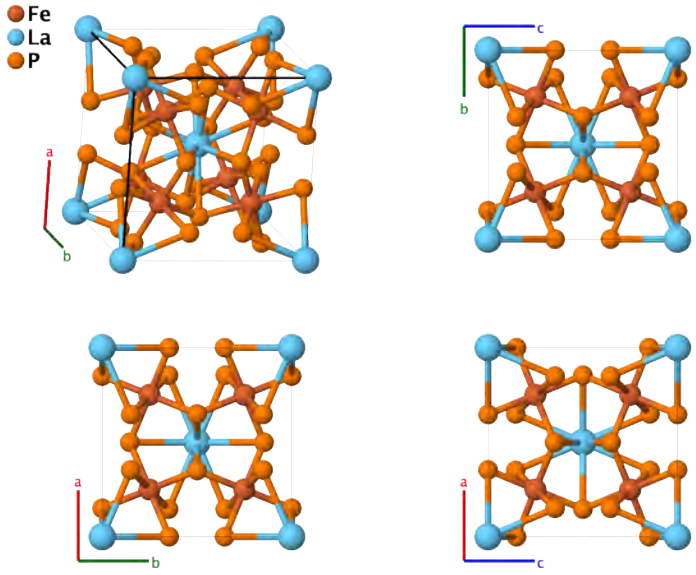
References:
- M. van Driel and H. J. Verweel, *Über die Struktur der Tripelnitrite*, Zeitschrift für Kristallographie - Crystalline Materials **95**, 308–314 (1936), doi:10.1524/zkri.1936.95.1.308.

Found in:
- C. Gottfried, ed., *Strukturbericht Band IV 1936* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1938).

Geometry files:
- CIF: pp. 1774
- POSCAR: pp. 1775

$LaFe_4P_{12}$ Structure: A4BC12_cI34_204_c_a_g

http://aflow.org/prototype-encyclopedia/A4BC12_cI34_204_c_a_g



Prototype	:	Fe_4LaP_{12}
AFLOW prototype label	:	A4BC12_cI34_204_c_a_g
***Strukturbericht* designation**	:	None
Pearson symbol	:	cI34
Space group number	:	204
Space group symbol	:	$Im\bar{3}$
AFLOW prototype command	:	`aflow --proto=A4BC12_cI34_204_c_a_g --params=`a, y_3, z_3

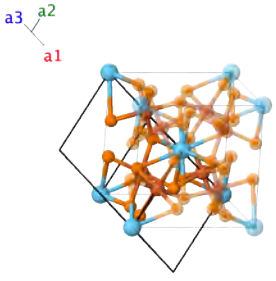
Other compounds with this structure

- $CeFe_4P_{12}$, $EuFe_4P_{12}$, $NdFe_4P_{12}$, $PrFe_4P_{12}$, $SmFe_4P_{12}$, $CeRu_4P_{12}$, $EuRu_4P_{12}$, $LaRu_4P_{12}$, $NdRu_4P_{12}$, $PrRu_4P_{12}$, $SmRu_4P_{12}$, $CeOs_4P_{12}$, $LaOs_4P_{12}$, $NdOs_4P_{12}$, $PrOs_4P_{12}$, $SmOs_4P_{12}$, $CeFe_4Sb_{12}$, $LaFe_4Sb_{12}$, $PrFe_4Sb_{12}$, $SmFe_4Sb_{12}$, $CeRu_4Sb_{12}$, $EuRu_4Sb_{12}$, $LaRu_4Sb_{12}$, $NdRu_4Sb_{12}$, $PrRu_4Sb_{12}$, $SmRu_4Sb_{12}$, $CeOs_4Sb_{12}$, $EuOs_4Sb_{12}$, $LaOs_4Sb_{12}$, $NdOs_4Sb_{12}$, $PrOs_4Sb_{12}$, and $SmOs_4Sb_{12}$

- This is a filled skutterudite ($CoAs_3$, $D0_2$) structure. Removing the rare-earth atom from the ($2a$) position yields the skutterudite structure.

Body-centered Cubic primitive vectors:

$$\mathbf{a}_1 = -\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}a\,\hat{\mathbf{x}} - \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} - \tfrac{1}{2}a\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($2a$)	La
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	($8c$)	Fe
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} - \frac{1}{4}a\,\hat{\mathbf{z}}$	($8c$)	Fe
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} - \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	($8c$)	Fe
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$-\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	($8c$)	Fe
$\mathbf{B}_6$	=	$(y_3 + z_3)\,\mathbf{a}_1 + z_3\,\mathbf{a}_2 + y_3\,\mathbf{a}_3$	=	$y_3 a\,\hat{\mathbf{y}} + z_3 a\,\hat{\mathbf{z}}$	($24g$)	P
$\mathbf{B}_7$	=	$(-y_3 + z_3)\,\mathbf{a}_1 + z_3\,\mathbf{a}_2 - y_3\,\mathbf{a}_3$	=	$-y_3 a\,\hat{\mathbf{y}} + z_3 a\,\hat{\mathbf{z}}$	($24g$)	P
$\mathbf{B}_8$	=	$(y_3 - z_3)\,\mathbf{a}_1 - z_3\,\mathbf{a}_2 + y_3\,\mathbf{a}_3$	=	$y_3 a\,\hat{\mathbf{y}} - z_3 a\,\hat{\mathbf{z}}$	($24g$)	P
$\mathbf{B}_9$	=	$(-y_3 - z_3)\,\mathbf{a}_1 - z_3\,\mathbf{a}_2 - y_3\,\mathbf{a}_3$	=	$-y_3 a\,\hat{\mathbf{y}} - z_3 a\,\hat{\mathbf{z}}$	($24g$)	P
$\mathbf{B}_{10}$	=	$y_3\,\mathbf{a}_1 + (y_3 + z_3)\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$z_3 a\,\hat{\mathbf{x}} + y_3 a\,\hat{\mathbf{z}}$	($24g$)	P
$\mathbf{B}_{11}$	=	$-y_3\,\mathbf{a}_1 + (-y_3 + z_3)\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$z_3 a\,\hat{\mathbf{x}} - y_3 a\,\hat{\mathbf{z}}$	($24g$)	P
$\mathbf{B}_{12}$	=	$y_3\,\mathbf{a}_1 + (y_3 - z_3)\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-z_3 a\,\hat{\mathbf{x}} + y_3 a\,\hat{\mathbf{z}}$	($24g$)	P
$\mathbf{B}_{13}$	=	$-y_3\,\mathbf{a}_1 + (-y_3 - z_3)\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-z_3 a\,\hat{\mathbf{x}} - y_3 a\,\hat{\mathbf{z}}$	($24g$)	P
$\mathbf{B}_{14}$	=	$z_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + (y_3 + z_3)\,\mathbf{a}_3$	=	$y_3 a\,\hat{\mathbf{x}} + z_3 a\,\hat{\mathbf{y}}$	($24g$)	P
$\mathbf{B}_{15}$	=	$z_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + (-y_3 + z_3)\,\mathbf{a}_3$	=	$-y_3 a\,\hat{\mathbf{x}} + z_3 a\,\hat{\mathbf{y}}$	($24g$)	P
$\mathbf{B}_{16}$	=	$-z_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + (y_3 - z_3)\,\mathbf{a}_3$	=	$y_3 a\,\hat{\mathbf{x}} - z_3 a\,\hat{\mathbf{y}}$	($24g$)	P
$\mathbf{B}_{17}$	=	$-z_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + (-y_3 - z_3)\,\mathbf{a}_3$	=	$-y_3 a\,\hat{\mathbf{x}} - z_3 a\,\hat{\mathbf{y}}$	($24g$)	P

References:

- W. Jeitschko and D. Braun, *$LaFe_4P_{12}$ with filled $CoAs_3$-type structure and isotypic lanthanoid-transition metal polyphosphides*, Acta Crystallogr. Sect. B Struct. Sci. **33**, 3401–3406 (1977), doi:10.1107/S056774087701108X.

Found in:

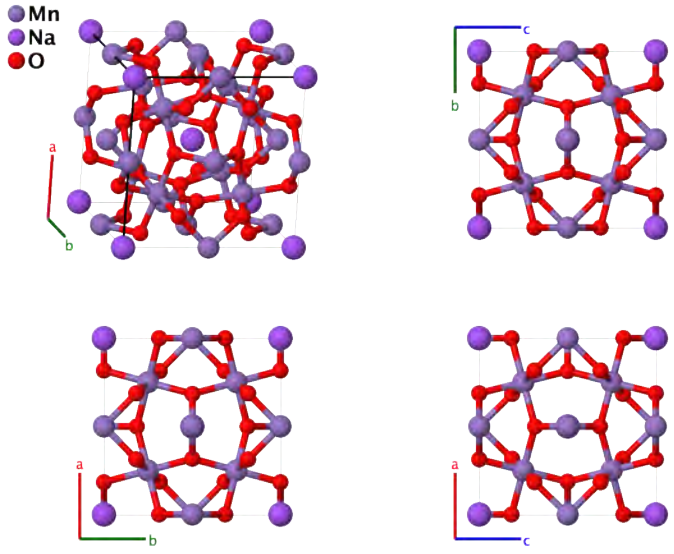
- C. R. Rotundu, *Novel Heavy Fermion Behavior in Praseodymium-based Materials: Experimental Study of $PrOs_4Sb_{12}$*, http://ufdc.ufl.edu/UFE0017620/00001/1j (2007). Ph. D. Thesis, University of Florida.
- D. J. Braun and W. Jeitschko, *Preparation and structural investigations of antimonides with the $LaFe_4P_{12}$ structure*, J. Less-Common Met. **72**, 147–156 (1980), doi:10.1016/0022-5088(80)90260-X.

Geometry files:

- CIF: pp. 1775
- POSCAR: pp. 1776

$NaMn_7O_{12}$ Structure: A7BC12_cI40_204_bc_a_g

http://aflow.org/prototype-encyclopedia/A7BC12_cI40_204_bc_a_g



Prototype	:	Mn_7NaO_{12}
AFLOW prototype label	:	A7BC12_cI40_204_bc_a_g
***Strukturbericht* designation**	:	None
Pearson symbol	:	cI40
Space group number	:	204
Space group symbol	:	$Im\bar{3}$
AFLOW prototype command	:	`aflow --proto=A7BC12_cI40_204_bc_a_g --params=`a, y_4, z_4

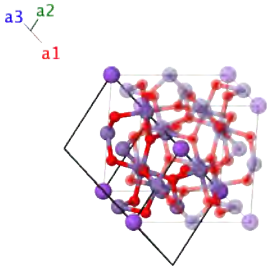
Other compounds with this structure

- $CaCu_3Ge_4O_{12}$ and $CaCu_3Mn_4O_{12}$

- This is a double perovskite structure, and is only stable above 3 Gbar and above room temperature, but it is metastable under ambient conditions (Gilioli, 2005ab). The actual composition of the measured sample is $Na_{0.95}Mn_{7.05}O_{12}$, with the excess manganese displacing some of the sodium atoms on the ($2a$) site.

Body-centered Cubic primitive vectors:

$$\mathbf{a}_1 = -\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} - \frac{1}{2}a\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(2*a*)	Na
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	(6*b*)	Mn I
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}}$	(6*b*)	Mn I
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{z}}$	(6*b*)	Mn I
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(8*c*)	Mn II
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} - \frac{1}{4}a\,\hat{\mathbf{z}}$	(8*c*)	Mn II
$\mathbf{B}_7$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} - \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(8*c*)	Mn II
$\mathbf{B}_8$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$-\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(8*c*)	Mn II
$\mathbf{B}_9$	=	$(y_4 + z_4)\,\mathbf{a}_1 + z_4\,\mathbf{a}_2 + y_4\,\mathbf{a}_3$	=	$y_4 a\,\hat{\mathbf{y}} + z_4 a\,\hat{\mathbf{z}}$	(24*g*)	O
$\mathbf{B}_{10}$	=	$(-y_4 + z_4)\,\mathbf{a}_1 + z_4\,\mathbf{a}_2 - y_4\,\mathbf{a}_3$	=	$-y_4 a\,\hat{\mathbf{y}} + z_4 a\,\hat{\mathbf{z}}$	(24*g*)	O
$\mathbf{B}_{11}$	=	$(y_4 - z_4)\,\mathbf{a}_1 - z_4\,\mathbf{a}_2 + y_4\,\mathbf{a}_3$	=	$y_4 a\,\hat{\mathbf{y}} - z_4 a\,\hat{\mathbf{z}}$	(24*g*)	O
$\mathbf{B}_{12}$	=	$(-y_4 - z_4)\,\mathbf{a}_1 - z_4\,\mathbf{a}_2 - y_4\,\mathbf{a}_3$	=	$-y_4 a\,\hat{\mathbf{y}} - z_4 a\,\hat{\mathbf{z}}$	(24*g*)	O
$\mathbf{B}_{13}$	=	$y_4\,\mathbf{a}_1 + (y_4 + z_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$z_4 a\,\hat{\mathbf{x}} + y_4 a\,\hat{\mathbf{z}}$	(24*g*)	O
$\mathbf{B}_{14}$	=	$-y_4\,\mathbf{a}_1 + (-y_4 + z_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$z_4 a\,\hat{\mathbf{x}} - y_4 a\,\hat{\mathbf{z}}$	(24*g*)	O
$\mathbf{B}_{15}$	=	$y_4\,\mathbf{a}_1 + (y_4 - z_4)\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-z_4 a\,\hat{\mathbf{x}} + y_4 a\,\hat{\mathbf{z}}$	(24*g*)	O
$\mathbf{B}_{16}$	=	$-y_4\,\mathbf{a}_1 + (-y_4 - z_4)\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-z_4 a\,\hat{\mathbf{x}} - y_4 a\,\hat{\mathbf{z}}$	(24*g*)	O
$\mathbf{B}_{17}$	=	$z_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + (y_4 + z_4)\,\mathbf{a}_3$	=	$y_4 a\,\hat{\mathbf{x}} + z_4 a\,\hat{\mathbf{y}}$	(24*g*)	O
$\mathbf{B}_{18}$	=	$z_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + (-y_4 + z_4)\,\mathbf{a}_3$	=	$-y_4 a\,\hat{\mathbf{x}} + z_4 a\,\hat{\mathbf{y}}$	(24*g*)	O
$\mathbf{B}_{19}$	=	$-z_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + (y_4 - z_4)\,\mathbf{a}_3$	=	$y_4 a\,\hat{\mathbf{x}} - z_4 a\,\hat{\mathbf{y}}$	(24*g*)	O
$\mathbf{B}_{20}$	=	$-z_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 + (-y_4 - z_4)\,\mathbf{a}_3$	=	$-y_4 a\,\hat{\mathbf{x}} - z_4 a\,\hat{\mathbf{y}}$	(24*g*)	O

References:

- E. Gilioli, G. Calestani, F. Licci, A. Gauzzi, F. Bolzoni, A. Prodi, and M. Marezio, *P – T phase diagram and single crystal structural refinement of* $NaMn_7O_{12}$, Solid State Sci. **7**, 746–752 (2005), doi:10.1016/j.solidstatesciences.2004.11.020.

Found in:

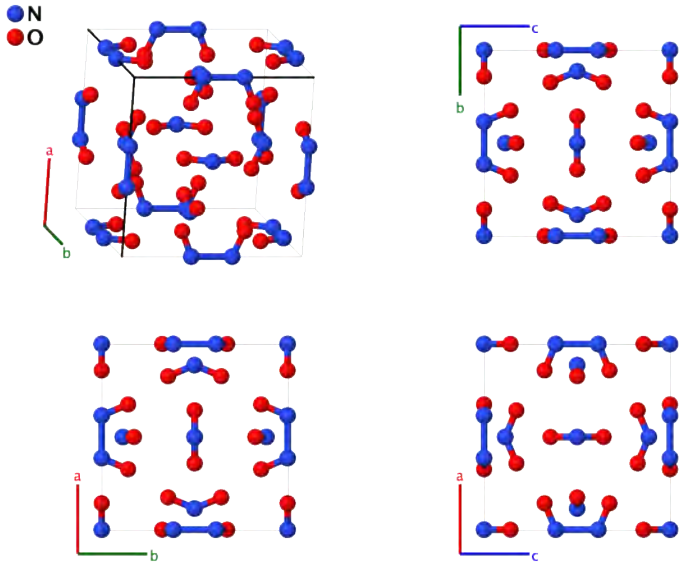
- E. Gilioli, F. Licci, G. Calestani, A. Prodi, A. Gauzzi, and G. Salviati, *Crystal growth and structural refinement of* $NaMn_7O_{12}$, Cryst. Res. Technol. **40**, 1072–1075 (2005), doi:10.1002/crat.200410489.

Geometry files:

- CIF: pp. 1776
- POSCAR: pp. 1777

NO_2 (Modern, *C*26) Structure: AB2_cI36_204_d_g

http://aflow.org/prototype-encyclopedia/AB2_cI36_204_d_g

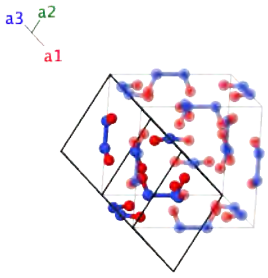


Prototype	:	NO_2
AFLOW prototype label	:	AB2_cI36_204_d_g
***Strukturbericht* designation**	:	*C*26
Pearson symbol	:	cI36
Space group number	:	204
Space group symbol	:	$Im\bar{3}$
AFLOW prototype command	:	`aflow --proto=AB2_cI36_204_d_g --params=`a, x_1, y_2, z_2

- (Hermann, 1937) listed two possible structures for the low temperature solid cubic phase of NO_2, which were given *Strukturbericht* designations $C26_a$ and $C26_b$, the only structures with Roman subscripts in the original series.
- $C26_a$ (AB2_cI36_199_b_c) was set in space group $I2_13$ #199. Hermann noted that this structure has a very short distance (1.88 Å) between oxygen atoms on different NO_2 molecules, and that this structure does not form the $(NO_2)_2$ aggregate molecule found in the $C26_b$ structure, making "making this proposed structure very unlikely."
- Recognizing this, (Hendricks, 1931) suggested that NO_2 was actually in space group $I23$ #197. (Hermann, 1997) gave this structure the $C26_b$ designation, but noted that based on Hendricks's atomic positions the space group was actually $Im\bar{3}$ #204.
- The modern accepted structure for NO_2 (AB2_cI36_204_d_g) is set in space group $Im\bar{3}$ #204, confirming Hendricks. We follow (Villars, 2005) and give this the *C*26 designation.
- We used the experimental data for this phase collected at 20 K.

Body-centered Cubic primitive vectors:

$$\mathbf{a}_1 = -\tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{y}} + \tfrac{1}{2}\,a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2}\,a\,\hat{\mathbf{y}} + \tfrac{1}{2}\,a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{y}} - \tfrac{1}{2}\,a\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_2 + x_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}}$	$(12d)$	N
$\mathbf{B}_2$	=	$-x_1\,\mathbf{a}_2 - x_1\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}}$	$(12d)$	N
$\mathbf{B}_3$	=	$x_1\,\mathbf{a}_1 + x_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{y}}$	$(12d)$	N
$\mathbf{B}_4$	=	$-x_1\,\mathbf{a}_1 - x_1\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{y}}$	$(12d)$	N
$\mathbf{B}_5$	=	$x_1\,\mathbf{a}_1 + x_1\,\mathbf{a}_2$	=	$x_1 a\,\hat{\mathbf{z}}$	$(12d)$	N
$\mathbf{B}_6$	=	$-x_1\,\mathbf{a}_1 - x_1\,\mathbf{a}_2$	=	$-x_1 a\,\hat{\mathbf{z}}$	$(12d)$	N
$\mathbf{B}_7$	=	$(y_2 + z_2)\,\mathbf{a}_1 + z_2\,\mathbf{a}_2 + y_2\,\mathbf{a}_3$	=	$y_2 a\,\hat{\mathbf{y}} + z_2 a\,\hat{\mathbf{z}}$	$(24g)$	O
$\mathbf{B}_8$	=	$(-y_2 + z_2)\,\mathbf{a}_1 + z_2\,\mathbf{a}_2 - y_2\,\mathbf{a}_3$	=	$-y_2 a\,\hat{\mathbf{y}} + z_2 a\,\hat{\mathbf{z}}$	$(24g)$	O
$\mathbf{B}_9$	=	$(y_2 - z_2)\,\mathbf{a}_1 - z_2\,\mathbf{a}_2 + y_2\,\mathbf{a}_3$	=	$y_2 a\,\hat{\mathbf{y}} - z_2 a\,\hat{\mathbf{z}}$	$(24g)$	O
$\mathbf{B}_{10}$	=	$(-y_2 - z_2)\,\mathbf{a}_1 - z_2\,\mathbf{a}_2 - y_2\,\mathbf{a}_3$	=	$-y_2 a\,\hat{\mathbf{y}} - z_2 a\,\hat{\mathbf{z}}$	$(24g)$	O
$\mathbf{B}_{11}$	=	$y_2\,\mathbf{a}_1 + (y_2 + z_2)\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$z_2 a\,\hat{\mathbf{x}} + y_2 a\,\hat{\mathbf{z}}$	$(24g)$	O
$\mathbf{B}_{12}$	=	$-y_2\,\mathbf{a}_1 + (-y_2 + z_2)\,\mathbf{a}_2 + z_2\,\mathbf{a}_3$	=	$z_2 a\,\hat{\mathbf{x}} - y_2 a\,\hat{\mathbf{z}}$	$(24g)$	O
$\mathbf{B}_{13}$	=	$y_2\,\mathbf{a}_1 + (y_2 - z_2)\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$-z_2 a\,\hat{\mathbf{x}} + y_2 a\,\hat{\mathbf{z}}$	$(24g)$	O
$\mathbf{B}_{14}$	=	$-y_2\,\mathbf{a}_1 + (-y_2 - z_2)\,\mathbf{a}_2 - z_2\,\mathbf{a}_3$	=	$-z_2 a\,\hat{\mathbf{x}} - y_2 a\,\hat{\mathbf{z}}$	$(24g)$	O
$\mathbf{B}_{15}$	=	$z_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + (y_2 + z_2)\,\mathbf{a}_3$	=	$y_2 a\,\hat{\mathbf{x}} + z_2 a\,\hat{\mathbf{y}}$	$(24g)$	O
$\mathbf{B}_{16}$	=	$z_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + (-y_2 + z_2)\,\mathbf{a}_3$	=	$-y_2 a\,\hat{\mathbf{x}} + z_2 a\,\hat{\mathbf{y}}$	$(24g)$	O
$\mathbf{B}_{17}$	=	$-z_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2 + (y_2 - z_2)\,\mathbf{a}_3$	=	$y_2 a\,\hat{\mathbf{x}} - z_2 a\,\hat{\mathbf{y}}$	$(24g)$	O
$\mathbf{B}_{18}$	=	$-z_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2 + (-y_2 - z_2)\,\mathbf{a}_3$	=	$-y_2 a\,\hat{\mathbf{x}} - z_2 a\,\hat{\mathbf{y}}$	$(24g)$	O

References:
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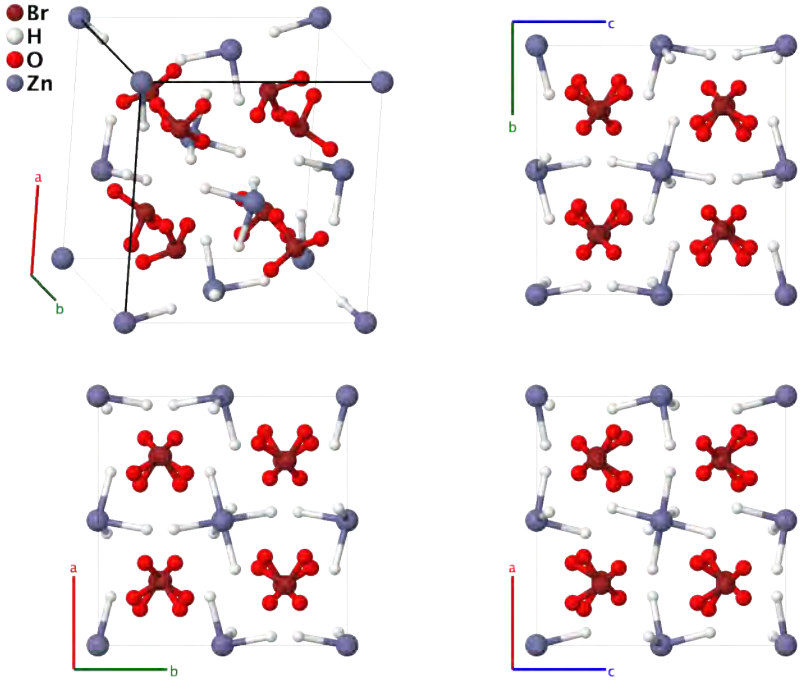
- P. Villars and K. Cenzual, eds., *Crystal Structure Data of Inorganic Compounds* (Springer-Verlag, Berlin, Heidelberg, 2005). Landolt-Bornstein Volume III 43A2.

Geometry files:

- CIF: pp. 1777
- POSCAR: pp. 1777

$Zn(BrO_3)_2{\cdot}6H_2O$ ($J1_{10}$) Structure: A2B6C6D_cP60_205_c_d_d_a

http://aflow.org/prototype-encyclopedia/A2B6C6D_cP60_205_c_d_d_a

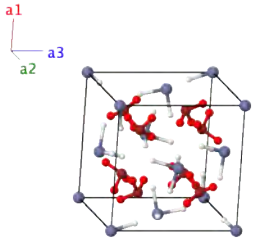


Prototype	:	$Br_2(H_2O)_6O_6Zn$
AFLOW prototype label	:	A2B6C6D_cP60_205_c_d_d_a
***Strukturbericht* designation**	:	$J1_{10}$
Pearson symbol	:	cP60
Space group number	:	205
Space group symbol	:	$Pa\bar{3}$
AFLOW prototype command	:	`aflow --proto=A2B6C6D_cP60_205_c_d_d_a --params=`$a, x_2, x_3, y_3, z_3, x_4, y_4, z_4$

- The positions of the hydrogen atoms in the water molecules were not determined, so we only provide the positions of the oxygen atom (labeled as H_2O).

Simple Cubic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = a\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(4*a*)	Zn
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(4*a*)	Zn
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(4*a*)	Zn
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$	(4*a*)	Zn
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$x_2a\,\hat{\mathbf{x}} + x_2a\,\hat{\mathbf{y}} + x_2a\,\hat{\mathbf{z}}$	(8*c*)	Br
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 - x_2\,\mathbf{a}_2 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} - x_2a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{z}}$	(8*c*)	Br
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{z}}$	(8*c*)	Br
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_2 - x_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{y}} - x_2a\,\hat{\mathbf{z}}$	(8*c*)	Br
$\mathbf{B}_9$	=	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 - x_2\,\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}} - x_2a\,\hat{\mathbf{y}} - x_2a\,\hat{\mathbf{z}}$	(8*c*)	Br
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + x_2\,\mathbf{a}_2 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + x_2a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{z}}$	(8*c*)	Br
$\mathbf{B}_{11}$	=	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_3$	=	$x_2a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{z}}$	(8*c*)	Br
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{y}} + x_2a\,\hat{\mathbf{z}}$	(8*c*)	Br
$\mathbf{B}_{13}$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + y_3a\,\hat{\mathbf{y}} + z_3a\,\hat{\mathbf{z}}$	(24*d*)	H_2O
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} - y_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	H_2O
$\mathbf{B}_{15}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	H_2O
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{y}} - z_3a\,\hat{\mathbf{z}}$	(24*d*)	H_2O
$\mathbf{B}_{17}$	=	$z_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + y_3\,\mathbf{a}_3$	=	$z_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + y_3a\,\hat{\mathbf{z}}$	(24*d*)	H_2O
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} + z_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 - y_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + z_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} - y_3a\,\hat{\mathbf{z}}$	(24*d*)	H_2O
$\mathbf{B}_{19}$	=	$\left(\frac{1}{2} - z_3\right)\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - z_3\right)a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	H_2O
$\mathbf{B}_{20}$	=	$-z_3\,\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_3$	=	$-z_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	H_2O
$\mathbf{B}_{21}$	=	$y_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$y_3a\,\hat{\mathbf{x}} + z_3a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(24*d*)	H_2O
$\mathbf{B}_{22}$	=	$-y_3\,\mathbf{a}_1 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_3$	=	$-y_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + z_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	H_2O
$\mathbf{B}_{23}$	=	$\left(\frac{1}{2} + y_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - z_3\right)a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(24*d*)	H_2O
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} - y_3\right)\mathbf{a}_1 - z_3\,\mathbf{a}_2 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{x}} - z_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	H_2O
$\mathbf{B}_{25}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - y_3a\,\hat{\mathbf{y}} - z_3a\,\hat{\mathbf{z}}$	(24*d*)	H_2O
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + y_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	H_2O
$\mathbf{B}_{27}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	H_2O
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{y}} + z_3a\,\hat{\mathbf{z}}$	(24*d*)	H_2O
$\mathbf{B}_{29}$	=	$-z_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - y_3\,\mathbf{a}_3$	=	$-z_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} - y_3a\,\hat{\mathbf{z}}$	(24*d*)	H_2O
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2} - z_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + y_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - z_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + y_3a\,\hat{\mathbf{z}}$	(24*d*)	H_2O
$\mathbf{B}_{31}$	=	$\left(\frac{1}{2} + z_3\right)\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + z_3\right)a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	H_2O
$\mathbf{B}_{32}$	=	$z_3\,\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_3$	=	$z_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	H_2O
$\mathbf{B}_{33}$	=	$-y_3\,\mathbf{a}_1 - z_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$-y_3a\,\hat{\mathbf{x}} - z_3a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(24*d*)	H_2O
$\mathbf{B}_{34}$	=	$y_3\,\mathbf{a}_1 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_3$	=	$y_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - z_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	H_2O
$\mathbf{B}_{35}$	=	$\left(\frac{1}{2} - y_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + z_3\right)a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(24*d*)	H_2O

$\mathbf{B}_{36}$	=	$\left(\frac{1}{2}+y_3\right)\mathbf{a}_1+z_3\,\mathbf{a}_2+\left(\frac{1}{2}-x_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_3\right)a\,\hat{\mathbf{x}}+z_3a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	H_2O
$\mathbf{B}_{37}$	=	$x_4\,\mathbf{a}_1+y_4\,\mathbf{a}_2+z_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}}+y_4a\,\hat{\mathbf{y}}+z_4a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{38}$	=	$\left(\frac{1}{2}-x_4\right)\mathbf{a}_1-y_4\,\mathbf{a}_2+\left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{x}}-y_4a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_4\right)a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{39}$	=	$-x_4\,\mathbf{a}_1+\left(\frac{1}{2}+y_4\right)\mathbf{a}_2+\left(\frac{1}{2}-z_4\right)\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_4\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_4\right)a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{40}$	=	$\left(\frac{1}{2}+x_4\right)\mathbf{a}_1+\left(\frac{1}{2}-y_4\right)\mathbf{a}_2-z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_4\right)a\,\hat{\mathbf{y}}-z_4a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{41}$	=	$z_4\,\mathbf{a}_1+x_4\,\mathbf{a}_2+y_4\,\mathbf{a}_3$	=	$z_4a\,\hat{\mathbf{x}}+x_4a\,\hat{\mathbf{y}}+y_4a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{42}$	=	$\left(\frac{1}{2}+z_4\right)\mathbf{a}_1+\left(\frac{1}{2}-x_4\right)\mathbf{a}_2-y_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+z_4\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{y}}-y_4a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{43}$	=	$\left(\frac{1}{2}-z_4\right)\mathbf{a}_1-x_4\,\mathbf{a}_2+\left(\frac{1}{2}+y_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-z_4\right)a\,\hat{\mathbf{x}}-x_4a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+y_4\right)a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{44}$	=	$-z_4\,\mathbf{a}_1+\left(\frac{1}{2}+x_4\right)\mathbf{a}_2+\left(\frac{1}{2}-y_4\right)\mathbf{a}_3$	=	$-z_4a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-y_4\right)a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{45}$	=	$y_4\,\mathbf{a}_1+z_4\,\mathbf{a}_2+x_4\,\mathbf{a}_3$	=	$y_4a\,\hat{\mathbf{x}}+z_4a\,\hat{\mathbf{y}}+x_4a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{46}$	=	$-y_4\,\mathbf{a}_1+\left(\frac{1}{2}+z_4\right)\mathbf{a}_2+\left(\frac{1}{2}-x_4\right)\mathbf{a}_3$	=	$-y_4a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+z_4\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{47}$	=	$\left(\frac{1}{2}+y_4\right)\mathbf{a}_1+\left(\frac{1}{2}-z_4\right)\mathbf{a}_2-x_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_4\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-z_4\right)a\,\hat{\mathbf{y}}-x_4a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{48}$	=	$\left(\frac{1}{2}-y_4\right)\mathbf{a}_1-z_4\,\mathbf{a}_2+\left(\frac{1}{2}+x_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_4\right)a\,\hat{\mathbf{x}}-z_4a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{49}$	=	$-x_4\,\mathbf{a}_1-y_4\,\mathbf{a}_2-z_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}}-y_4a\,\hat{\mathbf{y}}-z_4a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{50}$	=	$\left(\frac{1}{2}+x_4\right)\mathbf{a}_1+y_4\,\mathbf{a}_2+\left(\frac{1}{2}-z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{x}}+y_4a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_4\right)a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{51}$	=	$x_4\,\mathbf{a}_1+\left(\frac{1}{2}-y_4\right)\mathbf{a}_2+\left(\frac{1}{2}+z_4\right)\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_4\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_4\right)a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{52}$	=	$\left(\frac{1}{2}-x_4\right)\mathbf{a}_1+\left(\frac{1}{2}+y_4\right)\mathbf{a}_2+z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_4\right)a\,\hat{\mathbf{y}}+z_4a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{53}$	=	$-z_4\,\mathbf{a}_1-x_4\,\mathbf{a}_2-y_4\,\mathbf{a}_3$	=	$-z_4a\,\hat{\mathbf{x}}-x_4a\,\hat{\mathbf{y}}-y_4a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{54}$	=	$\left(\frac{1}{2}-z_4\right)\mathbf{a}_1+\left(\frac{1}{2}+x_4\right)\mathbf{a}_2+y_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-z_4\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{y}}+y_4a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{55}$	=	$\left(\frac{1}{2}+z_4\right)\mathbf{a}_1+x_4\,\mathbf{a}_2+\left(\frac{1}{2}-y_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+z_4\right)a\,\hat{\mathbf{x}}+x_4a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-y_4\right)a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{56}$	=	$z_4\,\mathbf{a}_1+\left(\frac{1}{2}-x_4\right)\mathbf{a}_2+\left(\frac{1}{2}+y_4\right)\mathbf{a}_3$	=	$z_4a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+y_4\right)a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{57}$	=	$-y_4\,\mathbf{a}_1-z_4\,\mathbf{a}_2-x_4\,\mathbf{a}_3$	=	$-y_4a\,\hat{\mathbf{x}}-z_4a\,\hat{\mathbf{y}}-x_4a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{58}$	=	$y_4\,\mathbf{a}_1+\left(\frac{1}{2}-z_4\right)\mathbf{a}_2+\left(\frac{1}{2}+x_4\right)\mathbf{a}_3$	=	$y_4a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-z_4\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{59}$	=	$\left(\frac{1}{2}-y_4\right)\mathbf{a}_1+\left(\frac{1}{2}+z_4\right)\mathbf{a}_2+x_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_4\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+z_4\right)a\,\hat{\mathbf{y}}+x_4a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{60}$	=	$\left(\frac{1}{2}+y_4\right)\mathbf{a}_1+z_4\,\mathbf{a}_2+\left(\frac{1}{2}-x_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_4\right)a\,\hat{\mathbf{x}}+z_4a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{z}}$	(24*d*)	O

References:

- S. H. Yü and C. A. Beevers, *The Crystal Structure of Zinc Bromate Hexahydrate [$Zn(BrO_3)_2\cdot 6H_2O$]*, Zeitschrift für Kristallographie - Crystalline Materials **95**, 426–434 (1936), doi:10.1524/zkri.1936.95.1.426.

Found in:

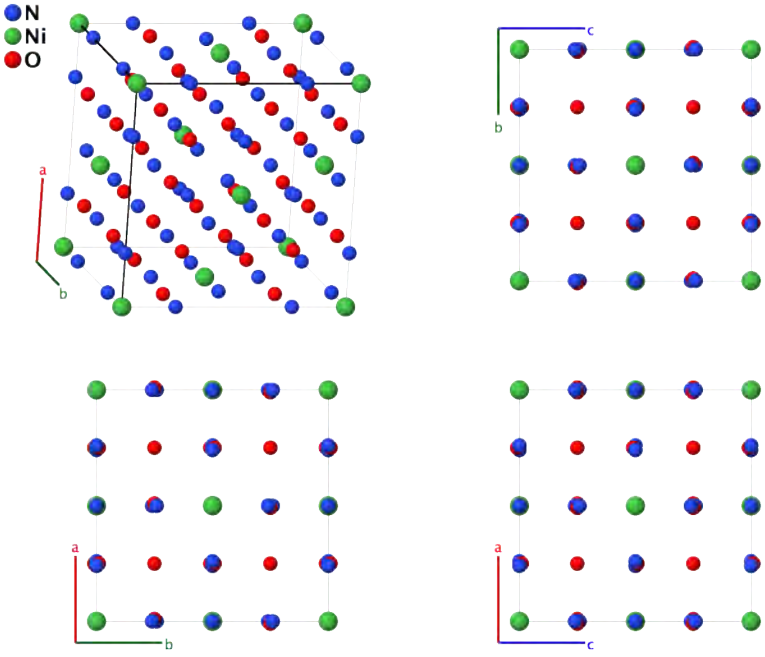
- C. Gottfried, ed., *Strukturbericht Band IV 1936* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1938).

Geometry files:

- CIF: pp. 1777
- POSCAR: pp. 1778

$H6_4$ [$Ni(NO_3)_2(NH_3)_6$] (*obsolete*) Structure: A2B6CD6_cP60_205_c_d_a_d

http://aflow.org/prototype-encyclopedia/A2B6CD6_cP60_205_c_d_a_d

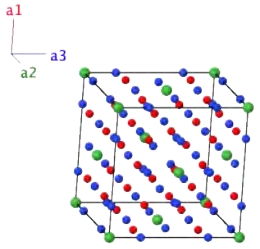


Prototype	:	$N_2(NH_3)_6NiO_6$
AFLOW prototype label	:	A2B6CD6_cP60_205_c_d_a_d
***Strukturbericht* designation**	:	$H6_4$
Pearson symbol	:	cP60
Space group number	:	205
Space group symbol	:	$Pa\bar{3}$
AFLOW prototype command	:	`aflow --proto=A2B6CD6_cP60_205_c_d_a_d --params=`$a, x_2, x_3, y_3, z_3, x_4, y_4, z_4$

- (Wyckoff, 1922) determined this approximate structure. In his paper, non-zero coordinates were $x_2 = 1/4$ for the nitrogen (8*c*) atoms, $x_3 = v$, "where v is somewhat less than 0.25," for the NH_3 (24*d*) molecules, and $x_4 = y_4 = 1/4$, $z_4 = v'$, where v' "should not deviate far from 0." If we take these coordinates as written the space group becomes $Fm\bar{3}m$ #225 rather than Wyckoff's $Pa\bar{3}$ #205, so we adjusted v and v' slightly to put the system in his space group.
- (Ewald, 1931) gave this the *Strukturbericht* designation $H61$, or $H6_1$ in later notation. (Hermann, 1937) moved it to $I1_4$ in their "list of type descriptions," but no other volume of *Strukturbericht* refers to it at all. Accordingly, we will designate this structure by its original label.
- This structure is an idealized appoximation to the true structure of $Ni(NO_3)_2(NH_3)_6$. (Bigoli, 1971) showed that the correct structure is triclinic, with space group $P\bar{1}$.
- The positions of the hydrogen atoms in the ammonia molecules were not determined, so we only provide the positions of the nitrogen atoms (labeled as NH_3).

Simple Cubic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= a\,\hat{\mathbf{x}}\\ \mathbf{a}_2 &= a\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= a\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(4*a*)	Ni
$\mathbf{B}_2$	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(4*a*)	Ni
$\mathbf{B}_3$	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(4*a*)	Ni
$\mathbf{B}_4$	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$	(4*a*)	Ni
$\mathbf{B}_5$	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	$x_2a\,\hat{\mathbf{x}} + x_2a\,\hat{\mathbf{y}} + x_2a\,\hat{\mathbf{z}}$	(8*c*)	N
$\mathbf{B}_6$	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 - x_2\,\mathbf{a}_2 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} - x_2a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{z}}$	(8*c*)	N
$\mathbf{B}_7$	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_3$	$-x_2a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{z}}$	(8*c*)	N
$\mathbf{B}_8$	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_2 - x_2\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{y}} - x_2a\,\hat{\mathbf{z}}$	(8*c*)	N
$\mathbf{B}_9$	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 - x_2\,\mathbf{a}_3$	$-x_2a\,\hat{\mathbf{x}} - x_2a\,\hat{\mathbf{y}} - x_2a\,\hat{\mathbf{z}}$	(8*c*)	N
$\mathbf{B}_{10}$	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + x_2\,\mathbf{a}_2 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + x_2a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{z}}$	(8*c*)	N
$\mathbf{B}_{11}$	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_3$	$x_2a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{z}}$	(8*c*)	N
$\mathbf{B}_{12}$	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_2 + x_2\,\mathbf{a}_3$	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{y}} + x_2a\,\hat{\mathbf{z}}$	(8*c*)	N
$\mathbf{B}_{13}$	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	$x_3a\,\hat{\mathbf{x}} + y_3a\,\hat{\mathbf{y}} + z_3a\,\hat{\mathbf{z}}$	(24*d*)	NH_3
$\mathbf{B}_{14}$	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} - y_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	NH_3
$\mathbf{B}_{15}$	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	$-x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	NH_3
$\mathbf{B}_{16}$	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 - z_3\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{y}} - z_3a\,\hat{\mathbf{z}}$	(24*d*)	NH_3
$\mathbf{B}_{17}$	$z_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + y_3\,\mathbf{a}_3$	$z_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + y_3a\,\hat{\mathbf{z}}$	(24*d*)	NH_3
$\mathbf{B}_{18}$	$\left(\frac{1}{2} + z_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 - y_3\,\mathbf{a}_3$	$\left(\frac{1}{2} + z_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} - y_3a\,\hat{\mathbf{z}}$	(24*d*)	NH_3
$\mathbf{B}_{19}$	$\left(\frac{1}{2} - z_3\right)\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_3$	$\left(\frac{1}{2} - z_3\right)a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	NH_3
$\mathbf{B}_{20}$	$-z_3\,\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_3$	$-z_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	NH_3
$\mathbf{B}_{21}$	$y_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	$y_3a\,\hat{\mathbf{x}} + z_3a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(24*d*)	NH_3
$\mathbf{B}_{22}$	$-y_3\,\mathbf{a}_1 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_3$	$-y_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + z_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	NH_3
$\mathbf{B}_{23}$	$\left(\frac{1}{2} + y_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_2 - x_3\,\mathbf{a}_3$	$\left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - z_3\right)a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(24*d*)	NH_3
$\mathbf{B}_{24}$	$\left(\frac{1}{2} - y_3\right)\mathbf{a}_1 - z_3\,\mathbf{a}_2 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_3$	$\left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{x}} - z_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	NH_3
$\mathbf{B}_{25}$	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	$-x_3a\,\hat{\mathbf{x}} - y_3a\,\hat{\mathbf{y}} - z_3a\,\hat{\mathbf{z}}$	(24*d*)	NH_3
$\mathbf{B}_{26}$	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + y_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	NH_3
$\mathbf{B}_{27}$	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	$x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	NH_3

$$
\begin{aligned}
\mathbf{B}_{28} &= \left(\tfrac{1}{2}-x_3\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_3\right)\mathbf{a}_2+z_3\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_3\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_3\right)a\,\hat{\mathbf{y}}+z_3a\,\hat{\mathbf{z}} && (24d) && \mathrm{NH_3}\\
\mathbf{B}_{29} &= -z_3\,\mathbf{a}_1-x_3\,\mathbf{a}_2-y_3\,\mathbf{a}_3 &&= -z_3a\,\hat{\mathbf{x}}-x_3a\,\hat{\mathbf{y}}-y_3a\,\hat{\mathbf{z}} && (24d) && \mathrm{NH_3}\\
\mathbf{B}_{30} &= \left(\tfrac{1}{2}-z_3\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_3\right)\mathbf{a}_2+y_3\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-z_3\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_3\right)a\,\hat{\mathbf{y}}+y_3a\,\hat{\mathbf{z}} && (24d) && \mathrm{NH_3}\\
\mathbf{B}_{31} &= \left(\tfrac{1}{2}+z_3\right)\mathbf{a}_1+x_3\,\mathbf{a}_2+\left(\tfrac{1}{2}-y_3\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+z_3\right)a\,\hat{\mathbf{x}}+x_3a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-y_3\right)a\,\hat{\mathbf{z}} && (24d) && \mathrm{NH_3}\\
\mathbf{B}_{32} &= z_3\,\mathbf{a}_1+\left(\tfrac{1}{2}-x_3\right)\mathbf{a}_2+\left(\tfrac{1}{2}+y_3\right)\mathbf{a}_3 &&= z_3a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_3\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+y_3\right)a\,\hat{\mathbf{z}} && (24d) && \mathrm{NH_3}\\
\mathbf{B}_{33} &= -y_3\,\mathbf{a}_1-z_3\,\mathbf{a}_2-x_3\,\mathbf{a}_3 &&= -y_3a\,\hat{\mathbf{x}}-z_3a\,\hat{\mathbf{y}}-x_3a\,\hat{\mathbf{z}} && (24d) && \mathrm{NH_3}\\
\mathbf{B}_{34} &= y_3\,\mathbf{a}_1+\left(\tfrac{1}{2}-z_3\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_3\right)\mathbf{a}_3 &&= y_3a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-z_3\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+x_3\right)a\,\hat{\mathbf{z}} && (24d) && \mathrm{NH_3}\\
\mathbf{B}_{35} &= \left(\tfrac{1}{2}-y_3\right)\mathbf{a}_1+\left(\tfrac{1}{2}+z_3\right)\mathbf{a}_2+x_3\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-y_3\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+z_3\right)a\,\hat{\mathbf{y}}+x_3a\,\hat{\mathbf{z}} && (24d) && \mathrm{NH_3}\\
\mathbf{B}_{36} &= \left(\tfrac{1}{2}+y_3\right)\mathbf{a}_1+z_3\,\mathbf{a}_2+\left(\tfrac{1}{2}-x_3\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+y_3\right)a\,\hat{\mathbf{x}}+z_3a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-x_3\right)a\,\hat{\mathbf{z}} && (24d) && \mathrm{NH_3}\\
\mathbf{B}_{37} &= x_4\,\mathbf{a}_1+y_4\,\mathbf{a}_2+z_4\,\mathbf{a}_3 &&= x_4a\,\hat{\mathbf{x}}+y_4a\,\hat{\mathbf{y}}+z_4a\,\hat{\mathbf{z}} && (24d) && \mathrm{O}\\
\mathbf{B}_{38} &= \left(\tfrac{1}{2}-x_4\right)\mathbf{a}_1-y_4\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_4\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_4\right)a\,\hat{\mathbf{x}}-y_4a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_4\right)a\,\hat{\mathbf{z}} && (24d) && \mathrm{O}\\
\mathbf{B}_{39} &= -x_4\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_4\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_4\right)\mathbf{a}_3 &&= -x_4a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_4\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_4\right)a\,\hat{\mathbf{z}} && (24d) && \mathrm{O}\\
\mathbf{B}_{40} &= \left(\tfrac{1}{2}+x_4\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_4\right)\mathbf{a}_2-z_4\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_4\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_4\right)a\,\hat{\mathbf{y}}-z_4a\,\hat{\mathbf{z}} && (24d) && \mathrm{O}\\
\mathbf{B}_{41} &= z_4\,\mathbf{a}_1+x_4\,\mathbf{a}_2+y_4\,\mathbf{a}_3 &&= z_4a\,\hat{\mathbf{x}}+x_4a\,\hat{\mathbf{y}}+y_4a\,\hat{\mathbf{z}} && (24d) && \mathrm{O}\\
\mathbf{B}_{42} &= \left(\tfrac{1}{2}+z_4\right)\mathbf{a}_1+\left(\tfrac{1}{2}-x_4\right)\mathbf{a}_2-y_4\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+z_4\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_4\right)a\,\hat{\mathbf{y}}-y_4a\,\hat{\mathbf{z}} && (24d) && \mathrm{O}\\
\mathbf{B}_{43} &= \left(\tfrac{1}{2}-z_4\right)\mathbf{a}_1-x_4\,\mathbf{a}_2+\left(\tfrac{1}{2}+y_4\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-z_4\right)a\,\hat{\mathbf{x}}-x_4a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+y_4\right)a\,\hat{\mathbf{z}} && (24d) && \mathrm{O}\\
\mathbf{B}_{44} &= -z_4\,\mathbf{a}_1+\left(\tfrac{1}{2}+x_4\right)\mathbf{a}_2+\left(\tfrac{1}{2}-y_4\right)\mathbf{a}_3 &&= -z_4a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_4\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-y_4\right)a\,\hat{\mathbf{z}} && (24d) && \mathrm{O}\\
\mathbf{B}_{45} &= y_4\,\mathbf{a}_1+z_4\,\mathbf{a}_2+x_4\,\mathbf{a}_3 &&= y_4a\,\hat{\mathbf{x}}+z_4a\,\hat{\mathbf{y}}+x_4a\,\hat{\mathbf{z}} && (24d) && \mathrm{O}\\
\mathbf{B}_{46} &= -y_4\,\mathbf{a}_1+\left(\tfrac{1}{2}+z_4\right)\mathbf{a}_2+\left(\tfrac{1}{2}-x_4\right)\mathbf{a}_3 &&= -y_4a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+z_4\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-x_4\right)a\,\hat{\mathbf{z}} && (24d) && \mathrm{O}\\
\mathbf{B}_{47} &= \left(\tfrac{1}{2}+y_4\right)\mathbf{a}_1+\left(\tfrac{1}{2}-z_4\right)\mathbf{a}_2-x_4\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+y_4\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-z_4\right)a\,\hat{\mathbf{y}}-x_4a\,\hat{\mathbf{z}} && (24d) && \mathrm{O}\\
\mathbf{B}_{48} &= \left(\tfrac{1}{2}-y_4\right)\mathbf{a}_1-z_4\,\mathbf{a}_2+\left(\tfrac{1}{2}+x_4\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-y_4\right)a\,\hat{\mathbf{x}}-z_4a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+x_4\right)a\,\hat{\mathbf{z}} && (24d) && \mathrm{O}\\
\mathbf{B}_{49} &= -x_4\,\mathbf{a}_1-y_4\,\mathbf{a}_2-z_4\,\mathbf{a}_3 &&= -x_4a\,\hat{\mathbf{x}}-y_4a\,\hat{\mathbf{y}}-z_4a\,\hat{\mathbf{z}} && (24d) && \mathrm{O}\\
\mathbf{B}_{50} &= \left(\tfrac{1}{2}+x_4\right)\mathbf{a}_1+y_4\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_4\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_4\right)a\,\hat{\mathbf{x}}+y_4a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_4\right)a\,\hat{\mathbf{z}} && (24d) && \mathrm{O}\\
\mathbf{B}_{51} &= x_4\,\mathbf{a}_1+\left(\tfrac{1}{2}-y_4\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_4\right)\mathbf{a}_3 &&= x_4a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_4\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_4\right)a\,\hat{\mathbf{z}} && (24d) && \mathrm{O}\\
\mathbf{B}_{52} &= \left(\tfrac{1}{2}-x_4\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_4\right)\mathbf{a}_2+z_4\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_4\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_4\right)a\,\hat{\mathbf{y}}+z_4a\,\hat{\mathbf{z}} && (24d) && \mathrm{O}\\
\mathbf{B}_{53} &= -z_4\,\mathbf{a}_1-x_4\,\mathbf{a}_2-y_4\,\mathbf{a}_3 &&= -z_4a\,\hat{\mathbf{x}}-x_4a\,\hat{\mathbf{y}}-y_4a\,\hat{\mathbf{z}} && (24d) && \mathrm{O}\\
\mathbf{B}_{54} &= \left(\tfrac{1}{2}-z_4\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_4\right)\mathbf{a}_2+y_4\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-z_4\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_4\right)a\,\hat{\mathbf{y}}+y_4a\,\hat{\mathbf{z}} && (24d) && \mathrm{O}\\
\mathbf{B}_{55} &= \left(\tfrac{1}{2}+z_4\right)\mathbf{a}_1+x_4\,\mathbf{a}_2+\left(\tfrac{1}{2}-y_4\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+z_4\right)a\,\hat{\mathbf{x}}+x_4a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-y_4\right)a\,\hat{\mathbf{z}} && (24d) && \mathrm{O}\\
\mathbf{B}_{56} &= z_4\,\mathbf{a}_1+\left(\tfrac{1}{2}-x_4\right)\mathbf{a}_2+\left(\tfrac{1}{2}+y_4\right)\mathbf{a}_3 &&= z_4a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_4\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+y_4\right)a\,\hat{\mathbf{z}} && (24d) && \mathrm{O}\\
\mathbf{B}_{57} &= -y_4\,\mathbf{a}_1-z_4\,\mathbf{a}_2-x_4\,\mathbf{a}_3 &&= -y_4a\,\hat{\mathbf{x}}-z_4a\,\hat{\mathbf{y}}-x_4a\,\hat{\mathbf{z}} && (24d) && \mathrm{O}\\
\mathbf{B}_{58} &= y_4\,\mathbf{a}_1+\left(\tfrac{1}{2}-z_4\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_4\right)\mathbf{a}_3 &&= y_4a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-z_4\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+x_4\right)a\,\hat{\mathbf{z}} && (24d) && \mathrm{O}\\
\mathbf{B}_{59} &= \left(\tfrac{1}{2}-y_4\right)\mathbf{a}_1+\left(\tfrac{1}{2}+z_4\right)\mathbf{a}_2+x_4\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-y_4\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+z_4\right)a\,\hat{\mathbf{y}}+x_4a\,\hat{\mathbf{z}} && (24d) && \mathrm{O}\\
\mathbf{B}_{60} &= \left(\tfrac{1}{2}+y_4\right)\mathbf{a}_1+z_4\,\mathbf{a}_2+\left(\tfrac{1}{2}-x_4\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+y_4\right)a\,\hat{\mathbf{x}}+z_4a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-x_4\right)a\,\hat{\mathbf{z}} && (24d) && \mathrm{O}
\end{aligned}
$$

References:

- R. W. G. Wyckoff, *The Composition and Crystal Structure of Nickel Nitrate Hexammoniate*, J. Am. Chem. Soc. **44**, 1260–1266 (1922), doi:10.1021/ja01427a010.

- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).
- F. Bigoli, A. Braibanti, A. Tiripicchio, and M. T. Camellini, *The crystal structures of nitrates of divalent hexaaquocations. III. Hexaaquonickel nitrate*, Acta Crystallogr. Sect. B Struct. Sci. **27**, 1427–1434 (1971), doi:10.1107/S0567740871004084.

Found in:

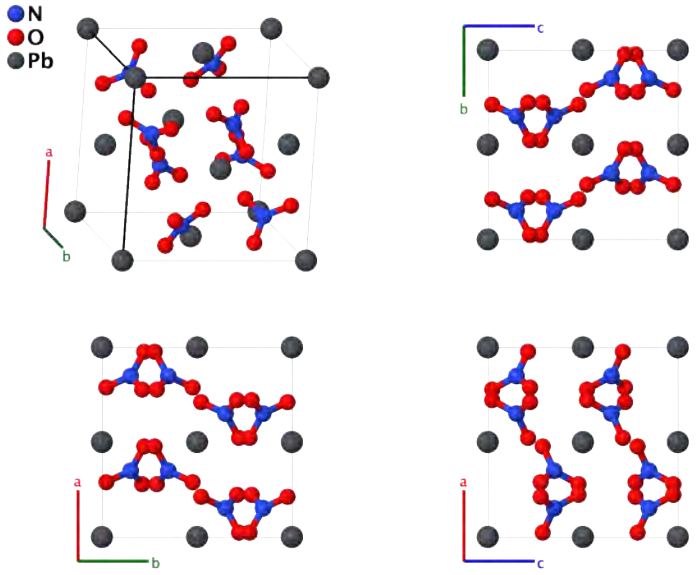
- P. P. Ewald and C. Hermann, eds., *Strukturbericht 1913-1928* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1931).

Geometry files:

- CIF: pp. 1778
- POSCAR: pp. 1779

$Pb(NO_3)_2$ ($G2_1$) Structure: A2B6C_cP36_205_c_d_a

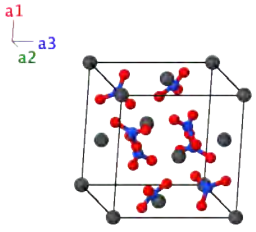
http://aflow.org/prototype-encyclopedia/A2B6C_cP36_205_c_d_a



Prototype	:	N_2O_6Pb
AFLOW prototype label	:	A2B6C_cP36_205_c_d_a
***Strukturbericht* designation**	:	$G2_1$
Pearson symbol	:	cP36
Space group number	:	205
Space group symbol	:	$Pa\bar{3}$
AFLOW prototype command	:	`aflow --proto=A2B6C_cP36_205_c_d_a --params=`a, x_2, x_3, y_3, z_3

Simple Cubic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = a\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($4a$)	Pb
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	($4a$)	Pb
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	($4a$)	Pb
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$	($4a$)	Pb

$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + x_2 a\,\hat{\mathbf{y}} + x_2 a\,\hat{\mathbf{z}}$	(8*c*)	N
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 - x_2\,\mathbf{a}_2 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} - x_2 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{z}}$	(8*c*)	N
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{z}}$	(8*c*)	N
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_2 - x_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{y}} - x_2 a\,\hat{\mathbf{z}}$	(8*c*)	N
$\mathbf{B}_9$	=	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 - x_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} - x_2 a\,\hat{\mathbf{y}} - x_2 a\,\hat{\mathbf{z}}$	(8*c*)	N
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + x_2\,\mathbf{a}_2 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + x_2 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{z}}$	(8*c*)	N
$\mathbf{B}_{11}$	=	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{z}}$	(8*c*)	N
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{y}} + x_2 a\,\hat{\mathbf{z}}$	(8*c*)	N
$\mathbf{B}_{13}$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 a\,\hat{\mathbf{y}} + z_3 a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} - y_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{15}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{y}} - z_3 a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{17}$	=	$z_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + y_3\,\mathbf{a}_3$	=	$z_3 a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}} + y_3 a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} + z_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 - y_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + z_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} - y_3 a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{19}$	=	$\left(\frac{1}{2} - z_3\right)\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - z_3\right)a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{20}$	=	$-z_3\,\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_3$	=	$-z_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{21}$	=	$y_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$y_3 a\,\hat{\mathbf{x}} + z_3 a\,\hat{\mathbf{y}} + x_3 a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{22}$	=	$-y_3\,\mathbf{a}_1 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_3$	=	$-y_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + z_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{23}$	=	$\left(\frac{1}{2} + y_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - z_3\right)a\,\hat{\mathbf{y}} - x_3 a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} - y_3\right)\mathbf{a}_1 - z_3\,\mathbf{a}_2 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{x}} - z_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{25}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 a\,\hat{\mathbf{y}} - z_3 a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + y_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{27}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{y}} + z_3 a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{29}$	=	$-z_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - y_3\,\mathbf{a}_3$	=	$-z_3 a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}} - y_3 a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2} - z_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + y_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - z_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + y_3 a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{31}$	=	$\left(\frac{1}{2} + z_3\right)\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + z_3\right)a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{32}$	=	$z_3\,\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_3$	=	$z_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{33}$	=	$-y_3\,\mathbf{a}_1 - z_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$-y_3 a\,\hat{\mathbf{x}} - z_3 a\,\hat{\mathbf{y}} - x_3 a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{34}$	=	$y_3\,\mathbf{a}_1 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_3$	=	$y_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - z_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{35}$	=	$\left(\frac{1}{2} - y_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + z_3\right)a\,\hat{\mathbf{y}} + x_3 a\,\hat{\mathbf{z}}$	(24*d*)	O
$\mathbf{B}_{36}$	=	$\left(\frac{1}{2} + y_3\right)\mathbf{a}_1 + z_3\,\mathbf{a}_2 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{x}} + z_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	O

References:

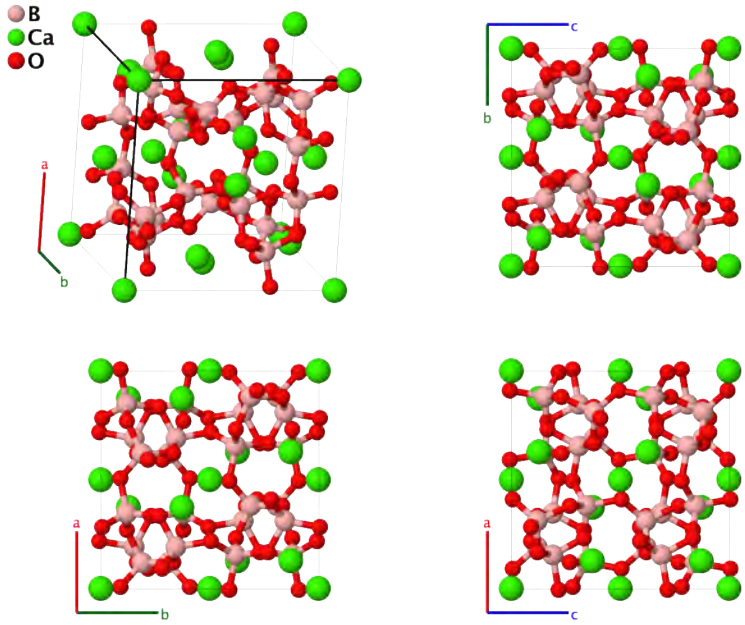
- H. Nowotny and G. Heger, *Structure Refinement of Lead Nitrate*, Acta Crystallogr. C **42**, 133–135 (1986), doi:10.1107/S0108270186097032.

Geometry files:

- CIF: pp. 1779
- POSCAR: pp. 1779

CaB_2O_4 (IV) Structure: A2BC4_cP84_205_d_ac_2d

http://aflow.org/prototype-encyclopedia/A2BC4_cP84_205_d_ac_2d

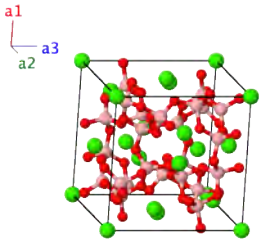


Prototype	:	B_2CaO_4
AFLOW prototype label	:	A2BC4_cP84_205_d_ac_2d
***Strukturbericht* designation**	:	None
Pearson symbol	:	cP84
Space group number	:	205
Space group symbol	:	$Pa\bar{3}$
AFLOW prototype command	:	`aflow --proto=A2BC4_cP84_205_d_ac_2d --params=`$a, x_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5$

- CaB_2O_4 exists in at least four phases (Marezio, 1969):
- I - The ground state, stable below 1.2 GPa, *Strukturbericht* $E3_2$.
- II – Orthorhombic high pressure structure, stable between 1.2 and 1.5 GPa, presumably calciborite.
- III – Orthorhombic high pressure structure, stable between 1.5 and 2.5 GPa.
- IV – Cubic high pressure structure, stable between 2.5 and 4.0 GPa (this structure).

Simple Cubic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = a\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(4*a*)	Ca I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(4*a*)	Ca I
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(4*a*)	Ca I
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$	(4*a*)	Ca I
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$x_2a\,\hat{\mathbf{x}} + x_2a\,\hat{\mathbf{y}} + x_2a\,\hat{\mathbf{z}}$	(8*c*)	Ca II
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 - x_2\,\mathbf{a}_2 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} - x_2a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{z}}$	(8*c*)	Ca II
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{z}}$	(8*c*)	Ca II
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_2 - x_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{y}} - x_2a\,\hat{\mathbf{z}}$	(8*c*)	Ca II
$\mathbf{B}_9$	=	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 - x_2\,\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}} - x_2a\,\hat{\mathbf{y}} - x_2a\,\hat{\mathbf{z}}$	(8*c*)	Ca II
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + x_2\,\mathbf{a}_2 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + x_2a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{z}}$	(8*c*)	Ca II
$\mathbf{B}_{11}$	=	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_3$	=	$x_2a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{z}}$	(8*c*)	Ca II
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{y}} + x_2a\,\hat{\mathbf{z}}$	(8*c*)	Ca II
$\mathbf{B}_{13}$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + y_3a\,\hat{\mathbf{y}} + z_3a\,\hat{\mathbf{z}}$	(24*d*)	B
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} - y_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	B
$\mathbf{B}_{15}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	B
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{y}} - z_3a\,\hat{\mathbf{z}}$	(24*d*)	B
$\mathbf{B}_{17}$	=	$z_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + y_3\,\mathbf{a}_3$	=	$z_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + y_3a\,\hat{\mathbf{z}}$	(24*d*)	B
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} + z_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 - y_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + z_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} - y_3a\,\hat{\mathbf{z}}$	(24*d*)	B
$\mathbf{B}_{19}$	=	$\left(\frac{1}{2} - z_3\right)\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - z_3\right)a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	B
$\mathbf{B}_{20}$	=	$-z_3\,\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_3$	=	$-z_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	B
$\mathbf{B}_{21}$	=	$y_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$y_3a\,\hat{\mathbf{x}} + z_3a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(24*d*)	B
$\mathbf{B}_{22}$	=	$-y_3\,\mathbf{a}_1 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_3$	=	$-y_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + z_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	B
$\mathbf{B}_{23}$	=	$\left(\frac{1}{2} + y_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - z_3\right)a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(24*d*)	B
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} - y_3\right)\mathbf{a}_1 - z_3\,\mathbf{a}_2 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{x}} - z_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	B
$\mathbf{B}_{25}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - y_3a\,\hat{\mathbf{y}} - z_3a\,\hat{\mathbf{z}}$	(24*d*)	B
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + y_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	B
$\mathbf{B}_{27}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	B
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{y}} + z_3a\,\hat{\mathbf{z}}$	(24*d*)	B
$\mathbf{B}_{29}$	=	$-z_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - y_3\,\mathbf{a}_3$	=	$-z_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} - y_3a\,\hat{\mathbf{z}}$	(24*d*)	B
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2} - z_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + y_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - z_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + y_3a\,\hat{\mathbf{z}}$	(24*d*)	B
$\mathbf{B}_{31}$	=	$\left(\frac{1}{2} + z_3\right)\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + z_3\right)a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	B
$\mathbf{B}_{32}$	=	$z_3\,\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_3$	=	$z_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	B
$\mathbf{B}_{33}$	=	$-y_3\,\mathbf{a}_1 - z_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$-y_3a\,\hat{\mathbf{x}} - z_3a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(24*d*)	B
$\mathbf{B}_{34}$	=	$y_3\,\mathbf{a}_1 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_3$	=	$y_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - z_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	B

$$
\begin{aligned}
\mathbf{B}_{35} &= \left(\tfrac{1}{2}-y_3\right)\mathbf{a}_1+\left(\tfrac{1}{2}+z_3\right)\mathbf{a}_2+x_3\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-y_3\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+z_3\right)a\,\hat{\mathbf{y}}+x_3a\,\hat{\mathbf{z}} && (24d) && \text{B}\\
\mathbf{B}_{36} &= \left(\tfrac{1}{2}+y_3\right)\mathbf{a}_1+z_3\,\mathbf{a}_2+\left(\tfrac{1}{2}-x_3\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+y_3\right)a\,\hat{\mathbf{x}}+z_3a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-x_3\right)a\,\hat{\mathbf{z}} && (24d) && \text{B}\\
\mathbf{B}_{37} &= x_4\,\mathbf{a}_1+y_4\,\mathbf{a}_2+z_4\,\mathbf{a}_3 &&= x_4a\,\hat{\mathbf{x}}+y_4a\,\hat{\mathbf{y}}+z_4a\,\hat{\mathbf{z}} && (24d) && \text{O I}\\
\mathbf{B}_{38} &= \left(\tfrac{1}{2}-x_4\right)\mathbf{a}_1-y_4\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_4\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_4\right)a\,\hat{\mathbf{x}}-y_4a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_4\right)a\,\hat{\mathbf{z}} && (24d) && \text{O I}\\
\mathbf{B}_{39} &= -x_4\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_4\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_4\right)\mathbf{a}_3 &&= -x_4a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_4\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_4\right)a\,\hat{\mathbf{z}} && (24d) && \text{O I}\\
\mathbf{B}_{40} &= \left(\tfrac{1}{2}+x_4\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_4\right)\mathbf{a}_2-z_4\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_4\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_4\right)a\,\hat{\mathbf{y}}-z_4a\,\hat{\mathbf{z}} && (24d) && \text{O I}\\
\mathbf{B}_{41} &= z_4\,\mathbf{a}_1+x_4\,\mathbf{a}_2+y_4\,\mathbf{a}_3 &&= z_4a\,\hat{\mathbf{x}}+x_4a\,\hat{\mathbf{y}}+y_4a\,\hat{\mathbf{z}} && (24d) && \text{O I}\\
\mathbf{B}_{42} &= \left(\tfrac{1}{2}+z_4\right)\mathbf{a}_1+\left(\tfrac{1}{2}-x_4\right)\mathbf{a}_2-y_4\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+z_4\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_4\right)a\,\hat{\mathbf{y}}-y_4a\,\hat{\mathbf{z}} && (24d) && \text{O I}\\
\mathbf{B}_{43} &= \left(\tfrac{1}{2}-z_4\right)\mathbf{a}_1-x_4\,\mathbf{a}_2+\left(\tfrac{1}{2}+y_4\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-z_4\right)a\,\hat{\mathbf{x}}-x_4a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+y_4\right)a\,\hat{\mathbf{z}} && (24d) && \text{O I}\\
\mathbf{B}_{44} &= -z_4\,\mathbf{a}_1+\left(\tfrac{1}{2}+x_4\right)\mathbf{a}_2+\left(\tfrac{1}{2}-y_4\right)\mathbf{a}_3 &&= -z_4a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_4\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-y_4\right)a\,\hat{\mathbf{z}} && (24d) && \text{O I}\\
\mathbf{B}_{45} &= y_4\,\mathbf{a}_1+z_4\,\mathbf{a}_2+x_4\,\mathbf{a}_3 &&= y_4a\,\hat{\mathbf{x}}+z_4a\,\hat{\mathbf{y}}+x_4a\,\hat{\mathbf{z}} && (24d) && \text{O I}\\
\mathbf{B}_{46} &= -y_4\,\mathbf{a}_1+\left(\tfrac{1}{2}+z_4\right)\mathbf{a}_2+\left(\tfrac{1}{2}-x_4\right)\mathbf{a}_3 &&= -y_4a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+z_4\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-x_4\right)a\,\hat{\mathbf{z}} && (24d) && \text{O I}\\
\mathbf{B}_{47} &= \left(\tfrac{1}{2}+y_4\right)\mathbf{a}_1+\left(\tfrac{1}{2}-z_4\right)\mathbf{a}_2-x_4\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+y_4\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-z_4\right)a\,\hat{\mathbf{y}}-x_4a\,\hat{\mathbf{z}} && (24d) && \text{O I}\\
\mathbf{B}_{48} &= \left(\tfrac{1}{2}-y_4\right)\mathbf{a}_1-z_4\,\mathbf{a}_2+\left(\tfrac{1}{2}+x_4\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-y_4\right)a\,\hat{\mathbf{x}}-z_4a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+x_4\right)a\,\hat{\mathbf{z}} && (24d) && \text{O I}\\
\mathbf{B}_{49} &= -x_4\,\mathbf{a}_1-y_4\,\mathbf{a}_2-z_4\,\mathbf{a}_3 &&= -x_4a\,\hat{\mathbf{x}}-y_4a\,\hat{\mathbf{y}}-z_4a\,\hat{\mathbf{z}} && (24d) && \text{O I}\\
\mathbf{B}_{50} &= \left(\tfrac{1}{2}+x_4\right)\mathbf{a}_1+y_4\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_4\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_4\right)a\,\hat{\mathbf{x}}+y_4a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_4\right)a\,\hat{\mathbf{z}} && (24d) && \text{O I}\\
\mathbf{B}_{51} &= x_4\,\mathbf{a}_1+\left(\tfrac{1}{2}-y_4\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_4\right)\mathbf{a}_3 &&= x_4a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_4\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_4\right)a\,\hat{\mathbf{z}} && (24d) && \text{O I}\\
\mathbf{B}_{52} &= \left(\tfrac{1}{2}-x_4\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_4\right)\mathbf{a}_2+z_4\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_4\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_4\right)a\,\hat{\mathbf{y}}+z_4a\,\hat{\mathbf{z}} && (24d) && \text{O I}\\
\mathbf{B}_{53} &= -z_4\,\mathbf{a}_1-x_4\,\mathbf{a}_2-y_4\,\mathbf{a}_3 &&= -z_4a\,\hat{\mathbf{x}}-x_4a\,\hat{\mathbf{y}}-y_4a\,\hat{\mathbf{z}} && (24d) && \text{O I}\\
\mathbf{B}_{54} &= \left(\tfrac{1}{2}-z_4\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_4\right)\mathbf{a}_2+y_4\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-z_4\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_4\right)a\,\hat{\mathbf{y}}+y_4a\,\hat{\mathbf{z}} && (24d) && \text{O I}\\
\mathbf{B}_{55} &= \left(\tfrac{1}{2}+z_4\right)\mathbf{a}_1+x_4\,\mathbf{a}_2+\left(\tfrac{1}{2}-y_4\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+z_4\right)a\,\hat{\mathbf{x}}+x_4a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-y_4\right)a\,\hat{\mathbf{z}} && (24d) && \text{O I}\\
\mathbf{B}_{56} &= z_4\,\mathbf{a}_1+\left(\tfrac{1}{2}-x_4\right)\mathbf{a}_2+\left(\tfrac{1}{2}+y_4\right)\mathbf{a}_3 &&= z_4a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_4\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+y_4\right)a\,\hat{\mathbf{z}} && (24d) && \text{O I}\\
\mathbf{B}_{57} &= -y_4\,\mathbf{a}_1-z_4\,\mathbf{a}_2-x_4\,\mathbf{a}_3 &&= -y_4a\,\hat{\mathbf{x}}-z_4a\,\hat{\mathbf{y}}-x_4a\,\hat{\mathbf{z}} && (24d) && \text{O I}\\
\mathbf{B}_{58} &= y_4\,\mathbf{a}_1+\left(\tfrac{1}{2}-z_4\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_4\right)\mathbf{a}_3 &&= y_4a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-z_4\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+x_4\right)a\,\hat{\mathbf{z}} && (24d) && \text{O I}\\
\mathbf{B}_{59} &= \left(\tfrac{1}{2}-y_4\right)\mathbf{a}_1+\left(\tfrac{1}{2}+z_4\right)\mathbf{a}_2+x_4\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-y_4\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+z_4\right)a\,\hat{\mathbf{y}}+x_4a\,\hat{\mathbf{z}} && (24d) && \text{O I}\\
\mathbf{B}_{60} &= \left(\tfrac{1}{2}+y_4\right)\mathbf{a}_1+z_4\,\mathbf{a}_2+\left(\tfrac{1}{2}-x_4\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+y_4\right)a\,\hat{\mathbf{x}}+z_4a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-x_4\right)a\,\hat{\mathbf{z}} && (24d) && \text{O I}\\
\mathbf{B}_{61} &= x_5\,\mathbf{a}_1+y_5\,\mathbf{a}_2+z_5\,\mathbf{a}_3 &&= x_5a\,\hat{\mathbf{x}}+y_5a\,\hat{\mathbf{y}}+z_5a\,\hat{\mathbf{z}} && (24d) && \text{O II}\\
\mathbf{B}_{62} &= \left(\tfrac{1}{2}-x_5\right)\mathbf{a}_1-y_5\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_5\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_5\right)a\,\hat{\mathbf{x}}-y_5a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_5\right)a\,\hat{\mathbf{z}} && (24d) && \text{O II}\\
\mathbf{B}_{63} &= -x_5\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_5\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_5\right)\mathbf{a}_3 &&= -x_5a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_5\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_5\right)a\,\hat{\mathbf{z}} && (24d) && \text{O II}\\
\mathbf{B}_{64} &= \left(\tfrac{1}{2}+x_5\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_5\right)\mathbf{a}_2-z_5\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_5\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_5\right)a\,\hat{\mathbf{y}}-z_5a\,\hat{\mathbf{z}} && (24d) && \text{O II}\\
\mathbf{B}_{65} &= z_5\,\mathbf{a}_1+x_5\,\mathbf{a}_2+y_5\,\mathbf{a}_3 &&= z_5a\,\hat{\mathbf{x}}+x_5a\,\hat{\mathbf{y}}+y_5a\,\hat{\mathbf{z}} && (24d) && \text{O II}\\
\mathbf{B}_{66} &= \left(\tfrac{1}{2}+z_5\right)\mathbf{a}_1+\left(\tfrac{1}{2}-x_5\right)\mathbf{a}_2-y_5\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+z_5\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_5\right)a\,\hat{\mathbf{y}}-y_5a\,\hat{\mathbf{z}} && (24d) && \text{O II}\\
\mathbf{B}_{67} &= \left(\tfrac{1}{2}-z_5\right)\mathbf{a}_1-x_5\,\mathbf{a}_2+\left(\tfrac{1}{2}+y_5\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-z_5\right)a\,\hat{\mathbf{x}}-x_5a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+y_5\right)a\,\hat{\mathbf{z}} && (24d) && \text{O II}\\
\mathbf{B}_{68} &= -z_5\,\mathbf{a}_1+\left(\tfrac{1}{2}+x_5\right)\mathbf{a}_2+\left(\tfrac{1}{2}-y_5\right)\mathbf{a}_3 &&= -z_5a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_5\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-y_5\right)a\,\hat{\mathbf{z}} && (24d) && \text{O II}\\
\mathbf{B}_{69} &= y_5\,\mathbf{a}_1+z_5\,\mathbf{a}_2+x_5\,\mathbf{a}_3 &&= y_5a\,\hat{\mathbf{x}}+z_5a\,\hat{\mathbf{y}}+x_5a\,\hat{\mathbf{z}} && (24d) && \text{O II}\\
\mathbf{B}_{70} &= -y_5\,\mathbf{a}_1+\left(\tfrac{1}{2}+z_5\right)\mathbf{a}_2+\left(\tfrac{1}{2}-x_5\right)\mathbf{a}_3 &&= -y_5a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+z_5\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-x_5\right)a\,\hat{\mathbf{z}} && (24d) && \text{O II}
\end{aligned}
$$

$\mathbf{B}_{71}$	=	$\left(\frac{1}{2}+y_5\right)\mathbf{a}_1+\left(\frac{1}{2}-z_5\right)\mathbf{a}_2-x_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-z_5\right)a\,\hat{\mathbf{y}}-x_5a\,\hat{\mathbf{z}}$	(24*d*)	O II
$\mathbf{B}_{72}$	=	$\left(\frac{1}{2}-y_5\right)\mathbf{a}_1-z_5\,\mathbf{a}_2+\left(\frac{1}{2}+x_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_5\right)a\,\hat{\mathbf{x}}-z_5a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{z}}$	(24*d*)	O II
$\mathbf{B}_{73}$	=	$-x_5\,\mathbf{a}_1-y_5\,\mathbf{a}_2-z_5\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}}-y_5a\,\hat{\mathbf{y}}-z_5a\,\hat{\mathbf{z}}$	(24*d*)	O II
$\mathbf{B}_{74}$	=	$\left(\frac{1}{2}+x_5\right)\mathbf{a}_1+y_5\,\mathbf{a}_2+\left(\frac{1}{2}-z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{x}}+y_5a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_5\right)a\,\hat{\mathbf{z}}$	(24*d*)	O II
$\mathbf{B}_{75}$	=	$x_5\,\mathbf{a}_1+\left(\frac{1}{2}-y_5\right)\mathbf{a}_2+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_5\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)a\,\hat{\mathbf{z}}$	(24*d*)	O II
$\mathbf{B}_{76}$	=	$\left(\frac{1}{2}-x_5\right)\mathbf{a}_1+\left(\frac{1}{2}+y_5\right)\mathbf{a}_2+z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_5\right)a\,\hat{\mathbf{y}}+z_5a\,\hat{\mathbf{z}}$	(24*d*)	O II
$\mathbf{B}_{77}$	=	$-z_5\,\mathbf{a}_1-x_5\,\mathbf{a}_2-y_5\,\mathbf{a}_3$	=	$-z_5a\,\hat{\mathbf{x}}-x_5a\,\hat{\mathbf{y}}-y_5a\,\hat{\mathbf{z}}$	(24*d*)	O II
$\mathbf{B}_{78}$	=	$\left(\frac{1}{2}-z_5\right)\mathbf{a}_1+\left(\frac{1}{2}+x_5\right)\mathbf{a}_2+y_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-z_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{y}}+y_5a\,\hat{\mathbf{z}}$	(24*d*)	O II
$\mathbf{B}_{79}$	=	$\left(\frac{1}{2}+z_5\right)\mathbf{a}_1+x_5\,\mathbf{a}_2+\left(\frac{1}{2}-y_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+z_5\right)a\,\hat{\mathbf{x}}+x_5a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-y_5\right)a\,\hat{\mathbf{z}}$	(24*d*)	O II
$\mathbf{B}_{80}$	=	$z_5\,\mathbf{a}_1+\left(\frac{1}{2}-x_5\right)\mathbf{a}_2+\left(\frac{1}{2}+y_5\right)\mathbf{a}_3$	=	$z_5a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+y_5\right)a\,\hat{\mathbf{z}}$	(24*d*)	O II
$\mathbf{B}_{81}$	=	$-y_5\,\mathbf{a}_1-z_5\,\mathbf{a}_2-x_5\,\mathbf{a}_3$	=	$-y_5a\,\hat{\mathbf{x}}-z_5a\,\hat{\mathbf{y}}-x_5a\,\hat{\mathbf{z}}$	(24*d*)	O II
$\mathbf{B}_{82}$	=	$y_5\,\mathbf{a}_1+\left(\frac{1}{2}-z_5\right)\mathbf{a}_2+\left(\frac{1}{2}+x_5\right)\mathbf{a}_3$	=	$y_5a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-z_5\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{z}}$	(24*d*)	O II
$\mathbf{B}_{83}$	=	$\left(\frac{1}{2}-y_5\right)\mathbf{a}_1+\left(\frac{1}{2}+z_5\right)\mathbf{a}_2+x_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+z_5\right)a\,\hat{\mathbf{y}}+x_5a\,\hat{\mathbf{z}}$	(24*d*)	O II
$\mathbf{B}_{84}$	=	$\left(\frac{1}{2}+y_5\right)\mathbf{a}_1+z_5\,\mathbf{a}_2+\left(\frac{1}{2}-x_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_5\right)a\,\hat{\mathbf{x}}+z_5a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{z}}$	(24*d*)	O II

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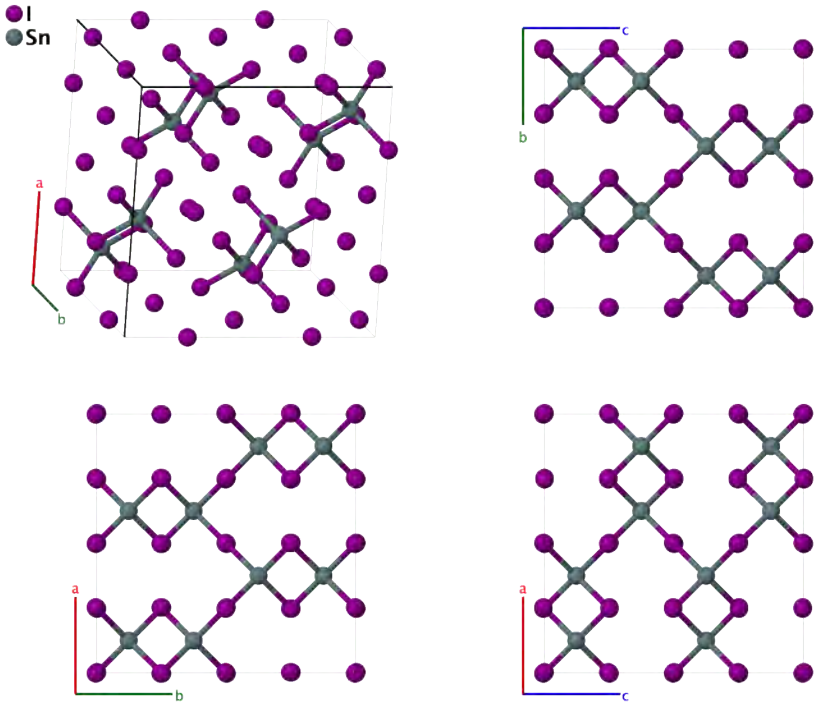
- M. Marezio, J. P. Remeika, and P. D. Dernier, *The crystal structure of the high-pressure phase CaB_2O_4(IV), and polymorphism in CaB_2O_4*, Acta Crystallogr. Sect. B Struct. Sci. **25**, 965–970 (1969), doi:10.1107/S0567740869003256.

Geometry files:

- CIF: pp. 1780
- POSCAR: pp. 1780

SnI_4 ($D1_1$) Structure: A4B_cP40_205_cd_c

http://aflow.org/prototype-encyclopedia/A4B_cP40_205_cd_c



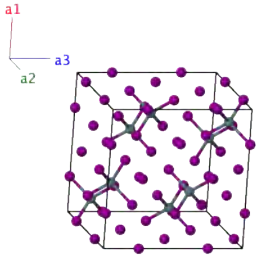
Prototype	:	I_4Sn
AFLOW prototype label	:	A4B_cP40_205_cd_c
***Strukturbericht* designation**	:	$D1_1$
Pearson symbol	:	cP40
Space group number	:	205
Space group symbol	:	$Pa\bar{3}$
AFLOW prototype command	:	`aflow --proto=A4B_cP40_205_cd_c --params=`$a, x_1, x_2, x_3, y_3, z_3$

Other compounds with this structure

- SiI_4, $TiBr_4$, TiI_4, and $Ni(CO)_4$

Simple Cubic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = a\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + x_1\,\mathbf{a}_2 + x_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + x_1 a\,\hat{\mathbf{y}} + x_1 a\,\hat{\mathbf{z}}$	(8*c*)	I I
$\mathbf{B}_2$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 - x_1\,\mathbf{a}_2 + \left(\frac{1}{2} + x_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right)a\,\hat{\mathbf{x}} - x_1 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_1\right)a\,\hat{\mathbf{z}}$	(8*c*)	I I
$\mathbf{B}_3$	=	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2} + x_1\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_1\right)\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_1\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_1\right)a\,\hat{\mathbf{z}}$	(8*c*)	I I
$\mathbf{B}_4$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_1\right)\mathbf{a}_2 - x_1\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_1\right)a\,\hat{\mathbf{y}} - x_1 a\,\hat{\mathbf{z}}$	(8*c*)	I I
$\mathbf{B}_5$	=	$-x_1\,\mathbf{a}_1 - x_1\,\mathbf{a}_2 - x_1\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} - x_1 a\,\hat{\mathbf{y}} - x_1 a\,\hat{\mathbf{z}}$	(8*c*)	I I
$\mathbf{B}_6$	=	$\left(\frac{1}{2} + x_1\right)\mathbf{a}_1 + x_1\,\mathbf{a}_2 + \left(\frac{1}{2} - x_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_1\right)a\,\hat{\mathbf{x}} + x_1 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_1\right)a\,\hat{\mathbf{z}}$	(8*c*)	I I
$\mathbf{B}_7$	=	$x_1\,\mathbf{a}_1 + \left(\frac{1}{2} - x_1\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_1\right)\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_1\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_1\right)a\,\hat{\mathbf{z}}$	(8*c*)	I I
$\mathbf{B}_8$	=	$\left(\frac{1}{2} - x_1\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_1\right)\mathbf{a}_2 + x_1\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_1\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_1\right)a\,\hat{\mathbf{y}} + x_1 a\,\hat{\mathbf{z}}$	(8*c*)	I I
$\mathbf{B}_9$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + x_2 a\,\hat{\mathbf{y}} + x_2 a\,\hat{\mathbf{z}}$	(8*c*)	Sn
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 - x_2\,\mathbf{a}_2 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} - x_2 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{z}}$	(8*c*)	Sn
$\mathbf{B}_{11}$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{z}}$	(8*c*)	Sn
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_2 - x_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{y}} - x_2 a\,\hat{\mathbf{z}}$	(8*c*)	Sn
$\mathbf{B}_{13}$	=	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 - x_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} - x_2 a\,\hat{\mathbf{y}} - x_2 a\,\hat{\mathbf{z}}$	(8*c*)	Sn
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + x_2\,\mathbf{a}_2 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + x_2 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{z}}$	(8*c*)	Sn
$\mathbf{B}_{15}$	=	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{z}}$	(8*c*)	Sn
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{y}} + x_2 a\,\hat{\mathbf{z}}$	(8*c*)	Sn
$\mathbf{B}_{17}$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + y_3 a\,\hat{\mathbf{y}} + z_3 a\,\hat{\mathbf{z}}$	(24*d*)	I II
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} - y_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	I II
$\mathbf{B}_{19}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	I II
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{y}} - z_3 a\,\hat{\mathbf{z}}$	(24*d*)	I II
$\mathbf{B}_{21}$	=	$z_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + y_3\,\mathbf{a}_3$	=	$z_3 a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}} + y_3 a\,\hat{\mathbf{z}}$	(24*d*)	I II
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} + z_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 - y_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + z_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} - y_3 a\,\hat{\mathbf{z}}$	(24*d*)	I II
$\mathbf{B}_{23}$	=	$\left(\frac{1}{2} - z_3\right)\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - z_3\right)a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	I II
$\mathbf{B}_{24}$	=	$-z_3\,\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_3$	=	$-z_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	I II
$\mathbf{B}_{25}$	=	$y_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$y_3 a\,\hat{\mathbf{x}} + z_3 a\,\hat{\mathbf{y}} + x_3 a\,\hat{\mathbf{z}}$	(24*d*)	I II
$\mathbf{B}_{26}$	=	$-y_3\,\mathbf{a}_1 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_3$	=	$-y_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + z_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	I II
$\mathbf{B}_{27}$	=	$\left(\frac{1}{2} + y_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - z_3\right)a\,\hat{\mathbf{y}} - x_3 a\,\hat{\mathbf{z}}$	(24*d*)	I II
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} - y_3\right)\mathbf{a}_1 - z_3\,\mathbf{a}_2 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{x}} - z_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	I II
$\mathbf{B}_{29}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} - y_3 a\,\hat{\mathbf{y}} - z_3 a\,\hat{\mathbf{z}}$	(24*d*)	I II
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + y_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	I II
$\mathbf{B}_{31}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	I II
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{y}} + z_3 a\,\hat{\mathbf{z}}$	(24*d*)	I II
$\mathbf{B}_{33}$	=	$-z_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - y_3\,\mathbf{a}_3$	=	$-z_3 a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}} - y_3 a\,\hat{\mathbf{z}}$	(24*d*)	I II
$\mathbf{B}_{34}$	=	$\left(\frac{1}{2} - z_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + y_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - z_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + y_3 a\,\hat{\mathbf{z}}$	(24*d*)	I II
$\mathbf{B}_{35}$	=	$\left(\frac{1}{2} + z_3\right)\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + z_3\right)a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	I II

$$\mathbf{B}_{36} = z_3\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + y_3\right)\mathbf{a}_3 = z_3 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_3\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + y_3\right) a\,\hat{\mathbf{z}} \quad (24d) \quad \text{I II}$$

$$\mathbf{B}_{37} = -y_3\,\mathbf{a}_1 - z_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3 = -y_3 a\,\hat{\mathbf{x}} - z_3 a\,\hat{\mathbf{y}} - x_3 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{I II}$$

$$\mathbf{B}_{38} = y_3\,\mathbf{a}_1 + \left(\tfrac{1}{2} - z_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_3\right)\mathbf{a}_3 = y_3 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - z_3\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_3\right) a\,\hat{\mathbf{z}} \quad (24d) \quad \text{I II}$$

$$\mathbf{B}_{39} = \left(\tfrac{1}{2} - y_3\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + z_3\right)\mathbf{a}_2 + x_3\,\mathbf{a}_3 = \left(\tfrac{1}{2} - y_3\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + z_3\right) a\,\hat{\mathbf{y}} + x_3 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{I II}$$

$$\mathbf{B}_{40} = \left(\tfrac{1}{2} + y_3\right)\mathbf{a}_1 + z_3\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_3\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + y_3\right) a\,\hat{\mathbf{x}} + z_3 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_3\right) a\,\hat{\mathbf{z}} \quad (24d) \quad \text{I II}$$

References:

- F. Meller and I. Fankuchen, *The crystal structure of tin tetraiodide*, Acta Cryst. **8**, 343–344 (1955), doi:10.1107/S0365110X55001035.

Found in:

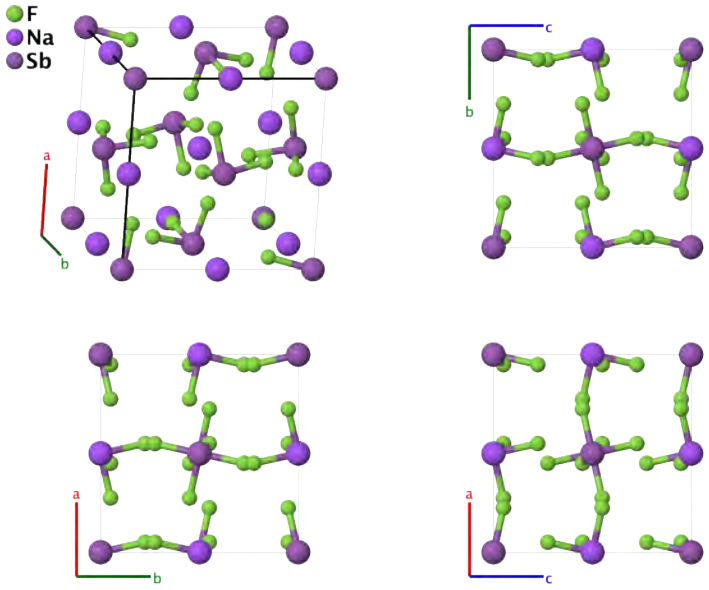
- Y. Fujii, M. Kowaka, and A. Onodera, *The pressure-induced metallic amorphous state of SnI_4. I. A novel crystal-to-amorphous transition studied by X-ray scattering*, J. Phys. C: Solid State Phys. **18**, 789–797 (1985), doi:10.1088/0022-3719/18/4/010.

Geometry files:

- CIF: pp. 1780
- POSCAR: pp. 1781

$NaSbF_6$ Structure: A6BC_cP32_205_d_b_a

http://aflow.org/prototype-encyclopedia/A6BC_cP32_205_d_b_a

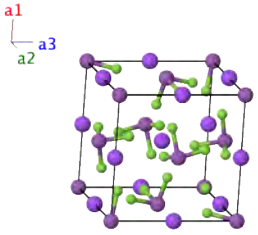


Prototype	:	F_6NaSb
AFLOW prototype label	:	A6BC_cP32_205_d_b_a
***Strukturbericht* designation**	:	None
Pearson symbol	:	cP32
Space group number	:	205
Space group symbol	:	$Pa\bar{3}$
AFLOW prototype command	:	`aflow --proto=A6BC_cP32_205_d_b_a --params=`a, x_3, y_3, z_3

- This is the pure fluorine end-point of the $NaSbF_{6-x}(OH)_x$ compounds. Although the replacement of fluoride by OH does not affect the shape of the Sb-$(F,OH)_6$ ions, it has a profound effect on the structure, as can be seen by looking at and $NaSb(OH)_6$ ($J1_{11}$) and $NaSbF_4(OH)_2$ ($J1_{12}$).

Simple Cubic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = a\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($4a$)	Sb

$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(4*a*)	Sb
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(4*a*)	Sb
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$	(4*a*)	Sb
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(4*b*)	Na
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{y}}$	(4*b*)	Na
$\mathbf{B}_7$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	(4*b*)	Na
$\mathbf{B}_8$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{z}}$	(4*b*)	Na
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + y_3a\,\hat{\mathbf{y}} + z_3a\,\hat{\mathbf{z}}$	(24*d*)	F
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} - y_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	F
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	F
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{y}} - z_3a\,\hat{\mathbf{z}}$	(24*d*)	F
$\mathbf{B}_{13}$	=	$z_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + y_3\,\mathbf{a}_3$	=	$z_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + y_3a\,\hat{\mathbf{z}}$	(24*d*)	F
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} + z_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 - y_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + z_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} - y_3a\,\hat{\mathbf{z}}$	(24*d*)	F
$\mathbf{B}_{15}$	=	$\left(\frac{1}{2} - z_3\right)\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - z_3\right)a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	F
$\mathbf{B}_{16}$	=	$-z_3\,\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_3$	=	$-z_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	F
$\mathbf{B}_{17}$	=	$y_3\,\mathbf{a}_1 + z_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$y_3a\,\hat{\mathbf{x}} + z_3a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(24*d*)	F
$\mathbf{B}_{18}$	=	$-y_3\,\mathbf{a}_1 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_3$	=	$-y_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + z_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	F
$\mathbf{B}_{19}$	=	$\left(\frac{1}{2} + y_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - z_3\right)a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(24*d*)	F
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} - y_3\right)\mathbf{a}_1 - z_3\,\mathbf{a}_2 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{x}} - z_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	F
$\mathbf{B}_{21}$	=	$-x_3\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 - z_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - y_3a\,\hat{\mathbf{y}} - z_3a\,\hat{\mathbf{z}}$	(24*d*)	F
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + y_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	F
$\mathbf{B}_{23}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	F
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + z_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{y}} + z_3a\,\hat{\mathbf{z}}$	(24*d*)	F
$\mathbf{B}_{25}$	=	$-z_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - y_3\,\mathbf{a}_3$	=	$-z_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} - y_3a\,\hat{\mathbf{z}}$	(24*d*)	F
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} - z_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + y_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - z_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + y_3a\,\hat{\mathbf{z}}$	(24*d*)	F
$\mathbf{B}_{27}$	=	$\left(\frac{1}{2} + z_3\right)\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + z_3\right)a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	F
$\mathbf{B}_{28}$	=	$z_3\,\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_3$	=	$z_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	F
$\mathbf{B}_{29}$	=	$-y_3\,\mathbf{a}_1 - z_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$-y_3a\,\hat{\mathbf{x}} - z_3a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(24*d*)	F
$\mathbf{B}_{30}$	=	$y_3\,\mathbf{a}_1 + \left(\frac{1}{2} - z_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_3$	=	$y_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - z_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	F
$\mathbf{B}_{31}$	=	$\left(\frac{1}{2} - y_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + z_3\right)\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + z_3\right)a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(24*d*)	F
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2} + y_3\right)\mathbf{a}_1 + z_3\,\mathbf{a}_2 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{x}} + z_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{z}}$	(24*d*)	F

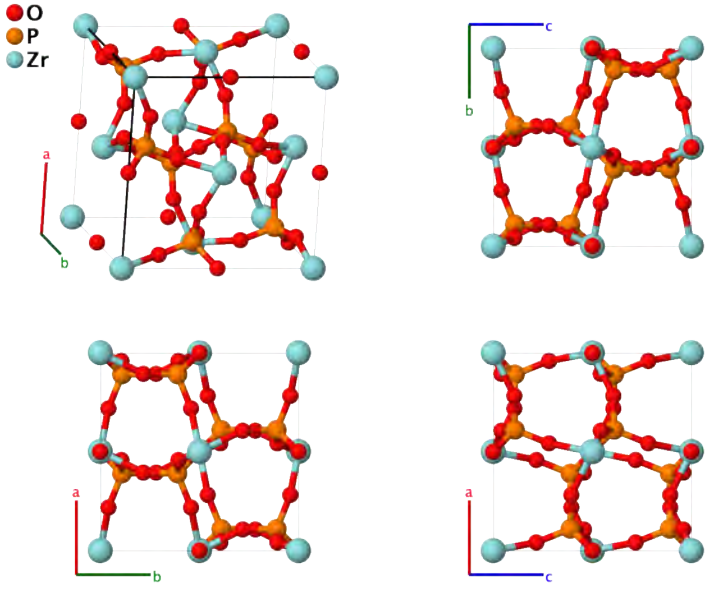
References:
- N. Schrewelius, *Röntgenuntersuchung der Verbindungen $NaSb(OH)_6$, $NaSbF_6$, $NaSbO_3$ und gleichartiger Stoffe*, Z. Anorg. Allg. Chem. **238**, 241–254 (1938), doi:10.1002/zaac.19382380209.

Found in:
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:
- CIF: pp. 1781
- POSCAR: pp. 1782

ZrP_2O_7 High-Temperature ($K6_1$) Structure: A7B2C_cP40_205_bd_c_a

http://aflow.org/prototype-encyclopedia/A7B2C_cP40_205_bd_c_a



Prototype	:	P_2O_7Zr
AFLOW prototype label	:	A7B2C_cP40_205_bd_c_a
***Strukturbericht* designation**	:	$K6_1$
Pearson symbol	:	cP40
Space group number	:	205
Space group symbol	:	$Pa\bar{3}$
AFLOW prototype command	:	`aflow --proto=A7B2C_cP40_205_bd_c_a --params=`a, x_3, x_4, y_4, z_4

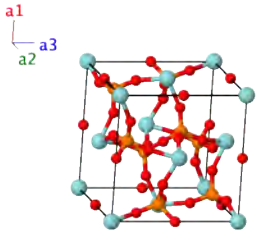
Other compounds with this structure

- AB_2O_7, (A = Si, Ge, Sn, Pb, Ti, Zr, Hf, Mo, W, Re, Ce, U, *etc.*; B = P, V, As)

- This is the high temperature form of all the structures listed. The low temperature structure depends on the composition. Below 290 °C, ZrP_2O_7 transforms to an orthorhombic structure, space group *Pbca* #61, with 136 unique crystallographic positions and 1080 atomic sites. See (Birkedal, 2006) and (Stinton, 2006) for more details.

Simple Cubic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = a\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(4*a*)	Zr
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(4*a*)	Zr
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(4*a*)	Zr
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$	(4*a*)	Zr
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(4*b*)	O I
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{y}}$	(4*b*)	O I
$\mathbf{B}_7$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	(4*b*)	O I
$\mathbf{B}_8$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{z}}$	(4*b*)	O I
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(8*c*)	P
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{z}}$	(8*c*)	P
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{z}}$	(8*c*)	P
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(8*c*)	P
$\mathbf{B}_{13}$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(8*c*)	P
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{z}}$	(8*c*)	P
$\mathbf{B}_{15}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{z}}$	(8*c*)	P
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(8*c*)	P
$\mathbf{B}_{17}$	=	$x_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + y_4a\,\hat{\mathbf{y}} + z_4a\,\hat{\mathbf{z}}$	(24*d*)	O II
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 - y_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} - y_4a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)a\,\hat{\mathbf{z}}$	(24*d*)	O II
$\mathbf{B}_{19}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)a\,\hat{\mathbf{z}}$	(24*d*)	O II
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)a\,\hat{\mathbf{y}} - z_4a\,\hat{\mathbf{z}}$	(24*d*)	O II
$\mathbf{B}_{21}$	=	$z_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + y_4\,\mathbf{a}_3$	=	$z_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + y_4a\,\hat{\mathbf{z}}$	(24*d*)	O II
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} + z_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_2 - y_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + z_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{y}} - y_4a\,\hat{\mathbf{z}}$	(24*d*)	O II
$\mathbf{B}_{23}$	=	$\left(\frac{1}{2} - z_4\right)\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - z_4\right)a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + y_4\right)a\,\hat{\mathbf{z}}$	(24*d*)	O II
$\mathbf{B}_{24}$	=	$-z_4\,\mathbf{a}_1 + \left(\frac{1}{2} + x_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_3$	=	$-z_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - y_4\right)a\,\hat{\mathbf{z}}$	(24*d*)	O II
$\mathbf{B}_{25}$	=	$y_4\,\mathbf{a}_1 + z_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$y_4a\,\hat{\mathbf{x}} + z_4a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	(24*d*)	O II
$\mathbf{B}_{26}$	=	$-y_4\,\mathbf{a}_1 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_3$	=	$-y_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + z_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{z}}$	(24*d*)	O II
$\mathbf{B}_{27}$	=	$\left(\frac{1}{2} + y_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - z_4\right)a\,\hat{\mathbf{y}} - x_4a\,\hat{\mathbf{z}}$	(24*d*)	O II
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} - y_4\right)\mathbf{a}_1 - z_4\,\mathbf{a}_2 + \left(\frac{1}{2} + x_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_4\right)a\,\hat{\mathbf{x}} - z_4a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{z}}$	(24*d*)	O II
$\mathbf{B}_{29}$	=	$-x_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} - y_4a\,\hat{\mathbf{y}} - z_4a\,\hat{\mathbf{z}}$	(24*d*)	O II
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + y_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + y_4a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)a\,\hat{\mathbf{z}}$	(24*d*)	O II
$\mathbf{B}_{31}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2} - y_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)a\,\hat{\mathbf{z}}$	(24*d*)	O II
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_4\right)a\,\hat{\mathbf{y}} + z_4a\,\hat{\mathbf{z}}$	(24*d*)	O II
$\mathbf{B}_{33}$	=	$-z_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - y_4\,\mathbf{a}_3$	=	$-z_4a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} - y_4a\,\hat{\mathbf{z}}$	(24*d*)	O II
$\mathbf{B}_{34}$	=	$\left(\frac{1}{2} - z_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_4\right)\mathbf{a}_2 + y_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - z_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{y}} + y_4a\,\hat{\mathbf{z}}$	(24*d*)	O II

$\mathbf{B}_{35}$	=	$\left(\frac{1}{2}+z_4\right)\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \left(\frac{1}{2}-y_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+z_4\right)a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-y_4\right)a\,\hat{\mathbf{z}}$	(24*d*)	O II
$\mathbf{B}_{36}$	=	$z_4\,\mathbf{a}_1 + \left(\frac{1}{2}-x_4\right)\mathbf{a}_2 + \left(\frac{1}{2}+y_4\right)\mathbf{a}_3$	=	$z_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+y_4\right)a\,\hat{\mathbf{z}}$	(24*d*)	O II
$\mathbf{B}_{37}$	=	$-y_4\,\mathbf{a}_1 - z_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$-y_4 a\,\hat{\mathbf{x}} - z_4 a\,\hat{\mathbf{y}} - x_4 a\,\hat{\mathbf{z}}$	(24*d*)	O II
$\mathbf{B}_{38}$	=	$y_4\,\mathbf{a}_1 + \left(\frac{1}{2}-z_4\right)\mathbf{a}_2 + \left(\frac{1}{2}+x_4\right)\mathbf{a}_3$	=	$y_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-z_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{z}}$	(24*d*)	O II
$\mathbf{B}_{39}$	=	$\left(\frac{1}{2}-y_4\right)\mathbf{a}_1 + \left(\frac{1}{2}+z_4\right)\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+z_4\right)a\,\hat{\mathbf{y}} + x_4 a\,\hat{\mathbf{z}}$	(24*d*)	O II
$\mathbf{B}_{40}$	=	$\left(\frac{1}{2}+y_4\right)\mathbf{a}_1 + z_4\,\mathbf{a}_2 + \left(\frac{1}{2}-x_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_4\right)a\,\hat{\mathbf{x}} + z_4 a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{z}}$	(24*d*)	O II

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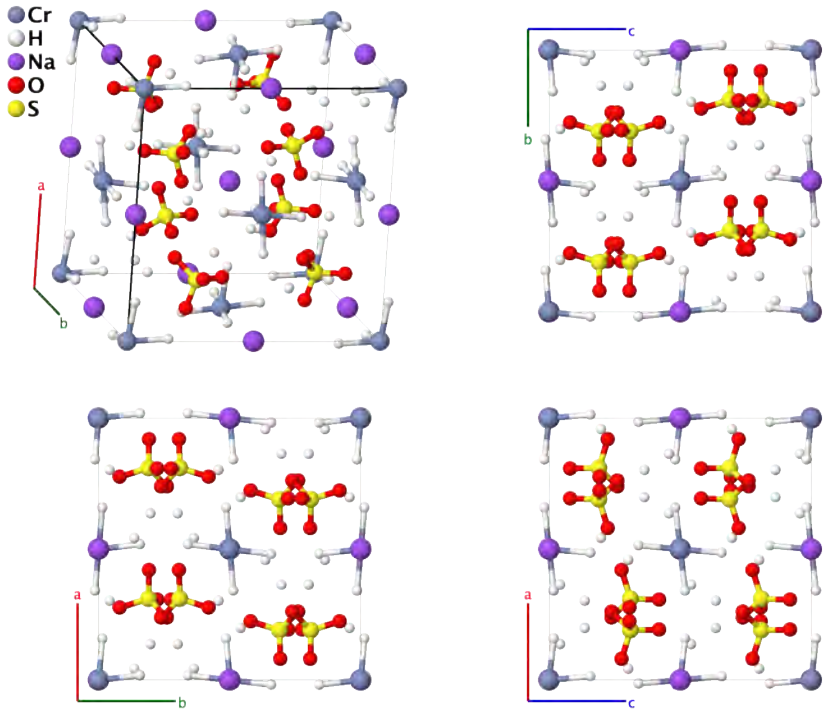
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1782
- POSCAR: pp. 1782

$NaCr(SO_4)_2{\cdot}12H_2O$ Alum Structure: AB12CD8E2_cP96_205_a_2d_b_cd_c

http://aflow.org/prototype-encyclopedia/AB12CD8E2_cP96_205_a_2d_b_cd_c



Prototype	:	$Cr(H_2O)_{12}NaO_8S_2$
AFLOW prototype label	:	AB12CD8E2_cP96_205_a_2d_b_cd_c
***Strukturbericht* designation**	:	None
Pearson symbol	:	cP96
Space group number	:	205
Space group symbol	:	$Pa\bar{3}$
AFLOW prototype command	:	`aflow --proto=AB12CD8E2_cP96_205_a_2d_b_cd_c --params=`$a, x_3, x_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7$

Other compounds with this structure

- $CsCr(SO_4)_2{\cdot}12H_2O$

- The alums have the general formula $AB(XO_4)_2{\cdot}12H_2O$, where A is a monovalent ion, B is a trivalent ion, and X is a chalcogen. In most cases atom B is aluminum and atom X is sulfur, leading to the name alum.
- All alums have their room-temperature form in space group $Pa\bar{3}$ #205, but the bonding between the A and B ions and the XO_4 complex can be quite different.
- (Lipson, 1935ab) described three general forms of alum based on the sizes of the monovalent ions. Each of these forms was given a *Strukturbericht* designation by (Gottfried, 1937):
  - α-alum, with intermediate sized ions, prototype $KAl(SO_4)_2{\cdot}12H_2O$, $H4_{13}$,

  - β-alum, with large ions, prototype $(NH_3CH_3)Al(SO_4)_2{\cdot}12H_2O$, $H4_{14}$, and
  - γ-alum, with small ions, prototype $NaAl(SO_4)_2{\cdot}12H_2O$, $H4_{15}$.

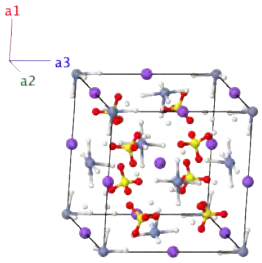
- This classification scheme is not complete, *e.g.*, (Ledsham, 1968) points out that $NaCr(SO_4)_2{\cdot}12H_2O$ (this structure) does not fit into any of these categories, and that the actual structure depends on the combination of monovalent and trivalent ions.

- As noted above, the $Pa\overline{3}$ structures of alum are the room temperature form. As the temperature decreases the alum structure may transform. For example, in the temperature range 150-170 K the β-alum $(NH_3CH_3)Al(SO_4)_2{\cdot}12H_2O$ transforms into an orthorhombic structure with fully ordered NH_3CH_3 ions.

- The positions of the hydrogen atoms in the water molecules were not determined, so we only provide the positions of the oxygen atoms (labeled as H_2O).

Simple Cubic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= a\,\hat{\mathbf{x}}\\ \mathbf{a}_2 &= a\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= a\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(4a)$	Cr
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	$(4a)$	Cr
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	$(4a)$	Cr
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$	$(4a)$	Cr
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	$(4b)$	Na
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{y}}$	$(4b)$	Na
$\mathbf{B}_7$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	$(4b)$	Na
$\mathbf{B}_8$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{z}}$	$(4b)$	Na
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	$(8c)$	O I
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{z}}$	$(8c)$	O I
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{z}}$	$(8c)$	O I
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	$(8c)$	O I
$\mathbf{B}_{13}$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	$(8c)$	O I
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{z}}$	$(8c)$	O I
$\mathbf{B}_{15}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{z}}$	$(8c)$	O I
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	$(8c)$	O I
$\mathbf{B}_{17}$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	$(8c)$	S
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \left(\frac{1}{2} + x_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{z}}$	$(8c)$	S

$\mathbf{B}_{19}$	$-x_4\,\mathbf{a}_1+\left(\frac{1}{2}+x_4\right)\mathbf{a}_2+\left(\frac{1}{2}-x_4\right)\mathbf{a}_3$	$-x_4a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{z}}$	(8*c*)	S
$\mathbf{B}_{20}$	$\left(\frac{1}{2}+x_4\right)\mathbf{a}_1+\left(\frac{1}{2}-x_4\right)\mathbf{a}_2-x_4\,\mathbf{a}_3$	$\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{y}}-x_4a\,\hat{\mathbf{z}}$	(8*c*)	S
$\mathbf{B}_{21}$	$-x_4\,\mathbf{a}_1-x_4\,\mathbf{a}_2-x_4\,\mathbf{a}_3$	$-x_4a\,\hat{\mathbf{x}}-x_4a\,\hat{\mathbf{y}}-x_4a\,\hat{\mathbf{z}}$	(8*c*)	S
$\mathbf{B}_{22}$	$\left(\frac{1}{2}+x_4\right)\mathbf{a}_1+x_4\,\mathbf{a}_2+\left(\frac{1}{2}-x_4\right)\mathbf{a}_3$	$\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{x}}+x_4a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{z}}$	(8*c*)	S
$\mathbf{B}_{23}$	$x_4\,\mathbf{a}_1+\left(\frac{1}{2}-x_4\right)\mathbf{a}_2+\left(\frac{1}{2}+x_4\right)\mathbf{a}_3$	$x_4a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{z}}$	(8*c*)	S
$\mathbf{B}_{24}$	$\left(\frac{1}{2}-x_4\right)\mathbf{a}_1+\left(\frac{1}{2}+x_4\right)\mathbf{a}_2+x_4\,\mathbf{a}_3$	$\left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{y}}+x_4a\,\hat{\mathbf{z}}$	(8*c*)	S
$\mathbf{B}_{25}$	$x_5\,\mathbf{a}_1+y_5\,\mathbf{a}_2+z_5\,\mathbf{a}_3$	$x_5a\,\hat{\mathbf{x}}+y_5a\,\hat{\mathbf{y}}+z_5a\,\hat{\mathbf{z}}$	(24*d*)	H_2O I
$\mathbf{B}_{26}$	$\left(\frac{1}{2}-x_5\right)\mathbf{a}_1-y_5\,\mathbf{a}_2+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	$\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{x}}-y_5a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)a\,\hat{\mathbf{z}}$	(24*d*)	H_2O I
$\mathbf{B}_{27}$	$-x_5\,\mathbf{a}_1+\left(\frac{1}{2}+y_5\right)\mathbf{a}_2+\left(\frac{1}{2}-z_5\right)\mathbf{a}_3$	$-x_5a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_5\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_5\right)a\,\hat{\mathbf{z}}$	(24*d*)	H_2O I
$\mathbf{B}_{28}$	$\left(\frac{1}{2}+x_5\right)\mathbf{a}_1+\left(\frac{1}{2}-y_5\right)\mathbf{a}_2-z_5\,\mathbf{a}_3$	$\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_5\right)a\,\hat{\mathbf{y}}-z_5a\,\hat{\mathbf{z}}$	(24*d*)	H_2O I
$\mathbf{B}_{29}$	$z_5\,\mathbf{a}_1+x_5\,\mathbf{a}_2+y_5\,\mathbf{a}_3$	$z_5a\,\hat{\mathbf{x}}+x_5a\,\hat{\mathbf{y}}+y_5a\,\hat{\mathbf{z}}$	(24*d*)	H_2O I
$\mathbf{B}_{30}$	$\left(\frac{1}{2}+z_5\right)\mathbf{a}_1+\left(\frac{1}{2}-x_5\right)\mathbf{a}_2-y_5\,\mathbf{a}_3$	$\left(\frac{1}{2}+z_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{y}}-y_5a\,\hat{\mathbf{z}}$	(24*d*)	H_2O I
$\mathbf{B}_{31}$	$\left(\frac{1}{2}-z_5\right)\mathbf{a}_1-x_5\,\mathbf{a}_2+\left(\frac{1}{2}+y_5\right)\mathbf{a}_3$	$\left(\frac{1}{2}-z_5\right)a\,\hat{\mathbf{x}}-x_5a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+y_5\right)a\,\hat{\mathbf{z}}$	(24*d*)	H_2O I
$\mathbf{B}_{32}$	$-z_5\,\mathbf{a}_1+\left(\frac{1}{2}+x_5\right)\mathbf{a}_2+\left(\frac{1}{2}-y_5\right)\mathbf{a}_3$	$-z_5a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-y_5\right)a\,\hat{\mathbf{z}}$	(24*d*)	H_2O I
$\mathbf{B}_{33}$	$y_5\,\mathbf{a}_1+z_5\,\mathbf{a}_2+x_5\,\mathbf{a}_3$	$y_5a\,\hat{\mathbf{x}}+z_5a\,\hat{\mathbf{y}}+x_5a\,\hat{\mathbf{z}}$	(24*d*)	H_2O I
$\mathbf{B}_{34}$	$-y_5\,\mathbf{a}_1+\left(\frac{1}{2}+z_5\right)\mathbf{a}_2+\left(\frac{1}{2}-x_5\right)\mathbf{a}_3$	$-y_5a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+z_5\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{z}}$	(24*d*)	H_2O I
$\mathbf{B}_{35}$	$\left(\frac{1}{2}+y_5\right)\mathbf{a}_1+\left(\frac{1}{2}-z_5\right)\mathbf{a}_2-x_5\,\mathbf{a}_3$	$\left(\frac{1}{2}+y_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-z_5\right)a\,\hat{\mathbf{y}}-x_5a\,\hat{\mathbf{z}}$	(24*d*)	H_2O I
$\mathbf{B}_{36}$	$\left(\frac{1}{2}-y_5\right)\mathbf{a}_1-z_5\,\mathbf{a}_2+\left(\frac{1}{2}+x_5\right)\mathbf{a}_3$	$\left(\frac{1}{2}-y_5\right)a\,\hat{\mathbf{x}}-z_5a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{z}}$	(24*d*)	H_2O I
$\mathbf{B}_{37}$	$-x_5\,\mathbf{a}_1-y_5\,\mathbf{a}_2-z_5\,\mathbf{a}_3$	$-x_5a\,\hat{\mathbf{x}}-y_5a\,\hat{\mathbf{y}}-z_5a\,\hat{\mathbf{z}}$	(24*d*)	H_2O I
$\mathbf{B}_{38}$	$\left(\frac{1}{2}+x_5\right)\mathbf{a}_1+y_5\,\mathbf{a}_2+\left(\frac{1}{2}-z_5\right)\mathbf{a}_3$	$\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{x}}+y_5a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_5\right)a\,\hat{\mathbf{z}}$	(24*d*)	H_2O I
$\mathbf{B}_{39}$	$x_5\,\mathbf{a}_1+\left(\frac{1}{2}-y_5\right)\mathbf{a}_2+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	$x_5a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_5\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)a\,\hat{\mathbf{z}}$	(24*d*)	H_2O I
$\mathbf{B}_{40}$	$\left(\frac{1}{2}-x_5\right)\mathbf{a}_1+\left(\frac{1}{2}+y_5\right)\mathbf{a}_2+z_5\,\mathbf{a}_3$	$\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_5\right)a\,\hat{\mathbf{y}}+z_5a\,\hat{\mathbf{z}}$	(24*d*)	H_2O I
$\mathbf{B}_{41}$	$-z_5\,\mathbf{a}_1-x_5\,\mathbf{a}_2-y_5\,\mathbf{a}_3$	$-z_5a\,\hat{\mathbf{x}}-x_5a\,\hat{\mathbf{y}}-y_5a\,\hat{\mathbf{z}}$	(24*d*)	H_2O I
$\mathbf{B}_{42}$	$\left(\frac{1}{2}-z_5\right)\mathbf{a}_1+\left(\frac{1}{2}+x_5\right)\mathbf{a}_2+y_5\,\mathbf{a}_3$	$\left(\frac{1}{2}-z_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{y}}+y_5a\,\hat{\mathbf{z}}$	(24*d*)	H_2O I
$\mathbf{B}_{43}$	$\left(\frac{1}{2}+z_5\right)\mathbf{a}_1+x_5\,\mathbf{a}_2+\left(\frac{1}{2}-y_5\right)\mathbf{a}_3$	$\left(\frac{1}{2}+z_5\right)a\,\hat{\mathbf{x}}+x_5a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-y_5\right)a\,\hat{\mathbf{z}}$	(24*d*)	H_2O I
$\mathbf{B}_{44}$	$z_5\,\mathbf{a}_1+\left(\frac{1}{2}-x_5\right)\mathbf{a}_2+\left(\frac{1}{2}+y_5\right)\mathbf{a}_3$	$z_5a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+y_5\right)a\,\hat{\mathbf{z}}$	(24*d*)	H_2O I
$\mathbf{B}_{45}$	$-y_5\,\mathbf{a}_1-z_5\,\mathbf{a}_2-x_5\,\mathbf{a}_3$	$-y_5a\,\hat{\mathbf{x}}-z_5a\,\hat{\mathbf{y}}-x_5a\,\hat{\mathbf{z}}$	(24*d*)	H_2O I
$\mathbf{B}_{46}$	$y_5\,\mathbf{a}_1+\left(\frac{1}{2}-z_5\right)\mathbf{a}_2+\left(\frac{1}{2}+x_5\right)\mathbf{a}_3$	$y_5a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-z_5\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{z}}$	(24*d*)	H_2O I
$\mathbf{B}_{47}$	$\left(\frac{1}{2}-y_5\right)\mathbf{a}_1+\left(\frac{1}{2}+z_5\right)\mathbf{a}_2+x_5\,\mathbf{a}_3$	$\left(\frac{1}{2}-y_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+z_5\right)a\,\hat{\mathbf{y}}+x_5a\,\hat{\mathbf{z}}$	(24*d*)	H_2O I
$\mathbf{B}_{48}$	$\left(\frac{1}{2}+y_5\right)\mathbf{a}_1+z_5\,\mathbf{a}_2+\left(\frac{1}{2}-x_5\right)\mathbf{a}_3$	$\left(\frac{1}{2}+y_5\right)a\,\hat{\mathbf{x}}+z_5a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{z}}$	(24*d*)	H_2O I
$\mathbf{B}_{49}$	$x_6\,\mathbf{a}_1+y_6\,\mathbf{a}_2+z_6\,\mathbf{a}_3$	$x_6a\,\hat{\mathbf{x}}+y_6a\,\hat{\mathbf{y}}+z_6a\,\hat{\mathbf{z}}$	(24*d*)	H_2O II
$\mathbf{B}_{50}$	$\left(\frac{1}{2}-x_6\right)\mathbf{a}_1-y_6\,\mathbf{a}_2+\left(\frac{1}{2}+z_6\right)\mathbf{a}_3$	$\left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{x}}-y_6a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_6\right)a\,\hat{\mathbf{z}}$	(24*d*)	H_2O II
$\mathbf{B}_{51}$	$-x_6\,\mathbf{a}_1+\left(\frac{1}{2}+y_6\right)\mathbf{a}_2+\left(\frac{1}{2}-z_6\right)\mathbf{a}_3$	$-x_6a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_6\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_6\right)a\,\hat{\mathbf{z}}$	(24*d*)	H_2O II
$\mathbf{B}_{52}$	$\left(\frac{1}{2}+x_6\right)\mathbf{a}_1+\left(\frac{1}{2}-y_6\right)\mathbf{a}_2-z_6\,\mathbf{a}_3$	$\left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_6\right)a\,\hat{\mathbf{y}}-z_6a\,\hat{\mathbf{z}}$	(24*d*)	H_2O II
$\mathbf{B}_{53}$	$z_6\,\mathbf{a}_1+x_6\,\mathbf{a}_2+y_6\,\mathbf{a}_3$	$z_6a\,\hat{\mathbf{x}}+x_6a\,\hat{\mathbf{y}}+y_6a\,\hat{\mathbf{z}}$	(24*d*)	H_2O II
$\mathbf{B}_{54}$	$\left(\frac{1}{2}+z_6\right)\mathbf{a}_1+\left(\frac{1}{2}-x_6\right)\mathbf{a}_2-y_6\,\mathbf{a}_3$	$\left(\frac{1}{2}+z_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{y}}-y_6a\,\hat{\mathbf{z}}$	(24*d*)	H_2O II

$$
\begin{aligned}
\mathbf{B}_{55} &= \left(\tfrac{1}{2}-z_6\right)\mathbf{a}_1 - x_6\,\mathbf{a}_2 + \left(\tfrac{1}{2}+y_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-z_6\right)a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+y_6\right)a\,\hat{\mathbf{z}} && (24d) && \mathrm{H_2O\ II}\\
\mathbf{B}_{56} &= -z_6\,\mathbf{a}_1 + \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-y_6\right)\mathbf{a}_3 &&= -z_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-y_6\right)a\,\hat{\mathbf{z}} && (24d) && \mathrm{H_2O\ II}\\
\mathbf{B}_{57} &= y_6\,\mathbf{a}_1 + z_6\,\mathbf{a}_2 + x_6\,\mathbf{a}_3 &&= y_6 a\,\hat{\mathbf{x}} + z_6 a\,\hat{\mathbf{y}} + x_6 a\,\hat{\mathbf{z}} && (24d) && \mathrm{H_2O\ II}\\
\mathbf{B}_{58} &= -y_6\,\mathbf{a}_1 + \left(\tfrac{1}{2}+z_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-x_6\right)\mathbf{a}_3 &&= -y_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+z_6\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{z}} && (24d) && \mathrm{H_2O\ II}\\
\mathbf{B}_{59} &= \left(\tfrac{1}{2}+y_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-z_6\right)\mathbf{a}_2 - x_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+y_6\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-z_6\right)a\,\hat{\mathbf{y}} - x_6 a\,\hat{\mathbf{z}} && (24d) && \mathrm{H_2O\ II}\\
\mathbf{B}_{60} &= \left(\tfrac{1}{2}-y_6\right)\mathbf{a}_1 - z_6\,\mathbf{a}_2 + \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-y_6\right)a\,\hat{\mathbf{x}} - z_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{z}} && (24d) && \mathrm{H_2O\ II}\\
\mathbf{B}_{61} &= -x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3 &&= -x_6 a\,\hat{\mathbf{x}} - y_6 a\,\hat{\mathbf{y}} - z_6 a\,\hat{\mathbf{z}} && (24d) && \mathrm{H_2O\ II}\\
\mathbf{B}_{62} &= \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_1 + y_6\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{x}} + y_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_6\right)a\,\hat{\mathbf{z}} && (24d) && \mathrm{H_2O\ II}\\
\mathbf{B}_{63} &= x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_6\right)\mathbf{a}_3 &&= x_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_6\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_6\right)a\,\hat{\mathbf{z}} && (24d) && \mathrm{H_2O\ II}\\
\mathbf{B}_{64} &= \left(\tfrac{1}{2}-x_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_6\right)\mathbf{a}_2 + z_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_6\right)a\,\hat{\mathbf{y}} + z_6 a\,\hat{\mathbf{z}} && (24d) && \mathrm{H_2O\ II}\\
\mathbf{B}_{65} &= -z_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 - y_6\,\mathbf{a}_3 &&= -z_6 a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} - y_6 a\,\hat{\mathbf{z}} && (24d) && \mathrm{H_2O\ II}\\
\mathbf{B}_{66} &= \left(\tfrac{1}{2}-z_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_2 + y_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-z_6\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{y}} + y_6 a\,\hat{\mathbf{z}} && (24d) && \mathrm{H_2O\ II}\\
\mathbf{B}_{67} &= \left(\tfrac{1}{2}+z_6\right)\mathbf{a}_1 + x_6\,\mathbf{a}_2 + \left(\tfrac{1}{2}-y_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+z_6\right)a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-y_6\right)a\,\hat{\mathbf{z}} && (24d) && \mathrm{H_2O\ II}\\
\mathbf{B}_{68} &= z_6\,\mathbf{a}_1 + \left(\tfrac{1}{2}-x_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+y_6\right)\mathbf{a}_3 &&= z_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+y_6\right)a\,\hat{\mathbf{z}} && (24d) && \mathrm{H_2O\ II}\\
\mathbf{B}_{69} &= -y_6\,\mathbf{a}_1 - z_6\,\mathbf{a}_2 - x_6\,\mathbf{a}_3 &&= -y_6 a\,\hat{\mathbf{x}} - z_6 a\,\hat{\mathbf{y}} - x_6 a\,\hat{\mathbf{z}} && (24d) && \mathrm{H_2O\ II}\\
\mathbf{B}_{70} &= y_6\,\mathbf{a}_1 + \left(\tfrac{1}{2}-z_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_3 &&= y_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-z_6\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{z}} && (24d) && \mathrm{H_2O\ II}\\
\mathbf{B}_{71} &= \left(\tfrac{1}{2}-y_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+z_6\right)\mathbf{a}_2 + x_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-y_6\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+z_6\right)a\,\hat{\mathbf{y}} + x_6 a\,\hat{\mathbf{z}} && (24d) && \mathrm{H_2O\ II}\\
\mathbf{B}_{72} &= \left(\tfrac{1}{2}+y_6\right)\mathbf{a}_1 + z_6\,\mathbf{a}_2 + \left(\tfrac{1}{2}-x_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+y_6\right)a\,\hat{\mathbf{x}} + z_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{z}} && (24d) && \mathrm{H_2O\ II}\\
\mathbf{B}_{73} &= x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 &&= x_7 a\,\hat{\mathbf{x}} + y_7 a\,\hat{\mathbf{y}} + z_7 a\,\hat{\mathbf{z}} && (24d) && \mathrm{O\ II}\\
\mathbf{B}_{74} &= \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{x}} - y_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{z}} && (24d) && \mathrm{O\ II}\\
\mathbf{B}_{75} &= -x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_3 &&= -x_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_7\right)a\,\hat{\mathbf{z}} && (24d) && \mathrm{O\ II}\\
\mathbf{B}_{76} &= \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_7\right)\mathbf{a}_2 - z_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_7\right)a\,\hat{\mathbf{y}} - z_7 a\,\hat{\mathbf{z}} && (24d) && \mathrm{O\ II}\\
\mathbf{B}_{77} &= z_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + y_7\,\mathbf{a}_3 &&= z_7 a\,\hat{\mathbf{x}} + x_7 a\,\hat{\mathbf{y}} + y_7 a\,\hat{\mathbf{z}} && (24d) && \mathrm{O\ II}\\
\mathbf{B}_{78} &= \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_2 - y_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{y}} - y_7 a\,\hat{\mathbf{z}} && (24d) && \mathrm{O\ II}\\
\mathbf{B}_{79} &= \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_1 - x_7\,\mathbf{a}_2 + \left(\tfrac{1}{2}+y_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-z_7\right)a\,\hat{\mathbf{x}} - x_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+y_7\right)a\,\hat{\mathbf{z}} && (24d) && \mathrm{O\ II}\\
\mathbf{B}_{80} &= -z_7\,\mathbf{a}_1 + \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-y_7\right)\mathbf{a}_3 &&= -z_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-y_7\right)a\,\hat{\mathbf{z}} && (24d) && \mathrm{O\ II}\\
\mathbf{B}_{81} &= y_7\,\mathbf{a}_1 + z_7\,\mathbf{a}_2 + x_7\,\mathbf{a}_3 &&= y_7 a\,\hat{\mathbf{x}} + z_7 a\,\hat{\mathbf{y}} + x_7 a\,\hat{\mathbf{z}} && (24d) && \mathrm{O\ II}\\
\mathbf{B}_{82} &= -y_7\,\mathbf{a}_1 + \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_3 &&= -y_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{z}} && (24d) && \mathrm{O\ II}\\
\mathbf{B}_{83} &= \left(\tfrac{1}{2}+y_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_2 - x_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+y_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-z_7\right)a\,\hat{\mathbf{y}} - x_7 a\,\hat{\mathbf{z}} && (24d) && \mathrm{O\ II}\\
\mathbf{B}_{84} &= \left(\tfrac{1}{2}-y_7\right)\mathbf{a}_1 - z_7\,\mathbf{a}_2 + \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-y_7\right)a\,\hat{\mathbf{x}} - z_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{z}} && (24d) && \mathrm{O\ II}\\
\mathbf{B}_{85} &= -x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3 &&= -x_7 a\,\hat{\mathbf{x}} - y_7 a\,\hat{\mathbf{y}} - z_7 a\,\hat{\mathbf{z}} && (24d) && \mathrm{O\ II}\\
\mathbf{B}_{86} &= \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1 + y_7\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{x}} + y_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_7\right)a\,\hat{\mathbf{z}} && (24d) && \mathrm{O\ II}\\
\mathbf{B}_{87} &= x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 &&= x_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{z}} && (24d) && \mathrm{O\ II}\\
\mathbf{B}_{88} &= \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_7\right)\mathbf{a}_2 + z_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_7\right)a\,\hat{\mathbf{y}} + z_7 a\,\hat{\mathbf{z}} && (24d) && \mathrm{O\ II}\\
\mathbf{B}_{89} &= -z_7\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 - y_7\,\mathbf{a}_3 &&= -z_7 a\,\hat{\mathbf{x}} - x_7 a\,\hat{\mathbf{y}} - y_7 a\,\hat{\mathbf{z}} && (24d) && \mathrm{O\ II}\\
\mathbf{B}_{90} &= \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_2 + y_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-z_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{y}} + y_7 a\,\hat{\mathbf{z}} && (24d) && \mathrm{O\ II}
\end{aligned}
$$

$$\mathbf{B}_{91} = \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_1 + x_7\,\mathbf{a}_2 + \left(\tfrac{1}{2}-y_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{x}} + x_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-y_7\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O II}$$

$$\mathbf{B}_{92} = z_7\,\mathbf{a}_1 + \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+y_7\right)\mathbf{a}_3 = z_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+y_7\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O II}$$

$$\mathbf{B}_{93} = -y_7\,\mathbf{a}_1 - z_7\,\mathbf{a}_2 - x_7\,\mathbf{a}_3 = -y_7 a\,\hat{\mathbf{x}} - z_7 a\,\hat{\mathbf{y}} - x_7 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O II}$$

$$\mathbf{B}_{94} = y_7\,\mathbf{a}_1 + \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_3 = y_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-z_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O II}$$

$$\mathbf{B}_{95} = \left(\tfrac{1}{2}-y_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_2 + x_7\,\mathbf{a}_3 = \left(\tfrac{1}{2}-y_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{y}} + x_7 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O II}$$

$$\mathbf{B}_{96} = \left(\tfrac{1}{2}+y_7\right)\mathbf{a}_1 + z_7\,\mathbf{a}_2 + \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+y_7\right)a\,\hat{\mathbf{x}} + z_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O II}$$

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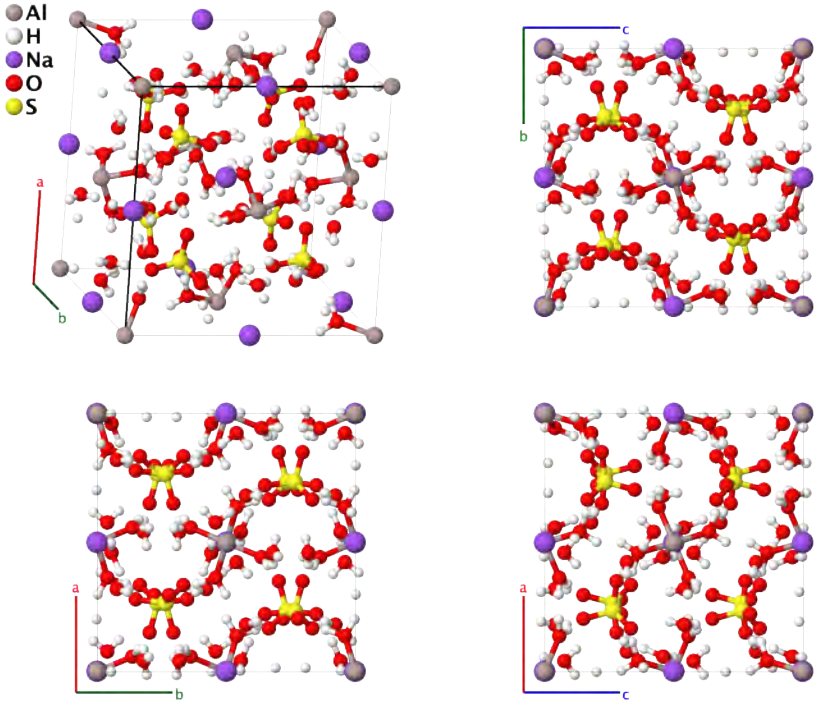
Geometry files:

- CIF: pp. 1782
- POSCAR: pp. 1783

γ-Alum [$AlNa(SO_4)_2{\cdot}12H_2O$, $H4_{15}$] Structure: AB24CD20E2_cP192_205_a_4d_b_c3d_c

http://aflow.org/prototype-encyclopedia/AB24CD20E2_cP192_205_a_4d_b_c3d_c



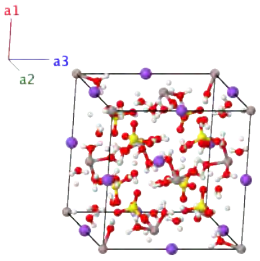
Prototype	:	$AlH_{24}NaO_{20}S_2$
AFLOW prototype label	:	AB24CD20E2_cP192_205_a_4d_b_c3d_c
***Strukturbericht* designation**	:	$H4_{15}$
Pearson symbol	:	cP192
Space group number	:	205
Space group symbol	:	$Pa\bar{3}$
AFLOW prototype command	:	`aflow --proto=AB24CD20E2_cP192_205_a_4d_b_c3d_c --params=`$a, x_3, x_4, x_5, y_5, z_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}$

- The alums have the general formula $AB(XO_4)_2{\cdot}12H_2O$, where A is a monovalent ion, B is a trivalent ion, and X is a chalcogen. In most cases atom B is aluminum and atom X is sulfur, leading to the name alum.
- All alums have their room-temperature form in space group $Pa\bar{3}$ #205, but the bonding between the A and B ions and the XO_4 complex can be quite different.
- (Lipson, 1935ab) described three general forms of alum based on the sizes of the monovalent ions. Each of these forms was given a *Strukturbericht* designation by (Gottfried, 1937):
  - α-alum, with intermediate sized ions, prototype $KAl(SO_4)_2{\cdot}12H_2O$, $H4_{13}$,
  - β-alum, with large ions, prototype $(NH_3CH_3)Al(SO_4)_2{\cdot}12H_2O$, $H4_{14}$, and
  - γ-alum, with small ions, prototype $NaAl(SO_4)_2{\cdot}12H_2O$, $H4_{15}$ (this structure).

- This classification scheme is not complete, *e.g.*, (Ledsham, 1968) points out that $NaCr(SO_4)_2{\cdot}12H_2O$ does not fit into any of these categories, and that the actual structure depends on the combination of monovalent and trivalent ions.
- As noted above, the $Pa\bar{3}$ structures of alum are the room temperature form. As the temperature decreases the alum structure may transform. For example, in the temperature range 150-170 K the β-alum $(NH_3CH_3)Al(SO_4)_2{\cdot}12H_2O$ transforms into an orthorhombic structure with fully ordered NH_3CH_3 ions.

Simple Cubic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= a\,\hat{\mathbf{x}}\\ \mathbf{a}_2 &= a\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= a\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(4*a*)	Al
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(4*a*)	Al
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(4*a*)	Al
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$	(4*a*)	Al
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(4*b*)	Na
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{y}}$	(4*b*)	Na
$\mathbf{B}_7$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	(4*b*)	Na
$\mathbf{B}_8$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{z}}$	(4*b*)	Na
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(8*c*)	O I
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2}-x_3\right)\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\frac{1}{2}+x_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{z}}$	(8*c*)	O I
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2}+x_3\right)\mathbf{a}_2 + \left(\frac{1}{2}-x_3\right)\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{z}}$	(8*c*)	O I
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2}+x_3\right)\mathbf{a}_1 + \left(\frac{1}{2}-x_3\right)\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(8*c*)	O I
$\mathbf{B}_{13}$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(8*c*)	O I
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2}+x_3\right)\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \left(\frac{1}{2}-x_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{z}}$	(8*c*)	O I
$\mathbf{B}_{15}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2}-x_3\right)\mathbf{a}_2 + \left(\frac{1}{2}+x_3\right)\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{z}}$	(8*c*)	O I
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2}-x_3\right)\mathbf{a}_1 + \left(\frac{1}{2}+x_3\right)\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(8*c*)	O I
$\mathbf{B}_{17}$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	(8*c*)	S
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2}-x_4\right)\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \left(\frac{1}{2}+x_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{z}}$	(8*c*)	S
$\mathbf{B}_{19}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2}+x_4\right)\mathbf{a}_2 + \left(\frac{1}{2}-x_4\right)\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{z}}$	(8*c*)	S
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2}+x_4\right)\mathbf{a}_1 + \left(\frac{1}{2}-x_4\right)\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{y}} - x_4a\,\hat{\mathbf{z}}$	(8*c*)	S
$\mathbf{B}_{21}$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} - x_4a\,\hat{\mathbf{z}}$	(8*c*)	S
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2}+x_4\right)\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \left(\frac{1}{2}-x_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{z}}$	(8*c*)	S
$\mathbf{B}_{23}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2}-x_4\right)\mathbf{a}_2 + \left(\frac{1}{2}+x_4\right)\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{z}}$	(8*c*)	S

$$
\begin{array}{lcccccc}
\mathbf{B}_{24} & = & \left(\frac{1}{2}-x_4\right)\mathbf{a}_1+\left(\frac{1}{2}+x_4\right)\mathbf{a}_2+x_4\,\mathbf{a}_3 & = & \left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{y}}+x_4a\,\hat{\mathbf{z}} & (8c) & \mathrm{S} \\
\mathbf{B}_{25} & = & x_5\,\mathbf{a}_1+y_5\,\mathbf{a}_2+z_5\,\mathbf{a}_3 & = & x_5a\,\hat{\mathbf{x}}+y_5a\,\hat{\mathbf{y}}+z_5a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\ I} \\
\mathbf{B}_{26} & = & \left(\frac{1}{2}-x_5\right)\mathbf{a}_1-y_5\,\mathbf{a}_2+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{x}}-y_5a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\ I} \\
\mathbf{B}_{27} & = & -x_5\,\mathbf{a}_1+\left(\frac{1}{2}+y_5\right)\mathbf{a}_2+\left(\frac{1}{2}-z_5\right)\mathbf{a}_3 & = & -x_5a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_5\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_5\right)a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\ I} \\
\mathbf{B}_{28} & = & \left(\frac{1}{2}+x_5\right)\mathbf{a}_1+\left(\frac{1}{2}-y_5\right)\mathbf{a}_2-z_5\,\mathbf{a}_3 & = & \left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_5\right)a\,\hat{\mathbf{y}}-z_5a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\ I} \\
\mathbf{B}_{29} & = & z_5\,\mathbf{a}_1+x_5\,\mathbf{a}_2+y_5\,\mathbf{a}_3 & = & z_5a\,\hat{\mathbf{x}}+x_5a\,\hat{\mathbf{y}}+y_5a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\ I} \\
\mathbf{B}_{30} & = & \left(\frac{1}{2}+z_5\right)\mathbf{a}_1+\left(\frac{1}{2}-x_5\right)\mathbf{a}_2-y_5\,\mathbf{a}_3 & = & \left(\frac{1}{2}+z_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{y}}-y_5a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\ I} \\
\mathbf{B}_{31} & = & \left(\frac{1}{2}-z_5\right)\mathbf{a}_1-x_5\,\mathbf{a}_2+\left(\frac{1}{2}+y_5\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-z_5\right)a\,\hat{\mathbf{x}}-x_5a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+y_5\right)a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\ I} \\
\mathbf{B}_{32} & = & -z_5\,\mathbf{a}_1+\left(\frac{1}{2}+x_5\right)\mathbf{a}_2+\left(\frac{1}{2}-y_5\right)\mathbf{a}_3 & = & -z_5a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-y_5\right)a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\ I} \\
\mathbf{B}_{33} & = & y_5\,\mathbf{a}_1+z_5\,\mathbf{a}_2+x_5\,\mathbf{a}_3 & = & y_5a\,\hat{\mathbf{x}}+z_5a\,\hat{\mathbf{y}}+x_5a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\ I} \\
\mathbf{B}_{34} & = & -y_5\,\mathbf{a}_1+\left(\frac{1}{2}+z_5\right)\mathbf{a}_2+\left(\frac{1}{2}-x_5\right)\mathbf{a}_3 & = & -y_5a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+z_5\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\ I} \\
\mathbf{B}_{35} & = & \left(\frac{1}{2}+y_5\right)\mathbf{a}_1+\left(\frac{1}{2}-z_5\right)\mathbf{a}_2-x_5\,\mathbf{a}_3 & = & \left(\frac{1}{2}+y_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-z_5\right)a\,\hat{\mathbf{y}}-x_5a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\ I} \\
\mathbf{B}_{36} & = & \left(\frac{1}{2}-y_5\right)\mathbf{a}_1-z_5\,\mathbf{a}_2+\left(\frac{1}{2}+x_5\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-y_5\right)a\,\hat{\mathbf{x}}-z_5a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\ I} \\
\mathbf{B}_{37} & = & -x_5\,\mathbf{a}_1-y_5\,\mathbf{a}_2-z_5\,\mathbf{a}_3 & = & -x_5a\,\hat{\mathbf{x}}-y_5a\,\hat{\mathbf{y}}-z_5a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\ I} \\
\mathbf{B}_{38} & = & \left(\frac{1}{2}+x_5\right)\mathbf{a}_1+y_5\,\mathbf{a}_2+\left(\frac{1}{2}-z_5\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{x}}+y_5a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_5\right)a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\ I} \\
\mathbf{B}_{39} & = & x_5\,\mathbf{a}_1+\left(\frac{1}{2}-y_5\right)\mathbf{a}_2+\left(\frac{1}{2}+z_5\right)\mathbf{a}_3 & = & x_5a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_5\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_5\right)a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\ I} \\
\mathbf{B}_{40} & = & \left(\frac{1}{2}-x_5\right)\mathbf{a}_1+\left(\frac{1}{2}+y_5\right)\mathbf{a}_2+z_5\,\mathbf{a}_3 & = & \left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_5\right)a\,\hat{\mathbf{y}}+z_5a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\ I} \\
\mathbf{B}_{41} & = & -z_5\,\mathbf{a}_1-x_5\,\mathbf{a}_2-y_5\,\mathbf{a}_3 & = & -z_5a\,\hat{\mathbf{x}}-x_5a\,\hat{\mathbf{y}}-y_5a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\ I} \\
\mathbf{B}_{42} & = & \left(\frac{1}{2}-z_5\right)\mathbf{a}_1+\left(\frac{1}{2}+x_5\right)\mathbf{a}_2+y_5\,\mathbf{a}_3 & = & \left(\frac{1}{2}-z_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{y}}+y_5a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\ I} \\
\mathbf{B}_{43} & = & \left(\frac{1}{2}+z_5\right)\mathbf{a}_1+x_5\,\mathbf{a}_2+\left(\frac{1}{2}-y_5\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+z_5\right)a\,\hat{\mathbf{x}}+x_5a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-y_5\right)a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\ I} \\
\mathbf{B}_{44} & = & z_5\,\mathbf{a}_1+\left(\frac{1}{2}-x_5\right)\mathbf{a}_2+\left(\frac{1}{2}+y_5\right)\mathbf{a}_3 & = & z_5a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+y_5\right)a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\ I} \\
\mathbf{B}_{45} & = & -y_5\,\mathbf{a}_1-z_5\,\mathbf{a}_2-x_5\,\mathbf{a}_3 & = & -y_5a\,\hat{\mathbf{x}}-z_5a\,\hat{\mathbf{y}}-x_5a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\ I} \\
\mathbf{B}_{46} & = & y_5\,\mathbf{a}_1+\left(\frac{1}{2}-z_5\right)\mathbf{a}_2+\left(\frac{1}{2}+x_5\right)\mathbf{a}_3 & = & y_5a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-z_5\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\ I} \\
\mathbf{B}_{47} & = & \left(\frac{1}{2}-y_5\right)\mathbf{a}_1+\left(\frac{1}{2}+z_5\right)\mathbf{a}_2+x_5\,\mathbf{a}_3 & = & \left(\frac{1}{2}-y_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+z_5\right)a\,\hat{\mathbf{y}}+x_5a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\ I} \\
\mathbf{B}_{48} & = & \left(\frac{1}{2}+y_5\right)\mathbf{a}_1+z_5\,\mathbf{a}_2+\left(\frac{1}{2}-x_5\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+y_5\right)a\,\hat{\mathbf{x}}+z_5a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\ I} \\
\mathbf{B}_{49} & = & x_6\,\mathbf{a}_1+y_6\,\mathbf{a}_2+z_6\,\mathbf{a}_3 & = & x_6a\,\hat{\mathbf{x}}+y_6a\,\hat{\mathbf{y}}+z_6a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\ II} \\
\mathbf{B}_{50} & = & \left(\frac{1}{2}-x_6\right)\mathbf{a}_1-y_6\,\mathbf{a}_2+\left(\frac{1}{2}+z_6\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{x}}-y_6a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_6\right)a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\ II} \\
\mathbf{B}_{51} & = & -x_6\,\mathbf{a}_1+\left(\frac{1}{2}+y_6\right)\mathbf{a}_2+\left(\frac{1}{2}-z_6\right)\mathbf{a}_3 & = & -x_6a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_6\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_6\right)a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\ II} \\
\mathbf{B}_{52} & = & \left(\frac{1}{2}+x_6\right)\mathbf{a}_1+\left(\frac{1}{2}-y_6\right)\mathbf{a}_2-z_6\,\mathbf{a}_3 & = & \left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_6\right)a\,\hat{\mathbf{y}}-z_6a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\ II} \\
\mathbf{B}_{53} & = & z_6\,\mathbf{a}_1+x_6\,\mathbf{a}_2+y_6\,\mathbf{a}_3 & = & z_6a\,\hat{\mathbf{x}}+x_6a\,\hat{\mathbf{y}}+y_6a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\ II} \\
\mathbf{B}_{54} & = & \left(\frac{1}{2}+z_6\right)\mathbf{a}_1+\left(\frac{1}{2}-x_6\right)\mathbf{a}_2-y_6\,\mathbf{a}_3 & = & \left(\frac{1}{2}+z_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{y}}-y_6a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\ II} \\
\mathbf{B}_{55} & = & \left(\frac{1}{2}-z_6\right)\mathbf{a}_1-x_6\,\mathbf{a}_2+\left(\frac{1}{2}+y_6\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-z_6\right)a\,\hat{\mathbf{x}}-x_6a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+y_6\right)a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\ II} \\
\mathbf{B}_{56} & = & -z_6\,\mathbf{a}_1+\left(\frac{1}{2}+x_6\right)\mathbf{a}_2+\left(\frac{1}{2}-y_6\right)\mathbf{a}_3 & = & -z_6a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-y_6\right)a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\ II} \\
\mathbf{B}_{57} & = & y_6\,\mathbf{a}_1+z_6\,\mathbf{a}_2+x_6\,\mathbf{a}_3 & = & y_6a\,\hat{\mathbf{x}}+z_6a\,\hat{\mathbf{y}}+x_6a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\ II} \\
\mathbf{B}_{58} & = & -y_6\,\mathbf{a}_1+\left(\frac{1}{2}+z_6\right)\mathbf{a}_2+\left(\frac{1}{2}-x_6\right)\mathbf{a}_3 & = & -y_6a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+z_6\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\ II} \\
\mathbf{B}_{59} & = & \left(\frac{1}{2}+y_6\right)\mathbf{a}_1+\left(\frac{1}{2}-z_6\right)\mathbf{a}_2-x_6\,\mathbf{a}_3 & = & \left(\frac{1}{2}+y_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-z_6\right)a\,\hat{\mathbf{y}}-x_6a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\ II}
\end{array}
$$

$$
\begin{aligned}
\mathbf{B}_{60} &= \left(\tfrac{1}{2}-y_6\right)\mathbf{a}_1 - z_6\,\mathbf{a}_2 + \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-y_6\right)a\,\hat{\mathbf{x}} - z_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{61} &= -x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3 &&= -x_6 a\,\hat{\mathbf{x}} - y_6 a\,\hat{\mathbf{y}} - z_6 a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{62} &= \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_1 + y_6\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{x}} + y_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_6\right)a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{63} &= x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_6\right)\mathbf{a}_3 &&= x_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_6\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_6\right)a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{64} &= \left(\tfrac{1}{2}-x_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_6\right)\mathbf{a}_2 + z_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_6\right)a\,\hat{\mathbf{y}} + z_6 a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{65} &= -z_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 - y_6\,\mathbf{a}_3 &&= -z_6 a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} - y_6 a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{66} &= \left(\tfrac{1}{2}-z_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_2 + y_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-z_6\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{y}} + y_6 a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{67} &= \left(\tfrac{1}{2}+z_6\right)\mathbf{a}_1 + x_6\,\mathbf{a}_2 + \left(\tfrac{1}{2}-y_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+z_6\right)a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-y_6\right)a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{68} &= z_6\,\mathbf{a}_1 + \left(\tfrac{1}{2}-x_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+y_6\right)\mathbf{a}_3 &&= z_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+y_6\right)a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{69} &= -y_6\,\mathbf{a}_1 - z_6\,\mathbf{a}_2 - x_6\,\mathbf{a}_3 &&= -y_6 a\,\hat{\mathbf{x}} - z_6 a\,\hat{\mathbf{y}} - x_6 a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{70} &= y_6\,\mathbf{a}_1 + \left(\tfrac{1}{2}-z_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_3 &&= y_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-z_6\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{71} &= \left(\tfrac{1}{2}-y_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+z_6\right)\mathbf{a}_2 + x_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-y_6\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+z_6\right)a\,\hat{\mathbf{y}} + x_6 a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{72} &= \left(\tfrac{1}{2}+y_6\right)\mathbf{a}_1 + z_6\,\mathbf{a}_2 + \left(\tfrac{1}{2}-x_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+y_6\right)a\,\hat{\mathbf{x}} + z_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{73} &= x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 &&= x_7 a\,\hat{\mathbf{x}} + y_7 a\,\hat{\mathbf{y}} + z_7 a\,\hat{\mathbf{z}} && (24d) && \text{H III}\\
\mathbf{B}_{74} &= \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{x}} - y_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{z}} && (24d) && \text{H III}\\
\mathbf{B}_{75} &= -x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_3 &&= -x_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_7\right)a\,\hat{\mathbf{z}} && (24d) && \text{H III}\\
\mathbf{B}_{76} &= \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_7\right)\mathbf{a}_2 - z_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_7\right)a\,\hat{\mathbf{y}} - z_7 a\,\hat{\mathbf{z}} && (24d) && \text{H III}\\
\mathbf{B}_{77} &= z_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + y_7\,\mathbf{a}_3 &&= z_7 a\,\hat{\mathbf{x}} + x_7 a\,\hat{\mathbf{y}} + y_7 a\,\hat{\mathbf{z}} && (24d) && \text{H III}\\
\mathbf{B}_{78} &= \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_2 - y_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{y}} - y_7 a\,\hat{\mathbf{z}} && (24d) && \text{H III}\\
\mathbf{B}_{79} &= \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_1 - x_7\,\mathbf{a}_2 + \left(\tfrac{1}{2}+y_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-z_7\right)a\,\hat{\mathbf{x}} - x_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+y_7\right)a\,\hat{\mathbf{z}} && (24d) && \text{H III}\\
\mathbf{B}_{80} &= -z_7\,\mathbf{a}_1 + \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-y_7\right)\mathbf{a}_3 &&= -z_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-y_7\right)a\,\hat{\mathbf{z}} && (24d) && \text{H III}\\
\mathbf{B}_{81} &= y_7\,\mathbf{a}_1 + z_7\,\mathbf{a}_2 + x_7\,\mathbf{a}_3 &&= y_7 a\,\hat{\mathbf{x}} + z_7 a\,\hat{\mathbf{y}} + x_7 a\,\hat{\mathbf{z}} && (24d) && \text{H III}\\
\mathbf{B}_{82} &= -y_7\,\mathbf{a}_1 + \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_3 &&= -y_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{z}} && (24d) && \text{H III}\\
\mathbf{B}_{83} &= \left(\tfrac{1}{2}+y_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_2 - x_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+y_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-z_7\right)a\,\hat{\mathbf{y}} - x_7 a\,\hat{\mathbf{z}} && (24d) && \text{H III}\\
\mathbf{B}_{84} &= \left(\tfrac{1}{2}-y_7\right)\mathbf{a}_1 - z_7\,\mathbf{a}_2 + \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-y_7\right)a\,\hat{\mathbf{x}} - z_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{z}} && (24d) && \text{H III}\\
\mathbf{B}_{85} &= -x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3 &&= -x_7 a\,\hat{\mathbf{x}} - y_7 a\,\hat{\mathbf{y}} - z_7 a\,\hat{\mathbf{z}} && (24d) && \text{H III}\\
\mathbf{B}_{86} &= \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1 + y_7\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{x}} + y_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_7\right)a\,\hat{\mathbf{z}} && (24d) && \text{H III}\\
\mathbf{B}_{87} &= x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 &&= x_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{z}} && (24d) && \text{H III}\\
\mathbf{B}_{88} &= \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_7\right)\mathbf{a}_2 + z_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_7\right)a\,\hat{\mathbf{y}} + z_7 a\,\hat{\mathbf{z}} && (24d) && \text{H III}\\
\mathbf{B}_{89} &= -z_7\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 - y_7\,\mathbf{a}_3 &&= -z_7 a\,\hat{\mathbf{x}} - x_7 a\,\hat{\mathbf{y}} - y_7 a\,\hat{\mathbf{z}} && (24d) && \text{H III}\\
\mathbf{B}_{90} &= \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_2 + y_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-z_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{y}} + y_7 a\,\hat{\mathbf{z}} && (24d) && \text{H III}\\
\mathbf{B}_{91} &= \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_1 + x_7\,\mathbf{a}_2 + \left(\tfrac{1}{2}-y_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{x}} + x_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-y_7\right)a\,\hat{\mathbf{z}} && (24d) && \text{H III}\\
\mathbf{B}_{92} &= z_7\,\mathbf{a}_1 + \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+y_7\right)\mathbf{a}_3 &&= z_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+y_7\right)a\,\hat{\mathbf{z}} && (24d) && \text{H III}\\
\mathbf{B}_{93} &= -y_7\,\mathbf{a}_1 - z_7\,\mathbf{a}_2 - x_7\,\mathbf{a}_3 &&= -y_7 a\,\hat{\mathbf{x}} - z_7 a\,\hat{\mathbf{y}} - x_7 a\,\hat{\mathbf{z}} && (24d) && \text{H III}\\
\mathbf{B}_{94} &= y_7\,\mathbf{a}_1 + \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_3 &&= y_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-z_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{z}} && (24d) && \text{H III}\\
\mathbf{B}_{95} &= \left(\tfrac{1}{2}-y_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_2 + x_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-y_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{y}} + x_7 a\,\hat{\mathbf{z}} && (24d) && \text{H III}
\end{aligned}
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\begin{aligned}
\mathbf{B}_{96} &= \left(\tfrac{1}{2}+y_7\right)\mathbf{a}_1 + z_7\,\mathbf{a}_2 + \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+y_7\right)a\,\hat{\mathbf{x}} + z_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{z}} && (24d) && \text{H III}\\
\mathbf{B}_{97} &= x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 &&= x_8 a\,\hat{\mathbf{x}} + y_8 a\,\hat{\mathbf{y}} + z_8 a\,\hat{\mathbf{z}} && (24d) && \text{H IV}\\
\mathbf{B}_{98} &= \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_1 - y_8\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_8\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{x}} - y_8 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_8\right)a\,\hat{\mathbf{z}} && (24d) && \text{H IV}\\
\mathbf{B}_{99} &= -x_8\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_8\right)\mathbf{a}_3 &&= -x_8 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_8\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_8\right)a\,\hat{\mathbf{z}} && (24d) && \text{H IV}\\
\mathbf{B}_{100} &= \left(\tfrac{1}{2}+x_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_8\right)\mathbf{a}_2 - z_8\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_8\right)a\,\hat{\mathbf{y}} - z_8 a\,\hat{\mathbf{z}} && (24d) && \text{H IV}\\
\mathbf{B}_{101} &= z_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 + y_8\,\mathbf{a}_3 &&= z_8 a\,\hat{\mathbf{x}} + x_8 a\,\hat{\mathbf{y}} + y_8 a\,\hat{\mathbf{z}} && (24d) && \text{H IV}\\
\mathbf{B}_{102} &= \left(\tfrac{1}{2}+z_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_2 - y_8\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+z_8\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{y}} - y_8 a\,\hat{\mathbf{z}} && (24d) && \text{H IV}\\
\mathbf{B}_{103} &= \left(\tfrac{1}{2}-z_8\right)\mathbf{a}_1 - x_8\,\mathbf{a}_2 + \left(\tfrac{1}{2}+y_8\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-z_8\right)a\,\hat{\mathbf{x}} - x_8 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+y_8\right)a\,\hat{\mathbf{z}} && (24d) && \text{H IV}\\
\mathbf{B}_{104} &= -z_8\,\mathbf{a}_1 + \left(\tfrac{1}{2}+x_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-y_8\right)\mathbf{a}_3 &&= -z_8 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-y_8\right)a\,\hat{\mathbf{z}} && (24d) && \text{H IV}\\
\mathbf{B}_{105} &= y_8\,\mathbf{a}_1 + z_8\,\mathbf{a}_2 + x_8\,\mathbf{a}_3 &&= y_8 a\,\hat{\mathbf{x}} + z_8 a\,\hat{\mathbf{y}} + x_8 a\,\hat{\mathbf{z}} && (24d) && \text{H IV}\\
\mathbf{B}_{106} &= -y_8\,\mathbf{a}_1 + \left(\tfrac{1}{2}+z_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_3 &&= -y_8 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+z_8\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{z}} && (24d) && \text{H IV}\\
\mathbf{B}_{107} &= \left(\tfrac{1}{2}+y_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-z_8\right)\mathbf{a}_2 - x_8\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+y_8\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-z_8\right)a\,\hat{\mathbf{y}} - x_8 a\,\hat{\mathbf{z}} && (24d) && \text{H IV}\\
\mathbf{B}_{108} &= \left(\tfrac{1}{2}-y_8\right)\mathbf{a}_1 - z_8\,\mathbf{a}_2 + \left(\tfrac{1}{2}+x_8\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-y_8\right)a\,\hat{\mathbf{x}} - z_8 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{z}} && (24d) && \text{H IV}\\
\mathbf{B}_{109} &= -x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3 &&= -x_8 a\,\hat{\mathbf{x}} - y_8 a\,\hat{\mathbf{y}} - z_8 a\,\hat{\mathbf{z}} && (24d) && \text{H IV}\\
\mathbf{B}_{110} &= \left(\tfrac{1}{2}+x_8\right)\mathbf{a}_1 + y_8\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_8\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{x}} + y_8 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_8\right)a\,\hat{\mathbf{z}} && (24d) && \text{H IV}\\
\mathbf{B}_{111} &= x_8\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_8\right)\mathbf{a}_3 &&= x_8 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_8\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_8\right)a\,\hat{\mathbf{z}} && (24d) && \text{H IV}\\
\mathbf{B}_{112} &= \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_8\right)\mathbf{a}_2 + z_8\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_8\right)a\,\hat{\mathbf{y}} + z_8 a\,\hat{\mathbf{z}} && (24d) && \text{H IV}\\
\mathbf{B}_{113} &= -z_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 - y_8\,\mathbf{a}_3 &&= -z_8 a\,\hat{\mathbf{x}} - x_8 a\,\hat{\mathbf{y}} - y_8 a\,\hat{\mathbf{z}} && (24d) && \text{H IV}\\
\mathbf{B}_{114} &= \left(\tfrac{1}{2}-z_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+x_8\right)\mathbf{a}_2 + y_8\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-z_8\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{y}} + y_8 a\,\hat{\mathbf{z}} && (24d) && \text{H IV}\\
\mathbf{B}_{115} &= \left(\tfrac{1}{2}+z_8\right)\mathbf{a}_1 + x_8\,\mathbf{a}_2 + \left(\tfrac{1}{2}-y_8\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+z_8\right)a\,\hat{\mathbf{x}} + x_8 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-y_8\right)a\,\hat{\mathbf{z}} && (24d) && \text{H IV}\\
\mathbf{B}_{116} &= z_8\,\mathbf{a}_1 + \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+y_8\right)\mathbf{a}_3 &&= z_8 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+y_8\right)a\,\hat{\mathbf{z}} && (24d) && \text{H IV}\\
\mathbf{B}_{117} &= -y_8\,\mathbf{a}_1 - z_8\,\mathbf{a}_2 - x_8\,\mathbf{a}_3 &&= -y_8 a\,\hat{\mathbf{x}} - z_8 a\,\hat{\mathbf{y}} - x_8 a\,\hat{\mathbf{z}} && (24d) && \text{H IV}\\
\mathbf{B}_{118} &= y_8\,\mathbf{a}_1 + \left(\tfrac{1}{2}-z_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+x_8\right)\mathbf{a}_3 &&= y_8 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-z_8\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{z}} && (24d) && \text{H IV}\\
\mathbf{B}_{119} &= \left(\tfrac{1}{2}-y_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+z_8\right)\mathbf{a}_2 + x_8\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-y_8\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+z_8\right)a\,\hat{\mathbf{y}} + x_8 a\,\hat{\mathbf{z}} && (24d) && \text{H IV}\\
\mathbf{B}_{120} &= \left(\tfrac{1}{2}+y_8\right)\mathbf{a}_1 + z_8\,\mathbf{a}_2 + \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+y_8\right)a\,\hat{\mathbf{x}} + z_8 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{z}} && (24d) && \text{H IV}\\
\mathbf{B}_{121} &= x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 &&= x_9 a\,\hat{\mathbf{x}} + y_9 a\,\hat{\mathbf{y}} + z_9 a\,\hat{\mathbf{z}} && (24d) && \text{O II}\\
\mathbf{B}_{122} &= \left(\tfrac{1}{2}-x_9\right)\mathbf{a}_1 - y_9\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_9\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_9\right)a\,\hat{\mathbf{x}} - y_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_9\right)a\,\hat{\mathbf{z}} && (24d) && \text{O II}\\
\mathbf{B}_{123} &= -x_9\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_9\right)\mathbf{a}_3 &&= -x_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_9\right)a\,\hat{\mathbf{z}} && (24d) && \text{O II}\\
\mathbf{B}_{124} &= \left(\tfrac{1}{2}+x_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_9\right)\mathbf{a}_2 - z_9\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_9\right)a\,\hat{\mathbf{y}} - z_9 a\,\hat{\mathbf{z}} && (24d) && \text{O II}\\
\mathbf{B}_{125} &= z_9\,\mathbf{a}_1 + x_9\,\mathbf{a}_2 + y_9\,\mathbf{a}_3 &&= z_9 a\,\hat{\mathbf{x}} + x_9 a\,\hat{\mathbf{y}} + y_9 a\,\hat{\mathbf{z}} && (24d) && \text{O II}\\
\mathbf{B}_{126} &= \left(\tfrac{1}{2}+z_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-x_9\right)\mathbf{a}_2 - y_9\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+z_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_9\right)a\,\hat{\mathbf{y}} - y_9 a\,\hat{\mathbf{z}} && (24d) && \text{O II}\\
\mathbf{B}_{127} &= \left(\tfrac{1}{2}-z_9\right)\mathbf{a}_1 - x_9\,\mathbf{a}_2 + \left(\tfrac{1}{2}+y_9\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-z_9\right)a\,\hat{\mathbf{x}} - x_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+y_9\right)a\,\hat{\mathbf{z}} && (24d) && \text{O II}\\
\mathbf{B}_{128} &= -z_9\,\mathbf{a}_1 + \left(\tfrac{1}{2}+x_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-y_9\right)\mathbf{a}_3 &&= -z_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-y_9\right)a\,\hat{\mathbf{z}} && (24d) && \text{O II}\\
\mathbf{B}_{129} &= y_9\,\mathbf{a}_1 + z_9\,\mathbf{a}_2 + x_9\,\mathbf{a}_3 &&= y_9 a\,\hat{\mathbf{x}} + z_9 a\,\hat{\mathbf{y}} + x_9 a\,\hat{\mathbf{z}} && (24d) && \text{O II}\\
\mathbf{B}_{130} &= -y_9\,\mathbf{a}_1 + \left(\tfrac{1}{2}+z_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-x_9\right)\mathbf{a}_3 &&= -y_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+z_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-x_9\right)a\,\hat{\mathbf{z}} && (24d) && \text{O II}\\
\mathbf{B}_{131} &= \left(\tfrac{1}{2}+y_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-z_9\right)\mathbf{a}_2 - x_9\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+y_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-z_9\right)a\,\hat{\mathbf{y}} - x_9 a\,\hat{\mathbf{z}} && (24d) && \text{O II}
\end{aligned}
$$

$$
\begin{aligned}
\mathbf{B}_{132} &= \left(\tfrac{1}{2}-y_9\right)\mathbf{a}_1 - z_9\,\mathbf{a}_2 + \left(\tfrac{1}{2}+x_9\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-y_9\right)a\,\hat{\mathbf{x}} - z_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_9\right)a\,\hat{\mathbf{z}} && (24d) && \text{O II}\\
\mathbf{B}_{133} &= -x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3 &&= -x_9 a\,\hat{\mathbf{x}} - y_9 a\,\hat{\mathbf{y}} - z_9 a\,\hat{\mathbf{z}} && (24d) && \text{O II}\\
\mathbf{B}_{134} &= \left(\tfrac{1}{2}+x_9\right)\mathbf{a}_1 + y_9\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_9\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_9\right)a\,\hat{\mathbf{x}} + y_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_9\right)a\,\hat{\mathbf{z}} && (24d) && \text{O II}\\
\mathbf{B}_{135} &= x_9\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_9\right)\mathbf{a}_3 &&= x_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_9\right)a\,\hat{\mathbf{z}} && (24d) && \text{O II}\\
\mathbf{B}_{136} &= \left(\tfrac{1}{2}-x_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_9\right)\mathbf{a}_2 + z_9\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_9\right)a\,\hat{\mathbf{y}} + z_9 a\,\hat{\mathbf{z}} && (24d) && \text{O II}\\
\mathbf{B}_{137} &= -z_9\,\mathbf{a}_1 - x_9\,\mathbf{a}_2 - y_9\,\mathbf{a}_3 &&= -z_9 a\,\hat{\mathbf{x}} - x_9 a\,\hat{\mathbf{y}} - y_9 a\,\hat{\mathbf{z}} && (24d) && \text{O II}\\
\mathbf{B}_{138} &= \left(\tfrac{1}{2}-z_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+x_9\right)\mathbf{a}_2 + y_9\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-z_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_9\right)a\,\hat{\mathbf{y}} + y_9 a\,\hat{\mathbf{z}} && (24d) && \text{O II}\\
\mathbf{B}_{139} &= \left(\tfrac{1}{2}+z_9\right)\mathbf{a}_1 + x_9\,\mathbf{a}_2 + \left(\tfrac{1}{2}-y_9\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+z_9\right)a\,\hat{\mathbf{x}} + x_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-y_9\right)a\,\hat{\mathbf{z}} && (24d) && \text{O II}\\
\mathbf{B}_{140} &= z_9\,\mathbf{a}_1 + \left(\tfrac{1}{2}-x_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+y_9\right)\mathbf{a}_3 &&= z_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+y_9\right)a\,\hat{\mathbf{z}} && (24d) && \text{O II}\\
\mathbf{B}_{141} &= -y_9\,\mathbf{a}_1 - z_9\,\mathbf{a}_2 - x_9\,\mathbf{a}_3 &&= -y_9 a\,\hat{\mathbf{x}} - z_9 a\,\hat{\mathbf{y}} - x_9 a\,\hat{\mathbf{z}} && (24d) && \text{O II}\\
\mathbf{B}_{142} &= y_9\,\mathbf{a}_1 + \left(\tfrac{1}{2}-z_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+x_9\right)\mathbf{a}_3 &&= y_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-z_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_9\right)a\,\hat{\mathbf{z}} && (24d) && \text{O II}\\
\mathbf{B}_{143} &= \left(\tfrac{1}{2}-y_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+z_9\right)\mathbf{a}_2 + x_9\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-y_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+z_9\right)a\,\hat{\mathbf{y}} + x_9 a\,\hat{\mathbf{z}} && (24d) && \text{O II}\\
\mathbf{B}_{144} &= \left(\tfrac{1}{2}+y_9\right)\mathbf{a}_1 + z_9\,\mathbf{a}_2 + \left(\tfrac{1}{2}-x_9\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+y_9\right)a\,\hat{\mathbf{x}} + z_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-x_9\right)a\,\hat{\mathbf{z}} && (24d) && \text{O II}\\
\mathbf{B}_{145} &= x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 &&= x_{10}a\,\hat{\mathbf{x}} + y_{10}a\,\hat{\mathbf{y}} + z_{10}a\,\hat{\mathbf{z}} && (24d) && \text{O III}\\
\mathbf{B}_{146} &= \left(\tfrac{1}{2}-x_{10}\right)\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{10}\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_{10}\right)a\,\hat{\mathbf{x}} - y_{10}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{10}\right)a\,\hat{\mathbf{z}} && (24d) && \text{O III}\\
\mathbf{B}_{147} &= -x_{10}\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_{10}\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_{10}\right)\mathbf{a}_3 &&= -x_{10}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_{10}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_{10}\right)a\,\hat{\mathbf{z}} && (24d) && \text{O III}\\
\mathbf{B}_{148} &= \left(\tfrac{1}{2}+x_{10}\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_{10}\right)\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_{10}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_{10}\right)a\,\hat{\mathbf{y}} - z_{10}a\,\hat{\mathbf{z}} && (24d) && \text{O III}\\
\mathbf{B}_{149} &= z_{10}\,\mathbf{a}_1 + x_{10}\,\mathbf{a}_2 + y_{10}\,\mathbf{a}_3 &&= z_{10}a\,\hat{\mathbf{x}} + x_{10}a\,\hat{\mathbf{y}} + y_{10}a\,\hat{\mathbf{z}} && (24d) && \text{O III}\\
\mathbf{B}_{150} &= \left(\tfrac{1}{2}+z_{10}\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-x_{10}\right)\mathbf{a}_2 - y_{10}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+z_{10}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_{10}\right)a\,\hat{\mathbf{y}} - y_{10}a\,\hat{\mathbf{z}} && (24d) && \text{O III}\\
\mathbf{B}_{151} &= \left(\tfrac{1}{2}-z_{10}\right)\mathbf{a}_1 - x_{10}\,\mathbf{a}_2 + \left(\tfrac{1}{2}+y_{10}\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-z_{10}\right)a\,\hat{\mathbf{x}} - x_{10}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+y_{10}\right)a\,\hat{\mathbf{z}} && (24d) && \text{O III}\\
\mathbf{B}_{152} &= -z_{10}\,\mathbf{a}_1 + \left(\tfrac{1}{2}+x_{10}\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-y_{10}\right)\mathbf{a}_3 &&= -z_{10}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_{10}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-y_{10}\right)a\,\hat{\mathbf{z}} && (24d) && \text{O III}\\
\mathbf{B}_{153} &= y_{10}\,\mathbf{a}_1 + z_{10}\,\mathbf{a}_2 + x_{10}\,\mathbf{a}_3 &&= y_{10}a\,\hat{\mathbf{x}} + z_{10}a\,\hat{\mathbf{y}} + x_{10}a\,\hat{\mathbf{z}} && (24d) && \text{O III}\\
\mathbf{B}_{154} &= -y_{10}\,\mathbf{a}_1 + \left(\tfrac{1}{2}+z_{10}\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-x_{10}\right)\mathbf{a}_3 &&= -y_{10}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+z_{10}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-x_{10}\right)a\,\hat{\mathbf{z}} && (24d) && \text{O III}\\
\mathbf{B}_{155} &= \left(\tfrac{1}{2}+y_{10}\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-z_{10}\right)\mathbf{a}_2 - x_{10}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+y_{10}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-z_{10}\right)a\,\hat{\mathbf{y}} - x_{10}a\,\hat{\mathbf{z}} && (24d) && \text{O III}\\
\mathbf{B}_{156} &= \left(\tfrac{1}{2}-y_{10}\right)\mathbf{a}_1 - z_{10}\,\mathbf{a}_2 + \left(\tfrac{1}{2}+x_{10}\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-y_{10}\right)a\,\hat{\mathbf{x}} - z_{10}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_{10}\right)a\,\hat{\mathbf{z}} && (24d) && \text{O III}\\
\mathbf{B}_{157} &= -x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 &&= -x_{10}a\,\hat{\mathbf{x}} - y_{10}a\,\hat{\mathbf{y}} - z_{10}a\,\hat{\mathbf{z}} && (24d) && \text{O III}\\
\mathbf{B}_{158} &= \left(\tfrac{1}{2}+x_{10}\right)\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_{10}\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_{10}\right)a\,\hat{\mathbf{x}} + y_{10}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_{10}\right)a\,\hat{\mathbf{z}} && (24d) && \text{O III}\\
\mathbf{B}_{159} &= x_{10}\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_{10}\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_{10}\right)\mathbf{a}_3 &&= x_{10}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_{10}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_{10}\right)a\,\hat{\mathbf{z}} && (24d) && \text{O III}\\
\mathbf{B}_{160} &= \left(\tfrac{1}{2}-x_{10}\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_{10}\right)\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_{10}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_{10}\right)a\,\hat{\mathbf{y}} + z_{10}a\,\hat{\mathbf{z}} && (24d) && \text{O III}\\
\mathbf{B}_{161} &= -z_{10}\,\mathbf{a}_1 - x_{10}\,\mathbf{a}_2 - y_{10}\,\mathbf{a}_3 &&= -z_{10}a\,\hat{\mathbf{x}} - x_{10}a\,\hat{\mathbf{y}} - y_{10}a\,\hat{\mathbf{z}} && (24d) && \text{O III}\\
\mathbf{B}_{162} &= \left(\tfrac{1}{2}-z_{10}\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+x_{10}\right)\mathbf{a}_2 + y_{10}\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-z_{10}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_{10}\right)a\,\hat{\mathbf{y}} + y_{10}a\,\hat{\mathbf{z}} && (24d) && \text{O III}\\
\mathbf{B}_{163} &= \left(\tfrac{1}{2}+z_{10}\right)\mathbf{a}_1 + x_{10}\,\mathbf{a}_2 + \left(\tfrac{1}{2}-y_{10}\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+z_{10}\right)a\,\hat{\mathbf{x}} + x_{10}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-y_{10}\right)a\,\hat{\mathbf{z}} && (24d) && \text{O III}\\
\mathbf{B}_{164} &= z_{10}\,\mathbf{a}_1 + \left(\tfrac{1}{2}-x_{10}\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+y_{10}\right)\mathbf{a}_3 &&= z_{10}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_{10}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+y_{10}\right)a\,\hat{\mathbf{z}} && (24d) && \text{O III}\\
\mathbf{B}_{165} &= -y_{10}\,\mathbf{a}_1 - z_{10}\,\mathbf{a}_2 - x_{10}\,\mathbf{a}_3 &&= -y_{10}a\,\hat{\mathbf{x}} - z_{10}a\,\hat{\mathbf{y}} - x_{10}a\,\hat{\mathbf{z}} && (24d) && \text{O III}
\end{aligned}
$$

$\mathbf{B}_{166}$	=	$y_{10}\,\mathbf{a}_1+\left(\frac{1}{2}-z_{10}\right)\,\mathbf{a}_2+\left(\frac{1}{2}+x_{10}\right)\,\mathbf{a}_3$	=	$y_{10}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-z_{10}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{z}}$	(24*d*)	O III
$\mathbf{B}_{167}$	=	$\left(\frac{1}{2}-y_{10}\right)\,\mathbf{a}_1+\left(\frac{1}{2}+z_{10}\right)\,\mathbf{a}_2+x_{10}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+z_{10}\right)a\,\hat{\mathbf{y}}+x_{10}a\,\hat{\mathbf{z}}$	(24*d*)	O III
$\mathbf{B}_{168}$	=	$\left(\frac{1}{2}+y_{10}\right)\,\mathbf{a}_1+z_{10}\,\mathbf{a}_2+\left(\frac{1}{2}-x_{10}\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_{10}\right)a\,\hat{\mathbf{x}}+z_{10}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{z}}$	(24*d*)	O III
$\mathbf{B}_{169}$	=	$x_{11}\,\mathbf{a}_1+y_{11}\,\mathbf{a}_2+z_{11}\,\mathbf{a}_3$	=	$x_{11}a\,\hat{\mathbf{x}}+y_{11}a\,\hat{\mathbf{y}}+z_{11}a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{170}$	=	$\left(\frac{1}{2}-x_{11}\right)\,\mathbf{a}_1-y_{11}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{11}\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{11}\right)a\,\hat{\mathbf{x}}-y_{11}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{11}\right)a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{171}$	=	$-x_{11}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{11}\right)\,\mathbf{a}_2+\left(\frac{1}{2}-z_{11}\right)\,\mathbf{a}_3$	=	$-x_{11}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{11}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{11}\right)a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{172}$	=	$\left(\frac{1}{2}+x_{11}\right)\,\mathbf{a}_1+\left(\frac{1}{2}-y_{11}\right)\,\mathbf{a}_2-z_{11}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{11}\right)a\,\hat{\mathbf{y}}-z_{11}a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{173}$	=	$z_{11}\,\mathbf{a}_1+x_{11}\,\mathbf{a}_2+y_{11}\,\mathbf{a}_3$	=	$z_{11}a\,\hat{\mathbf{x}}+x_{11}a\,\hat{\mathbf{y}}+y_{11}a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{174}$	=	$\left(\frac{1}{2}+z_{11}\right)\,\mathbf{a}_1+\left(\frac{1}{2}-x_{11}\right)\,\mathbf{a}_2-y_{11}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+z_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_{11}\right)a\,\hat{\mathbf{y}}-y_{11}a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{175}$	=	$\left(\frac{1}{2}-z_{11}\right)\,\mathbf{a}_1-x_{11}\,\mathbf{a}_2+\left(\frac{1}{2}+y_{11}\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-z_{11}\right)a\,\hat{\mathbf{x}}-x_{11}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+y_{11}\right)a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{176}$	=	$-z_{11}\,\mathbf{a}_1+\left(\frac{1}{2}+x_{11}\right)\,\mathbf{a}_2+\left(\frac{1}{2}-y_{11}\right)\,\mathbf{a}_3$	=	$-z_{11}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_{11}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-y_{11}\right)a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{177}$	=	$y_{11}\,\mathbf{a}_1+z_{11}\,\mathbf{a}_2+x_{11}\,\mathbf{a}_3$	=	$y_{11}a\,\hat{\mathbf{x}}+z_{11}a\,\hat{\mathbf{y}}+x_{11}a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{178}$	=	$-y_{11}\,\mathbf{a}_1+\left(\frac{1}{2}+z_{11}\right)\,\mathbf{a}_2+\left(\frac{1}{2}-x_{11}\right)\,\mathbf{a}_3$	=	$-y_{11}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+z_{11}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-x_{11}\right)a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{179}$	=	$\left(\frac{1}{2}+y_{11}\right)\,\mathbf{a}_1+\left(\frac{1}{2}-z_{11}\right)\,\mathbf{a}_2-x_{11}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-z_{11}\right)a\,\hat{\mathbf{y}}-x_{11}a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{180}$	=	$\left(\frac{1}{2}-y_{11}\right)\,\mathbf{a}_1-z_{11}\,\mathbf{a}_2+\left(\frac{1}{2}+x_{11}\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_{11}\right)a\,\hat{\mathbf{x}}-z_{11}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+x_{11}\right)a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{181}$	=	$-x_{11}\,\mathbf{a}_1-y_{11}\,\mathbf{a}_2-z_{11}\,\mathbf{a}_3$	=	$-x_{11}a\,\hat{\mathbf{x}}-y_{11}a\,\hat{\mathbf{y}}-z_{11}a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{182}$	=	$\left(\frac{1}{2}+x_{11}\right)\,\mathbf{a}_1+y_{11}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{11}\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_{11}\right)a\,\hat{\mathbf{x}}+y_{11}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{11}\right)a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{183}$	=	$x_{11}\,\mathbf{a}_1+\left(\frac{1}{2}-y_{11}\right)\,\mathbf{a}_2+\left(\frac{1}{2}+z_{11}\right)\,\mathbf{a}_3$	=	$x_{11}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{11}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{11}\right)a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{184}$	=	$\left(\frac{1}{2}-x_{11}\right)\,\mathbf{a}_1+\left(\frac{1}{2}+y_{11}\right)\,\mathbf{a}_2+z_{11}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{11}\right)a\,\hat{\mathbf{y}}+z_{11}a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{185}$	=	$-z_{11}\,\mathbf{a}_1-x_{11}\,\mathbf{a}_2-y_{11}\,\mathbf{a}_3$	=	$-z_{11}a\,\hat{\mathbf{x}}-x_{11}a\,\hat{\mathbf{y}}-y_{11}a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{186}$	=	$\left(\frac{1}{2}-z_{11}\right)\,\mathbf{a}_1+\left(\frac{1}{2}+x_{11}\right)\,\mathbf{a}_2+y_{11}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-z_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_{11}\right)a\,\hat{\mathbf{y}}+y_{11}a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{187}$	=	$\left(\frac{1}{2}+z_{11}\right)\,\mathbf{a}_1+x_{11}\,\mathbf{a}_2+\left(\frac{1}{2}-y_{11}\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+z_{11}\right)a\,\hat{\mathbf{x}}+x_{11}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-y_{11}\right)a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{188}$	=	$z_{11}\,\mathbf{a}_1+\left(\frac{1}{2}-x_{11}\right)\,\mathbf{a}_2+\left(\frac{1}{2}+y_{11}\right)\,\mathbf{a}_3$	=	$z_{11}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_{11}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+y_{11}\right)a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{189}$	=	$-y_{11}\,\mathbf{a}_1-z_{11}\,\mathbf{a}_2-x_{11}\,\mathbf{a}_3$	=	$-y_{11}a\,\hat{\mathbf{x}}-z_{11}a\,\hat{\mathbf{y}}-x_{11}a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{190}$	=	$y_{11}\,\mathbf{a}_1+\left(\frac{1}{2}-z_{11}\right)\,\mathbf{a}_2+\left(\frac{1}{2}+x_{11}\right)\,\mathbf{a}_3$	=	$y_{11}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-z_{11}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+x_{11}\right)a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{191}$	=	$\left(\frac{1}{2}-y_{11}\right)\,\mathbf{a}_1+\left(\frac{1}{2}+z_{11}\right)\,\mathbf{a}_2+x_{11}\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+z_{11}\right)a\,\hat{\mathbf{y}}+x_{11}a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{192}$	=	$\left(\frac{1}{2}+y_{11}\right)\,\mathbf{a}_1+z_{11}\,\mathbf{a}_2+\left(\frac{1}{2}-x_{11}\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_{11}\right)a\,\hat{\mathbf{x}}+z_{11}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-x_{11}\right)a\,\hat{\mathbf{z}}$	(24*d*)	O IV

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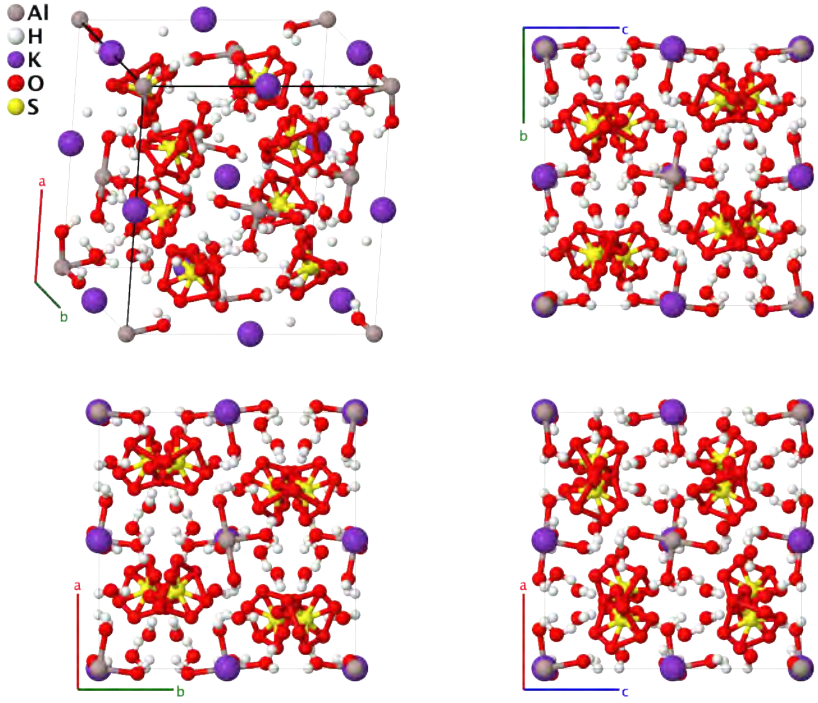
Geometry files:

- CIF: pp. 1783
- POSCAR: pp. 1784

α-Alum [$KAl(SO_4)_2 \cdot 12H_2O$, $H4_{13}$] Structure: AB24CD28E2_cP224_205_a_4d_b_2c4d_c

http://aflow.org/prototype-encyclopedia/AB24CD28E2_cP224_205_a_4d_b_2c4d_c



Prototype	:	$AlH_{24}KO_{20}S_2$
AFLOW prototype label	:	AB24CD28E2_cP224_205_a_4d_b_2c4d_c
***Strukturbericht* designation**	:	$H4_{13}$
Pearson symbol	:	cP224
Space group number	:	205
Space group symbol	:	$Pa\bar{3}$
AFLOW prototype command	:	`aflow --proto=AB24CD28E2_cP224_205_a_4d_b_2c4d_c --params=`$a, x_3, x_4, x_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}$

Other compounds with this structure

- $NH_4Al(SO_4)_2 \cdot 12H_2O$, $KCr(SO_4)_2 \cdot 12H_2O$, $RbAl(SO_4)_2 \cdot 12H_2O$, and $KAl(SeO_4)_2 \cdot 12H_2O$

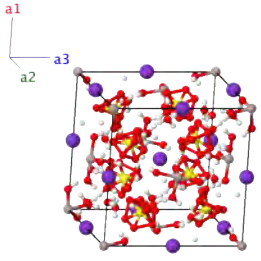
- The alums have the general formula $AB(XO_4)_2 \cdot 12H_2O$, where A is a monovalent ion, B is a trivalent ion, and X is a chalcogen. In most cases atom B is aluminum and atom X is sulfur, leading to the name alum.
- All alums have their room-temperature form in space group $Pa\bar{3}$ #205, but the bonding between the A and B ions and the XO_4 complex can be quite different.
- (Lipson, 1935ab) described three general forms of alum based on the sizes of the monovalent ions. Each of these forms was given a *Strukturbericht* designation by (Gottfried, 1937):
  - α-alum, with intermediate sized ions, prototype $KAl(SO_4)_2 \cdot 12H_2O$, $H4_{13}$ (this structure),

  - β-alum, with large ions, prototype $(NH_3CH_3)Al(SO_4)_2{\cdot}12H_2O$, $H4_{14}$, and
  - γ-alum, with small ions, prototype $NaAl(SO_4)_2{\cdot}12H_2O$, $H4_{15}$.

- This classification scheme is not complete, *e.g.*, (Ledsham, 1968) points out that $NaCr(SO_4)_2{\cdot}12H_2O$ does not fit into any of these categories, and that the actual structure depends on the combination of monovalent and trivalent ions.
- As noted above, the $Pa\bar{3}$ structures of alum are the room temperature form. As the temperature decreases the alum structure may transform. For example, in the temperature range 150-170 K the β-alum $(NH_3CH_3)Al(SO_4)_2{\cdot}12H_2O$ transforms into an orthorhombic structure with fully ordered NH_3CH_3 ions.
- (Ewald, 1931) designated the α-alum structure determined by (Cork, 1927) as $H4_2$, but when the superior structural determination of (Beevers, 1934) appeared this type was abandoned, and the new structure was given the designation $H4_{13}$. (Nyburg, 2000) were able to determine the positions of the hydrogen atoms, so we use their improved structure as our prototype.
- The oxygen atoms around the sulfur are statistically distributed in two ways: the probability of the O-I and O-III sites being occupied is 78.7%, while the probability of the O-II/O-IV combination being occupied is 21.3%.

Simple Cubic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = a\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(4a)$	Al
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	$(4a)$	Al
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	$(4a)$	Al
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$	$(4a)$	Al
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	$(4b)$	K
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{y}}$	$(4b)$	K
$\mathbf{B}_7$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	$(4b)$	K
$\mathbf{B}_8$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{z}}$	$(4b)$	K
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	$(8c)$	O I
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{z}}$	$(8c)$	O I
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{z}}$	$(8c)$	O I
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	$(8c)$	O I
$\mathbf{B}_{13}$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	$(8c)$	O I
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{z}}$	$(8c)$	O I
$\mathbf{B}_{15}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{z}}$	$(8c)$	O I
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	$(8c)$	O I

$$
\begin{aligned}
\mathbf{B}_{17} &= x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3 &&= x_4 a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}} + x_4 a\,\hat{\mathbf{z}} && (8c) && \text{O II}\\
\mathbf{B}_{18} &= \left(\tfrac{1}{2} - x_4\right)\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_4\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_4\right) a\,\hat{\mathbf{z}} && (8c) && \text{O II}\\
\mathbf{B}_{19} &= -x_4\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_4\right)\mathbf{a}_3 &&= -x_4 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_4\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_4\right) a\,\hat{\mathbf{z}} && (8c) && \text{O II}\\
\mathbf{B}_{20} &= \left(\tfrac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_4\right)\mathbf{a}_2 - x_4\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_4\right) a\,\hat{\mathbf{y}} - x_4 a\,\hat{\mathbf{z}} && (8c) && \text{O II}\\
\mathbf{B}_{21} &= -x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3 &&= -x_4 a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}} - x_4 a\,\hat{\mathbf{z}} && (8c) && \text{O II}\\
\mathbf{B}_{22} &= \left(\tfrac{1}{2} + x_4\right)\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_4\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_4\right) a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_4\right) a\,\hat{\mathbf{z}} && (8c) && \text{O II}\\
\mathbf{B}_{23} &= x_4\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_4\right)\mathbf{a}_3 &&= x_4 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_4\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_4\right) a\,\hat{\mathbf{z}} && (8c) && \text{O II}\\
\mathbf{B}_{24} &= \left(\tfrac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_4\right)\mathbf{a}_2 + x_4\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_4\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_4\right) a\,\hat{\mathbf{y}} + x_4 a\,\hat{\mathbf{z}} && (8c) && \text{O II}\\
\mathbf{B}_{25} &= x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3 &&= x_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}} && (8c) && \text{S}\\
\mathbf{B}_{26} &= \left(\tfrac{1}{2} - x_5\right)\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_5\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_5\right) a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_5\right) a\,\hat{\mathbf{z}} && (8c) && \text{S}\\
\mathbf{B}_{27} &= -x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_5\right)\mathbf{a}_3 &&= -x_5 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_5\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_5\right) a\,\hat{\mathbf{z}} && (8c) && \text{S}\\
\mathbf{B}_{28} &= \left(\tfrac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_5\right)\mathbf{a}_2 - x_5\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_5\right) a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}} && (8c) && \text{S}\\
\mathbf{B}_{29} &= -x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - x_5\,\mathbf{a}_3 &&= -x_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}} && (8c) && \text{S}\\
\mathbf{B}_{30} &= \left(\tfrac{1}{2} + x_5\right)\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_5\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_5\right) a\,\hat{\mathbf{z}} && (8c) && \text{S}\\
\mathbf{B}_{31} &= x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_5\right)\mathbf{a}_3 &&= x_5 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_5\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_5\right) a\,\hat{\mathbf{z}} && (8c) && \text{S}\\
\mathbf{B}_{32} &= \left(\tfrac{1}{2} - x_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_5\right)\mathbf{a}_2 + x_5\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_5\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_5\right) a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}} && (8c) && \text{S}\\
\mathbf{B}_{33} &= x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 &&= x_6 a\,\hat{\mathbf{x}} + y_6 a\,\hat{\mathbf{y}} + z_6 a\,\hat{\mathbf{z}} && (24d) && \text{H I}\\
\mathbf{B}_{34} &= \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} - y_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right) a\,\hat{\mathbf{z}} && (24d) && \text{H I}\\
\mathbf{B}_{35} &= -x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_3 &&= -x_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_6\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_6\right) a\,\hat{\mathbf{z}} && (24d) && \text{H I}\\
\mathbf{B}_{36} &= \left(\tfrac{1}{2} + x_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_6\right) a\,\hat{\mathbf{y}} - z_6 a\,\hat{\mathbf{z}} && (24d) && \text{H I}\\
\mathbf{B}_{37} &= z_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + y_6\,\mathbf{a}_3 &&= z_6 a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} + y_6 a\,\hat{\mathbf{z}} && (24d) && \text{H I}\\
\mathbf{B}_{38} &= \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_2 - y_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} + z_6\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_6\right) a\,\hat{\mathbf{y}} - y_6 a\,\hat{\mathbf{z}} && (24d) && \text{H I}\\
\mathbf{B}_{39} &= \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_1 - x_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} + y_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - z_6\right) a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + y_6\right) a\,\hat{\mathbf{z}} && (24d) && \text{H I}\\
\mathbf{B}_{40} &= -z_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - y_6\right)\mathbf{a}_3 &&= -z_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_6\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - y_6\right) a\,\hat{\mathbf{z}} && (24d) && \text{H I}\\
\mathbf{B}_{41} &= y_6\,\mathbf{a}_1 + z_6\,\mathbf{a}_2 + x_6\,\mathbf{a}_3 &&= y_6 a\,\hat{\mathbf{x}} + z_6 a\,\hat{\mathbf{y}} + x_6 a\,\hat{\mathbf{z}} && (24d) && \text{H I}\\
\mathbf{B}_{42} &= -y_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_3 &&= -y_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + z_6\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_6\right) a\,\hat{\mathbf{z}} && (24d) && \text{H I}\\
\mathbf{B}_{43} &= \left(\tfrac{1}{2} + y_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_2 - x_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} + y_6\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - z_6\right) a\,\hat{\mathbf{y}} - x_6 a\,\hat{\mathbf{z}} && (24d) && \text{H I}\\
\mathbf{B}_{44} &= \left(\tfrac{1}{2} - y_6\right)\mathbf{a}_1 - z_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - y_6\right) a\,\hat{\mathbf{x}} - z_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_6\right) a\,\hat{\mathbf{z}} && (24d) && \text{H I}\\
\mathbf{B}_{45} &= -x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3 &&= -x_6 a\,\hat{\mathbf{x}} - y_6 a\,\hat{\mathbf{y}} - z_6 a\,\hat{\mathbf{z}} && (24d) && \text{H I}\\
\mathbf{B}_{46} &= \left(\tfrac{1}{2} + x_6\right)\mathbf{a}_1 + y_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} + y_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_6\right) a\,\hat{\mathbf{z}} && (24d) && \text{H I}\\
\mathbf{B}_{47} &= x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 &&= x_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_6\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right) a\,\hat{\mathbf{z}} && (24d) && \text{H I}\\
\mathbf{B}_{48} &= \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_6\right)\mathbf{a}_2 + z_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_6\right) a\,\hat{\mathbf{y}} + z_6 a\,\hat{\mathbf{z}} && (24d) && \text{H I}\\
\mathbf{B}_{49} &= -z_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 - y_6\,\mathbf{a}_3 &&= -z_6 a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} - y_6 a\,\hat{\mathbf{z}} && (24d) && \text{H I}\\
\mathbf{B}_{50} &= \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_6\right)\mathbf{a}_2 + y_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} - z_6\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_6\right) a\,\hat{\mathbf{y}} + y_6 a\,\hat{\mathbf{z}} && (24d) && \text{H I}\\
\mathbf{B}_{51} &= \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_1 + x_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} - y_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} + z_6\right) a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - y_6\right) a\,\hat{\mathbf{z}} && (24d) && \text{H I}\\
\mathbf{B}_{52} &= z_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + y_6\right)\mathbf{a}_3 &&= z_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_6\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + y_6\right) a\,\hat{\mathbf{z}} && (24d) && \text{H I}
\end{aligned}
$$

$$
\begin{aligned}
\mathbf{B}_{53} &= -y_6\,\mathbf{a}_1 - z_6\,\mathbf{a}_2 - x_6\,\mathbf{a}_3 &&= -y_6 a\,\hat{\mathbf{x}} - z_6 a\,\hat{\mathbf{y}} - x_6 a\,\hat{\mathbf{z}} && (24d) && \text{H I}\\
\mathbf{B}_{54} &= y_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_6\right)\mathbf{a}_3 &&= y_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - z_6\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_6\right) a\,\hat{\mathbf{z}} && (24d) && \text{H I}\\
\mathbf{B}_{55} &= \left(\tfrac{1}{2} - y_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_2 + x_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} - y_6\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + z_6\right) a\,\hat{\mathbf{y}} + x_6 a\,\hat{\mathbf{z}} && (24d) && \text{H I}\\
\mathbf{B}_{56} &= \left(\tfrac{1}{2} + y_6\right)\mathbf{a}_1 + z_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} + y_6\right) a\,\hat{\mathbf{x}} + z_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_6\right) a\,\hat{\mathbf{z}} && (24d) && \text{H I}\\
\mathbf{B}_{57} &= x_7\,\mathbf{a}_1 + y_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 &&= x_7 a\,\hat{\mathbf{x}} + y_7 a\,\hat{\mathbf{y}} + z_7 a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{58} &= \left(\tfrac{1}{2} - x_7\right)\mathbf{a}_1 - y_7\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_7\right) a\,\hat{\mathbf{x}} - y_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_7\right) a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{59} &= -x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_7\right)\mathbf{a}_3 &&= -x_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_7\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_7\right) a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{60} &= \left(\tfrac{1}{2} + x_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_7\right)\mathbf{a}_2 - z_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_7\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_7\right) a\,\hat{\mathbf{y}} - z_7 a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{61} &= z_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + y_7\,\mathbf{a}_3 &&= z_7 a\,\hat{\mathbf{x}} + x_7 a\,\hat{\mathbf{y}} + y_7 a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{62} &= \left(\tfrac{1}{2} + z_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_7\right)\mathbf{a}_2 - y_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} + z_7\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_7\right) a\,\hat{\mathbf{y}} - y_7 a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{63} &= \left(\tfrac{1}{2} - z_7\right)\mathbf{a}_1 - x_7\,\mathbf{a}_2 + \left(\tfrac{1}{2} + y_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - z_7\right) a\,\hat{\mathbf{x}} - x_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + y_7\right) a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{64} &= -z_7\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - y_7\right)\mathbf{a}_3 &&= -z_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_7\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - y_7\right) a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{65} &= y_7\,\mathbf{a}_1 + z_7\,\mathbf{a}_2 + x_7\,\mathbf{a}_3 &&= y_7 a\,\hat{\mathbf{x}} + z_7 a\,\hat{\mathbf{y}} + x_7 a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{66} &= -y_7\,\mathbf{a}_1 + \left(\tfrac{1}{2} + z_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_7\right)\mathbf{a}_3 &&= -y_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + z_7\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_7\right) a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{67} &= \left(\tfrac{1}{2} + y_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - z_7\right)\mathbf{a}_2 - x_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} + y_7\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - z_7\right) a\,\hat{\mathbf{y}} - x_7 a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{68} &= \left(\tfrac{1}{2} - y_7\right)\mathbf{a}_1 - z_7\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - y_7\right) a\,\hat{\mathbf{x}} - z_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_7\right) a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{69} &= -x_7\,\mathbf{a}_1 - y_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3 &&= -x_7 a\,\hat{\mathbf{x}} - y_7 a\,\hat{\mathbf{y}} - z_7 a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{70} &= \left(\tfrac{1}{2} + x_7\right)\mathbf{a}_1 + y_7\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_7\right) a\,\hat{\mathbf{x}} + y_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_7\right) a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{71} &= x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_7\right)\mathbf{a}_3 &&= x_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_7\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_7\right) a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{72} &= \left(\tfrac{1}{2} - x_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_7\right)\mathbf{a}_2 + z_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_7\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_7\right) a\,\hat{\mathbf{y}} + z_7 a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{73} &= -z_7\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 - y_7\,\mathbf{a}_3 &&= -z_7 a\,\hat{\mathbf{x}} - x_7 a\,\hat{\mathbf{y}} - y_7 a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{74} &= \left(\tfrac{1}{2} - z_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_7\right)\mathbf{a}_2 + y_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} - z_7\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_7\right) a\,\hat{\mathbf{y}} + y_7 a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{75} &= \left(\tfrac{1}{2} + z_7\right)\mathbf{a}_1 + x_7\,\mathbf{a}_2 + \left(\tfrac{1}{2} - y_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} + z_7\right) a\,\hat{\mathbf{x}} + x_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - y_7\right) a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{76} &= z_7\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + y_7\right)\mathbf{a}_3 &&= z_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_7\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + y_7\right) a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{77} &= -y_7\,\mathbf{a}_1 - z_7\,\mathbf{a}_2 - x_7\,\mathbf{a}_3 &&= -y_7 a\,\hat{\mathbf{x}} - z_7 a\,\hat{\mathbf{y}} - x_7 a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{78} &= y_7\,\mathbf{a}_1 + \left(\tfrac{1}{2} - z_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_7\right)\mathbf{a}_3 &&= y_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - z_7\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_7\right) a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{79} &= \left(\tfrac{1}{2} - y_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + z_7\right)\mathbf{a}_2 + x_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} - y_7\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + z_7\right) a\,\hat{\mathbf{y}} + x_7 a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{80} &= \left(\tfrac{1}{2} + y_7\right)\mathbf{a}_1 + z_7\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} + y_7\right) a\,\hat{\mathbf{x}} + z_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_7\right) a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{81} &= x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 &&= x_8 a\,\hat{\mathbf{x}} + y_8 a\,\hat{\mathbf{y}} + z_8 a\,\hat{\mathbf{z}} && (24d) && \text{H III}\\
\mathbf{B}_{82} &= \left(\tfrac{1}{2} - x_8\right)\mathbf{a}_1 - y_8\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_8\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_8\right) a\,\hat{\mathbf{x}} - y_8 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_8\right) a\,\hat{\mathbf{z}} && (24d) && \text{H III}\\
\mathbf{B}_{83} &= -x_8\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_8\right)\mathbf{a}_3 &&= -x_8 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_8\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_8\right) a\,\hat{\mathbf{z}} && (24d) && \text{H III}\\
\mathbf{B}_{84} &= \left(\tfrac{1}{2} + x_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_8\right)\mathbf{a}_2 - z_8\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_8\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_8\right) a\,\hat{\mathbf{y}} - z_8 a\,\hat{\mathbf{z}} && (24d) && \text{H III}\\
\mathbf{B}_{85} &= z_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 + y_8\,\mathbf{a}_3 &&= z_8 a\,\hat{\mathbf{x}} + x_8 a\,\hat{\mathbf{y}} + y_8 a\,\hat{\mathbf{z}} && (24d) && \text{H III}\\
\mathbf{B}_{86} &= \left(\tfrac{1}{2} + z_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_8\right)\mathbf{a}_2 - y_8\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} + z_8\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_8\right) a\,\hat{\mathbf{y}} - y_8 a\,\hat{\mathbf{z}} && (24d) && \text{H III}\\
\mathbf{B}_{87} &= \left(\tfrac{1}{2} - z_8\right)\mathbf{a}_1 - x_8\,\mathbf{a}_2 + \left(\tfrac{1}{2} + y_8\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - z_8\right) a\,\hat{\mathbf{x}} - x_8 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + y_8\right) a\,\hat{\mathbf{z}} && (24d) && \text{H III}\\
\mathbf{B}_{88} &= -z_8\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - y_8\right)\mathbf{a}_3 &&= -z_8 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_8\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - y_8\right) a\,\hat{\mathbf{z}} && (24d) && \text{H III}
\end{aligned}
$$

$$\mathbf{B}_{89} = y_8\,\mathbf{a}_1 + z_8\,\mathbf{a}_2 + x_8\,\mathbf{a}_3 = y_8 a\,\hat{\mathbf{x}} + z_8 a\,\hat{\mathbf{y}} + x_8 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H III}$$

$$\mathbf{B}_{90} = -y_8\,\mathbf{a}_1 + \left(\tfrac{1}{2} + z_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_8\right)\mathbf{a}_3 = -y_8 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + z_8\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_8\right) a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H III}$$

$$\mathbf{B}_{91} = \left(\tfrac{1}{2} + y_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - z_8\right)\mathbf{a}_2 - x_8\,\mathbf{a}_3 = \left(\tfrac{1}{2} + y_8\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - z_8\right) a\,\hat{\mathbf{y}} - x_8 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H III}$$

$$\mathbf{B}_{92} = \left(\tfrac{1}{2} - y_8\right)\mathbf{a}_1 - z_8\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - y_8\right) a\,\hat{\mathbf{x}} - z_8 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_8\right) a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H III}$$

$$\mathbf{B}_{93} = -x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3 = -x_8 a\,\hat{\mathbf{x}} - y_8 a\,\hat{\mathbf{y}} - z_8 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H III}$$

$$\mathbf{B}_{94} = \left(\tfrac{1}{2} + x_8\right)\mathbf{a}_1 + y_8\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_8\right) a\,\hat{\mathbf{x}} + y_8 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_8\right) a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H III}$$

$$\mathbf{B}_{95} = x_8\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_8\right)\mathbf{a}_3 = x_8 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_8\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_8\right) a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H III}$$

$$\mathbf{B}_{96} = \left(\tfrac{1}{2} - x_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_8\right)\mathbf{a}_2 + z_8\,\mathbf{a}_3 = \left(\tfrac{1}{2} - x_8\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_8\right) a\,\hat{\mathbf{y}} + z_8 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H III}$$

$$\mathbf{B}_{97} = -z_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 - y_8\,\mathbf{a}_3 = -z_8 a\,\hat{\mathbf{x}} - x_8 a\,\hat{\mathbf{y}} - y_8 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H III}$$

$$\mathbf{B}_{98} = \left(\tfrac{1}{2} - z_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_8\right)\mathbf{a}_2 + y_8\,\mathbf{a}_3 = \left(\tfrac{1}{2} - z_8\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_8\right) a\,\hat{\mathbf{y}} + y_8 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H III}$$

$$\mathbf{B}_{99} = \left(\tfrac{1}{2} + z_8\right)\mathbf{a}_1 + x_8\,\mathbf{a}_2 + \left(\tfrac{1}{2} - y_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + z_8\right) a\,\hat{\mathbf{x}} + x_8 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - y_8\right) a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H III}$$

$$\mathbf{B}_{100} = z_8\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + y_8\right)\mathbf{a}_3 = z_8 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_8\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + y_8\right) a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H III}$$

$$\mathbf{B}_{101} = -y_8\,\mathbf{a}_1 - z_8\,\mathbf{a}_2 - x_8\,\mathbf{a}_3 = -y_8 a\,\hat{\mathbf{x}} - z_8 a\,\hat{\mathbf{y}} - x_8 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H III}$$

$$\mathbf{B}_{102} = y_8\,\mathbf{a}_1 + \left(\tfrac{1}{2} - z_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_8\right)\mathbf{a}_3 = y_8 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - z_8\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_8\right) a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H III}$$

$$\mathbf{B}_{103} = \left(\tfrac{1}{2} - y_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + z_8\right)\mathbf{a}_2 + x_8\,\mathbf{a}_3 = \left(\tfrac{1}{2} - y_8\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + z_8\right) a\,\hat{\mathbf{y}} + x_8 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H III}$$

$$\mathbf{B}_{104} = \left(\tfrac{1}{2} + y_8\right)\mathbf{a}_1 + z_8\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + y_8\right) a\,\hat{\mathbf{x}} + z_8 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_8\right) a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H III}$$

$$\mathbf{B}_{105} = x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 = x_9 a\,\hat{\mathbf{x}} + y_9 a\,\hat{\mathbf{y}} + z_9 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$\mathbf{B}_{106} = \left(\tfrac{1}{2} - x_9\right)\mathbf{a}_1 - y_9\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_9\right) a\,\hat{\mathbf{x}} - y_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_9\right) a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$\mathbf{B}_{107} = -x_9\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_9\right)\mathbf{a}_3 = -x_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_9\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_9\right) a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$\mathbf{B}_{108} = \left(\tfrac{1}{2} + x_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_9\right)\mathbf{a}_2 - z_9\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_9\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_9\right) a\,\hat{\mathbf{y}} - z_9 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$\mathbf{B}_{109} = z_9\,\mathbf{a}_1 + x_9\,\mathbf{a}_2 + y_9\,\mathbf{a}_3 = z_9 a\,\hat{\mathbf{x}} + x_9 a\,\hat{\mathbf{y}} + y_9 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$\mathbf{B}_{110} = \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_9\right)\mathbf{a}_2 - y_9\,\mathbf{a}_3 = \left(\tfrac{1}{2} + z_9\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_9\right) a\,\hat{\mathbf{y}} - y_9 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$\mathbf{B}_{111} = \left(\tfrac{1}{2} - z_9\right)\mathbf{a}_1 - x_9\,\mathbf{a}_2 + \left(\tfrac{1}{2} + y_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - z_9\right) a\,\hat{\mathbf{x}} - x_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + y_9\right) a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$\mathbf{B}_{112} = -z_9\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - y_9\right)\mathbf{a}_3 = -z_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_9\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - y_9\right) a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$\mathbf{B}_{113} = y_9\,\mathbf{a}_1 + z_9\,\mathbf{a}_2 + x_9\,\mathbf{a}_3 = y_9 a\,\hat{\mathbf{x}} + z_9 a\,\hat{\mathbf{y}} + x_9 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$\mathbf{B}_{114} = -y_9\,\mathbf{a}_1 + \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_9\right)\mathbf{a}_3 = -y_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + z_9\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_9\right) a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$\mathbf{B}_{115} = \left(\tfrac{1}{2} + y_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - z_9\right)\mathbf{a}_2 - x_9\,\mathbf{a}_3 = \left(\tfrac{1}{2} + y_9\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - z_9\right) a\,\hat{\mathbf{y}} - x_9 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$\mathbf{B}_{116} = \left(\tfrac{1}{2} - y_9\right)\mathbf{a}_1 - z_9\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - y_9\right) a\,\hat{\mathbf{x}} - z_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_9\right) a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$\mathbf{B}_{117} = -x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3 = -x_9 a\,\hat{\mathbf{x}} - y_9 a\,\hat{\mathbf{y}} - z_9 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$\mathbf{B}_{118} = \left(\tfrac{1}{2} + x_9\right)\mathbf{a}_1 + y_9\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_9\right) a\,\hat{\mathbf{x}} + y_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_9\right) a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$\mathbf{B}_{119} = x_9\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_3 = x_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_9\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_9\right) a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$\mathbf{B}_{120} = \left(\tfrac{1}{2} - x_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_9\right)\mathbf{a}_2 + z_9\,\mathbf{a}_3 = \left(\tfrac{1}{2} - x_9\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_9\right) a\,\hat{\mathbf{y}} + z_9 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$\mathbf{B}_{121} = -z_9\,\mathbf{a}_1 - x_9\,\mathbf{a}_2 - y_9\,\mathbf{a}_3 = -z_9 a\,\hat{\mathbf{x}} - x_9 a\,\hat{\mathbf{y}} - y_9 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$\mathbf{B}_{122} = \left(\tfrac{1}{2} - z_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_9\right)\mathbf{a}_2 + y_9\,\mathbf{a}_3 = \left(\tfrac{1}{2} - z_9\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_9\right) a\,\hat{\mathbf{y}} + y_9 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$\mathbf{B}_{123} = \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_1 + x_9\,\mathbf{a}_2 + \left(\tfrac{1}{2} - y_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + z_9\right) a\,\hat{\mathbf{x}} + x_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - y_9\right) a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$\mathbf{B}_{124} = z_9\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + y_9\right)\mathbf{a}_3 = z_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_9\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + y_9\right) a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$
\begin{array}{lcccll}
\mathbf{B}_{125} & = & -y_9\,\mathbf{a}_1 - z_9\,\mathbf{a}_2 - x_9\,\mathbf{a}_3 & = & -y_9 a\,\hat{\mathbf{x}} - z_9 a\,\hat{\mathbf{y}} - x_9 a\,\hat{\mathbf{z}} & (24d)\ \text{H IV} \\
\mathbf{B}_{126} & = & y_9\,\mathbf{a}_1 + \left(\tfrac{1}{2} - z_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_9\right)\mathbf{a}_3 & = & y_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - z_9\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_9\right) a\,\hat{\mathbf{z}} & (24d)\ \text{H IV} \\
\mathbf{B}_{127} & = & \left(\tfrac{1}{2} - y_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_2 + x_9\,\mathbf{a}_3 & = & \left(\tfrac{1}{2} - y_9\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + z_9\right) a\,\hat{\mathbf{y}} + x_9 a\,\hat{\mathbf{z}} & (24d)\ \text{H IV} \\
\mathbf{B}_{128} & = & \left(\tfrac{1}{2} + y_9\right)\mathbf{a}_1 + z_9\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_9\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} + y_9\right) a\,\hat{\mathbf{x}} + z_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_9\right) a\,\hat{\mathbf{z}} & (24d)\ \text{H IV} \\
\mathbf{B}_{129} & = & x_{10}\,\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 & = & x_{10} a\,\hat{\mathbf{x}} + y_{10} a\,\hat{\mathbf{y}} + z_{10} a\,\hat{\mathbf{z}} & (24d)\ \text{O III} \\
\mathbf{B}_{130} & = & \left(\tfrac{1}{2} - x_{10}\right)\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{10}\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} - x_{10}\right) a\,\hat{\mathbf{x}} - y_{10} a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{10}\right) a\,\hat{\mathbf{z}} & (24d)\ \text{O III} \\
\mathbf{B}_{131} & = & -x_{10}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{10}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{10}\right)\mathbf{a}_3 & = & -x_{10} a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{10}\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{10}\right) a\,\hat{\mathbf{z}} & (24d)\ \text{O III} \\
\mathbf{B}_{132} & = & \left(\tfrac{1}{2} + x_{10}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{10}\right)\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 & = & \left(\tfrac{1}{2} + x_{10}\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{10}\right) a\,\hat{\mathbf{y}} - z_{10} a\,\hat{\mathbf{z}} & (24d)\ \text{O III} \\
\mathbf{B}_{133} & = & z_{10}\,\mathbf{a}_1 + x_{10}\,\mathbf{a}_2 + y_{10}\,\mathbf{a}_3 & = & z_{10} a\,\hat{\mathbf{x}} + x_{10} a\,\hat{\mathbf{y}} + y_{10} a\,\hat{\mathbf{z}} & (24d)\ \text{O III} \\
\mathbf{B}_{134} & = & \left(\tfrac{1}{2} + z_{10}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_{10}\right)\mathbf{a}_2 - y_{10}\,\mathbf{a}_3 & = & \left(\tfrac{1}{2} + z_{10}\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_{10}\right) a\,\hat{\mathbf{y}} - y_{10} a\,\hat{\mathbf{z}} & (24d)\ \text{O III} \\
\mathbf{B}_{135} & = & \left(\tfrac{1}{2} - z_{10}\right)\mathbf{a}_1 - x_{10}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + y_{10}\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} - z_{10}\right) a\,\hat{\mathbf{x}} - x_{10} a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + y_{10}\right) a\,\hat{\mathbf{z}} & (24d)\ \text{O III} \\
\mathbf{B}_{136} & = & -z_{10}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_{10}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - y_{10}\right)\mathbf{a}_3 & = & -z_{10} a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_{10}\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - y_{10}\right) a\,\hat{\mathbf{z}} & (24d)\ \text{O III} \\
\mathbf{B}_{137} & = & y_{10}\,\mathbf{a}_1 + z_{10}\,\mathbf{a}_2 + x_{10}\,\mathbf{a}_3 & = & y_{10} a\,\hat{\mathbf{x}} + z_{10} a\,\hat{\mathbf{y}} + x_{10} a\,\hat{\mathbf{z}} & (24d)\ \text{O III} \\
\mathbf{B}_{138} & = & -y_{10}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + z_{10}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_{10}\right)\mathbf{a}_3 & = & -y_{10} a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + z_{10}\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_{10}\right) a\,\hat{\mathbf{z}} & (24d)\ \text{O III} \\
\mathbf{B}_{139} & = & \left(\tfrac{1}{2} + y_{10}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - z_{10}\right)\mathbf{a}_2 - x_{10}\,\mathbf{a}_3 & = & \left(\tfrac{1}{2} + y_{10}\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - z_{10}\right) a\,\hat{\mathbf{y}} - x_{10} a\,\hat{\mathbf{z}} & (24d)\ \text{O III} \\
\mathbf{B}_{140} & = & \left(\tfrac{1}{2} - y_{10}\right)\mathbf{a}_1 - z_{10}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_{10}\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} - y_{10}\right) a\,\hat{\mathbf{x}} - z_{10} a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_{10}\right) a\,\hat{\mathbf{z}} & (24d)\ \text{O III} \\
\mathbf{B}_{141} & = & -x_{10}\,\mathbf{a}_1 - y_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 & = & -x_{10} a\,\hat{\mathbf{x}} - y_{10} a\,\hat{\mathbf{y}} - z_{10} a\,\hat{\mathbf{z}} & (24d)\ \text{O III} \\
\mathbf{B}_{142} & = & \left(\tfrac{1}{2} + x_{10}\right)\mathbf{a}_1 + y_{10}\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{10}\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} + x_{10}\right) a\,\hat{\mathbf{x}} + y_{10} a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{10}\right) a\,\hat{\mathbf{z}} & (24d)\ \text{O III} \\
\mathbf{B}_{143} & = & x_{10}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{10}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{10}\right)\mathbf{a}_3 & = & x_{10} a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{10}\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{10}\right) a\,\hat{\mathbf{z}} & (24d)\ \text{O III} \\
\mathbf{B}_{144} & = & \left(\tfrac{1}{2} - x_{10}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{10}\right)\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 & = & \left(\tfrac{1}{2} - x_{10}\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{10}\right) a\,\hat{\mathbf{y}} + z_{10} a\,\hat{\mathbf{z}} & (24d)\ \text{O III} \\
\mathbf{B}_{145} & = & -z_{10}\,\mathbf{a}_1 - x_{10}\,\mathbf{a}_2 - y_{10}\,\mathbf{a}_3 & = & -z_{10} a\,\hat{\mathbf{x}} - x_{10} a\,\hat{\mathbf{y}} - y_{10} a\,\hat{\mathbf{z}} & (24d)\ \text{O III} \\
\mathbf{B}_{146} & = & \left(\tfrac{1}{2} - z_{10}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_{10}\right)\mathbf{a}_2 + y_{10}\,\mathbf{a}_3 & = & \left(\tfrac{1}{2} - z_{10}\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_{10}\right) a\,\hat{\mathbf{y}} + y_{10} a\,\hat{\mathbf{z}} & (24d)\ \text{O III} \\
\mathbf{B}_{147} & = & \left(\tfrac{1}{2} + z_{10}\right)\mathbf{a}_1 + x_{10}\,\mathbf{a}_2 + \left(\tfrac{1}{2} - y_{10}\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} + z_{10}\right) a\,\hat{\mathbf{x}} + x_{10} a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - y_{10}\right) a\,\hat{\mathbf{z}} & (24d)\ \text{O III} \\
\mathbf{B}_{148} & = & z_{10}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_{10}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + y_{10}\right)\mathbf{a}_3 & = & z_{10} a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_{10}\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + y_{10}\right) a\,\hat{\mathbf{z}} & (24d)\ \text{O III} \\
\mathbf{B}_{149} & = & -y_{10}\,\mathbf{a}_1 - z_{10}\,\mathbf{a}_2 - x_{10}\,\mathbf{a}_3 & = & -y_{10} a\,\hat{\mathbf{x}} - z_{10} a\,\hat{\mathbf{y}} - x_{10} a\,\hat{\mathbf{z}} & (24d)\ \text{O III} \\
\mathbf{B}_{150} & = & y_{10}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - z_{10}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_{10}\right)\mathbf{a}_3 & = & y_{10} a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - z_{10}\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_{10}\right) a\,\hat{\mathbf{z}} & (24d)\ \text{O III} \\
\mathbf{B}_{151} & = & \left(\tfrac{1}{2} - y_{10}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + z_{10}\right)\mathbf{a}_2 + x_{10}\,\mathbf{a}_3 & = & \left(\tfrac{1}{2} - y_{10}\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + z_{10}\right) a\,\hat{\mathbf{y}} + x_{10} a\,\hat{\mathbf{z}} & (24d)\ \text{O III} \\
\mathbf{B}_{152} & = & \left(\tfrac{1}{2} + y_{10}\right)\mathbf{a}_1 + z_{10}\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_{10}\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} + y_{10}\right) a\,\hat{\mathbf{x}} + z_{10} a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_{10}\right) a\,\hat{\mathbf{z}} & (24d)\ \text{O III} \\
\mathbf{B}_{153} & = & x_{11}\,\mathbf{a}_1 + y_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 & = & x_{11} a\,\hat{\mathbf{x}} + y_{11} a\,\hat{\mathbf{y}} + z_{11} a\,\hat{\mathbf{z}} & (24d)\ \text{O IV} \\
\mathbf{B}_{154} & = & \left(\tfrac{1}{2} - x_{11}\right)\mathbf{a}_1 - y_{11}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_{11}\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} - x_{11}\right) a\,\hat{\mathbf{x}} - y_{11} a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_{11}\right) a\,\hat{\mathbf{z}} & (24d)\ \text{O IV} \\
\mathbf{B}_{155} & = & -x_{11}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_{11}\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_{11}\right)\mathbf{a}_3 & = & -x_{11} a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_{11}\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_{11}\right) a\,\hat{\mathbf{z}} & (24d)\ \text{O IV} \\
\mathbf{B}_{156} & = & \left(\tfrac{1}{2} + x_{11}\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_{11}\right)\mathbf{a}_2 - z_{11}\,\mathbf{a}_3 & = & \left(\tfrac{1}{2} + x_{11}\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_{11}\right) a\,\hat{\mathbf{y}} - z_{11} a\,\hat{\mathbf{z}} & (24d)\ \text{O IV} \\
\mathbf{B}_{157} & = & z_{11}\,\mathbf{a}_1 + x_{11}\,\mathbf{a}_2 + y_{11}\,\mathbf{a}_3 & = & z_{11} a\,\hat{\mathbf{x}} + x_{11} a\,\hat{\mathbf{y}} + y_{11} a\,\hat{\mathbf{z}} & (24d)\ \text{O IV}
\end{array}
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\begin{array}{lcccccc}
\mathbf{B}_{158} & = & \left(\frac{1}{2}+z_{11}\right)\mathbf{a}_1+\left(\frac{1}{2}-x_{11}\right)\mathbf{a}_2-y_{11}\,\mathbf{a}_3 & = & \left(\frac{1}{2}+z_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_{11}\right)a\,\hat{\mathbf{y}}-y_{11}a\,\hat{\mathbf{z}} & (24d) & \text{O IV}\\
\mathbf{B}_{159} & = & \left(\frac{1}{2}-z_{11}\right)\mathbf{a}_1-x_{11}\,\mathbf{a}_2+\left(\frac{1}{2}+y_{11}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-z_{11}\right)a\,\hat{\mathbf{x}}-x_{11}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+y_{11}\right)a\,\hat{\mathbf{z}} & (24d) & \text{O IV}\\
\mathbf{B}_{160} & = & -z_{11}\,\mathbf{a}_1+\left(\frac{1}{2}+x_{11}\right)\mathbf{a}_2+\left(\frac{1}{2}-y_{11}\right)\mathbf{a}_3 & = & -z_{11}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_{11}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-y_{11}\right)a\,\hat{\mathbf{z}} & (24d) & \text{O IV}\\
\mathbf{B}_{161} & = & y_{11}\,\mathbf{a}_1+z_{11}\,\mathbf{a}_2+x_{11}\,\mathbf{a}_3 & = & y_{11}a\,\hat{\mathbf{x}}+z_{11}a\,\hat{\mathbf{y}}+x_{11}a\,\hat{\mathbf{z}} & (24d) & \text{O IV}\\
\mathbf{B}_{162} & = & -y_{11}\,\mathbf{a}_1+\left(\frac{1}{2}+z_{11}\right)\mathbf{a}_2+\left(\frac{1}{2}-x_{11}\right)\mathbf{a}_3 & = & -y_{11}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+z_{11}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-x_{11}\right)a\,\hat{\mathbf{z}} & (24d) & \text{O IV}\\
\mathbf{B}_{163} & = & \left(\frac{1}{2}+y_{11}\right)\mathbf{a}_1+\left(\frac{1}{2}-z_{11}\right)\mathbf{a}_2-x_{11}\,\mathbf{a}_3 & = & \left(\frac{1}{2}+y_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-z_{11}\right)a\,\hat{\mathbf{y}}-x_{11}a\,\hat{\mathbf{z}} & (24d) & \text{O IV}\\
\mathbf{B}_{164} & = & \left(\frac{1}{2}-y_{11}\right)\mathbf{a}_1-z_{11}\,\mathbf{a}_2+\left(\frac{1}{2}+x_{11}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-y_{11}\right)a\,\hat{\mathbf{x}}-z_{11}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+x_{11}\right)a\,\hat{\mathbf{z}} & (24d) & \text{O IV}\\
\mathbf{B}_{165} & = & -x_{11}\,\mathbf{a}_1-y_{11}\,\mathbf{a}_2-z_{11}\,\mathbf{a}_3 & = & -x_{11}a\,\hat{\mathbf{x}}-y_{11}a\,\hat{\mathbf{y}}-z_{11}a\,\hat{\mathbf{z}} & (24d) & \text{O IV}\\
\mathbf{B}_{166} & = & \left(\frac{1}{2}+x_{11}\right)\mathbf{a}_1+y_{11}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{11}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{11}\right)a\,\hat{\mathbf{x}}+y_{11}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{11}\right)a\,\hat{\mathbf{z}} & (24d) & \text{O IV}\\
\mathbf{B}_{167} & = & x_{11}\,\mathbf{a}_1+\left(\frac{1}{2}-y_{11}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{11}\right)\mathbf{a}_3 & = & x_{11}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{11}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{11}\right)a\,\hat{\mathbf{z}} & (24d) & \text{O IV}\\
\mathbf{B}_{168} & = & \left(\frac{1}{2}-x_{11}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{11}\right)\mathbf{a}_2+z_{11}\,\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{11}\right)a\,\hat{\mathbf{y}}+z_{11}a\,\hat{\mathbf{z}} & (24d) & \text{O IV}\\
\mathbf{B}_{169} & = & -z_{11}\,\mathbf{a}_1-x_{11}\,\mathbf{a}_2-y_{11}\,\mathbf{a}_3 & = & -z_{11}a\,\hat{\mathbf{x}}-x_{11}a\,\hat{\mathbf{y}}-y_{11}a\,\hat{\mathbf{z}} & (24d) & \text{O IV}\\
\mathbf{B}_{170} & = & \left(\frac{1}{2}-z_{11}\right)\mathbf{a}_1+\left(\frac{1}{2}+x_{11}\right)\mathbf{a}_2+y_{11}\,\mathbf{a}_3 & = & \left(\frac{1}{2}-z_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_{11}\right)a\,\hat{\mathbf{y}}+y_{11}a\,\hat{\mathbf{z}} & (24d) & \text{O IV}\\
\mathbf{B}_{171} & = & \left(\frac{1}{2}+z_{11}\right)\mathbf{a}_1+x_{11}\,\mathbf{a}_2+\left(\frac{1}{2}-y_{11}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+z_{11}\right)a\,\hat{\mathbf{x}}+x_{11}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-y_{11}\right)a\,\hat{\mathbf{z}} & (24d) & \text{O IV}\\
\mathbf{B}_{172} & = & z_{11}\,\mathbf{a}_1+\left(\frac{1}{2}-x_{11}\right)\mathbf{a}_2+\left(\frac{1}{2}+y_{11}\right)\mathbf{a}_3 & = & z_{11}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_{11}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+y_{11}\right)a\,\hat{\mathbf{z}} & (24d) & \text{O IV}\\
\mathbf{B}_{173} & = & -y_{11}\,\mathbf{a}_1-z_{11}\,\mathbf{a}_2-x_{11}\,\mathbf{a}_3 & = & -y_{11}a\,\hat{\mathbf{x}}-z_{11}a\,\hat{\mathbf{y}}-x_{11}a\,\hat{\mathbf{z}} & (24d) & \text{O IV}\\
\mathbf{B}_{174} & = & y_{11}\,\mathbf{a}_1+\left(\frac{1}{2}-z_{11}\right)\mathbf{a}_2+\left(\frac{1}{2}+x_{11}\right)\mathbf{a}_3 & = & y_{11}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-z_{11}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+x_{11}\right)a\,\hat{\mathbf{z}} & (24d) & \text{O IV}\\
\mathbf{B}_{175} & = & \left(\frac{1}{2}-y_{11}\right)\mathbf{a}_1+\left(\frac{1}{2}+z_{11}\right)\mathbf{a}_2+x_{11}\,\mathbf{a}_3 & = & \left(\frac{1}{2}-y_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+z_{11}\right)a\,\hat{\mathbf{y}}+x_{11}a\,\hat{\mathbf{z}} & (24d) & \text{O IV}\\
\mathbf{B}_{176} & = & \left(\frac{1}{2}+y_{11}\right)\mathbf{a}_1+z_{11}\,\mathbf{a}_2+\left(\frac{1}{2}-x_{11}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+y_{11}\right)a\,\hat{\mathbf{x}}+z_{11}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-x_{11}\right)a\,\hat{\mathbf{z}} & (24d) & \text{O IV}\\
\mathbf{B}_{177} & = & x_{12}\,\mathbf{a}_1+y_{12}\,\mathbf{a}_2+z_{12}\,\mathbf{a}_3 & = & x_{12}a\,\hat{\mathbf{x}}+y_{12}a\,\hat{\mathbf{y}}+z_{12}a\,\hat{\mathbf{z}} & (24d) & \text{O V}\\
\mathbf{B}_{178} & = & \left(\frac{1}{2}-x_{12}\right)\mathbf{a}_1-y_{12}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{12}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{12}\right)a\,\hat{\mathbf{x}}-y_{12}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{12}\right)a\,\hat{\mathbf{z}} & (24d) & \text{O V}\\
\mathbf{B}_{179} & = & -x_{12}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{12}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{12}\right)\mathbf{a}_3 & = & -x_{12}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{12}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{12}\right)a\,\hat{\mathbf{z}} & (24d) & \text{O V}\\
\mathbf{B}_{180} & = & \left(\frac{1}{2}+x_{12}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{12}\right)\mathbf{a}_2-z_{12}\,\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{12}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{12}\right)a\,\hat{\mathbf{y}}-z_{12}a\,\hat{\mathbf{z}} & (24d) & \text{O V}\\
\mathbf{B}_{181} & = & z_{12}\,\mathbf{a}_1+x_{12}\,\mathbf{a}_2+y_{12}\,\mathbf{a}_3 & = & z_{12}a\,\hat{\mathbf{x}}+x_{12}a\,\hat{\mathbf{y}}+y_{12}a\,\hat{\mathbf{z}} & (24d) & \text{O V}\\
\mathbf{B}_{182} & = & \left(\frac{1}{2}+z_{12}\right)\mathbf{a}_1+\left(\frac{1}{2}-x_{12}\right)\mathbf{a}_2-y_{12}\,\mathbf{a}_3 & = & \left(\frac{1}{2}+z_{12}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_{12}\right)a\,\hat{\mathbf{y}}-y_{12}a\,\hat{\mathbf{z}} & (24d) & \text{O V}\\
\mathbf{B}_{183} & = & \left(\frac{1}{2}-z_{12}\right)\mathbf{a}_1-x_{12}\,\mathbf{a}_2+\left(\frac{1}{2}+y_{12}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-z_{12}\right)a\,\hat{\mathbf{x}}-x_{12}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+y_{12}\right)a\,\hat{\mathbf{z}} & (24d) & \text{O V}\\
\mathbf{B}_{184} & = & -z_{12}\,\mathbf{a}_1+\left(\frac{1}{2}+x_{12}\right)\mathbf{a}_2+\left(\frac{1}{2}-y_{12}\right)\mathbf{a}_3 & = & -z_{12}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_{12}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-y_{12}\right)a\,\hat{\mathbf{z}} & (24d) & \text{O V}\\
\mathbf{B}_{185} & = & y_{12}\,\mathbf{a}_1+z_{12}\,\mathbf{a}_2+x_{12}\,\mathbf{a}_3 & = & y_{12}a\,\hat{\mathbf{x}}+z_{12}a\,\hat{\mathbf{y}}+x_{12}a\,\hat{\mathbf{z}} & (24d) & \text{O V}\\
\mathbf{B}_{186} & = & -y_{12}\,\mathbf{a}_1+\left(\frac{1}{2}+z_{12}\right)\mathbf{a}_2+\left(\frac{1}{2}-x_{12}\right)\mathbf{a}_3 & = & -y_{12}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+z_{12}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-x_{12}\right)a\,\hat{\mathbf{z}} & (24d) & \text{O V}\\
\mathbf{B}_{187} & = & \left(\frac{1}{2}+y_{12}\right)\mathbf{a}_1+\left(\frac{1}{2}-z_{12}\right)\mathbf{a}_2-x_{12}\,\mathbf{a}_3 & = & \left(\frac{1}{2}+y_{12}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-z_{12}\right)a\,\hat{\mathbf{y}}-x_{12}a\,\hat{\mathbf{z}} & (24d) & \text{O V}\\
\mathbf{B}_{188} & = & \left(\frac{1}{2}-y_{12}\right)\mathbf{a}_1-z_{12}\,\mathbf{a}_2+\left(\frac{1}{2}+x_{12}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-y_{12}\right)a\,\hat{\mathbf{x}}-z_{12}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+x_{12}\right)a\,\hat{\mathbf{z}} & (24d) & \text{O V}\\
\mathbf{B}_{189} & = & -x_{12}\,\mathbf{a}_1-y_{12}\,\mathbf{a}_2-z_{12}\,\mathbf{a}_3 & = & -x_{12}a\,\hat{\mathbf{x}}-y_{12}a\,\hat{\mathbf{y}}-z_{12}a\,\hat{\mathbf{z}} & (24d) & \text{O V}
\end{array}
$$

$$\mathbf{B}_{190} = \left(\tfrac{1}{2}+x_{12}\right)\mathbf{a}_1+y_{12}\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{12}\right)a\,\hat{\mathbf{x}}+y_{12}a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{12}\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O V}$$

$$\mathbf{B}_{191} = x_{12}\,\mathbf{a}_1+\left(\tfrac{1}{2}-y_{12}\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_{12}\right)\mathbf{a}_3 = x_{12}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{12}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{12}\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O V}$$

$$\mathbf{B}_{192} = \left(\tfrac{1}{2}-x_{12}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{12}\right)\mathbf{a}_2+z_{12}\,\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{12}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{12}\right)a\,\hat{\mathbf{y}}+z_{12}a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O V}$$

$$\mathbf{B}_{193} = -z_{12}\,\mathbf{a}_1-x_{12}\,\mathbf{a}_2-y_{12}\,\mathbf{a}_3 = -z_{12}a\,\hat{\mathbf{x}}-x_{12}a\,\hat{\mathbf{y}}-y_{12}a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O V}$$

$$\mathbf{B}_{194} = \left(\tfrac{1}{2}-z_{12}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_{12}\right)\mathbf{a}_2+y_{12}\,\mathbf{a}_3 = \left(\tfrac{1}{2}-z_{12}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_{12}\right)a\,\hat{\mathbf{y}}+y_{12}a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O V}$$

$$\mathbf{B}_{195} = \left(\tfrac{1}{2}+z_{12}\right)\mathbf{a}_1+x_{12}\,\mathbf{a}_2+\left(\tfrac{1}{2}-y_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+z_{12}\right)a\,\hat{\mathbf{x}}+x_{12}a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-y_{12}\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O V}$$

$$\mathbf{B}_{196} = z_{12}\,\mathbf{a}_1+\left(\tfrac{1}{2}-x_{12}\right)\mathbf{a}_2+\left(\tfrac{1}{2}+y_{12}\right)\mathbf{a}_3 = z_{12}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_{12}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+y_{12}\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O V}$$

$$\mathbf{B}_{197} = -y_{12}\,\mathbf{a}_1-z_{12}\,\mathbf{a}_2-x_{12}\,\mathbf{a}_3 = -y_{12}a\,\hat{\mathbf{x}}-z_{12}a\,\hat{\mathbf{y}}-x_{12}a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O V}$$

$$\mathbf{B}_{198} = y_{12}\,\mathbf{a}_1+\left(\tfrac{1}{2}-z_{12}\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_{12}\right)\mathbf{a}_3 = y_{12}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-z_{12}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+x_{12}\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O V}$$

$$\mathbf{B}_{199} = \left(\tfrac{1}{2}-y_{12}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+z_{12}\right)\mathbf{a}_2+x_{12}\,\mathbf{a}_3 = \left(\tfrac{1}{2}-y_{12}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+z_{12}\right)a\,\hat{\mathbf{y}}+x_{12}a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O V}$$

$$\mathbf{B}_{200} = \left(\tfrac{1}{2}+y_{12}\right)\mathbf{a}_1+z_{12}\,\mathbf{a}_2+\left(\tfrac{1}{2}-x_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+y_{12}\right)a\,\hat{\mathbf{x}}+z_{12}a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-x_{12}\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O V}$$

$$\mathbf{B}_{201} = x_{13}\,\mathbf{a}_1+y_{13}\,\mathbf{a}_2+z_{13}\,\mathbf{a}_3 = x_{13}a\,\hat{\mathbf{x}}+y_{13}a\,\hat{\mathbf{y}}+z_{13}a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O VI}$$

$$\mathbf{B}_{202} = \left(\tfrac{1}{2}-x_{13}\right)\mathbf{a}_1-y_{13}\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{13}\right)a\,\hat{\mathbf{x}}-y_{13}a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{13}\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O VI}$$

$$\mathbf{B}_{203} = -x_{13}\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_{13}\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_{13}\right)\mathbf{a}_3 = -x_{13}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{13}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{13}\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O VI}$$

$$\mathbf{B}_{204} = \left(\tfrac{1}{2}+x_{13}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{13}\right)\mathbf{a}_2-z_{13}\,\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{13}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{13}\right)a\,\hat{\mathbf{y}}-z_{13}a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O VI}$$

$$\mathbf{B}_{205} = z_{13}\,\mathbf{a}_1+x_{13}\,\mathbf{a}_2+y_{13}\,\mathbf{a}_3 = z_{13}a\,\hat{\mathbf{x}}+x_{13}a\,\hat{\mathbf{y}}+y_{13}a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O VI}$$

$$\mathbf{B}_{206} = \left(\tfrac{1}{2}+z_{13}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-x_{13}\right)\mathbf{a}_2-y_{13}\,\mathbf{a}_3 = \left(\tfrac{1}{2}+z_{13}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_{13}\right)a\,\hat{\mathbf{y}}-y_{13}a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O VI}$$

$$\mathbf{B}_{207} = \left(\tfrac{1}{2}-z_{13}\right)\mathbf{a}_1-x_{13}\,\mathbf{a}_2+\left(\tfrac{1}{2}+y_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-z_{13}\right)a\,\hat{\mathbf{x}}-x_{13}a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+y_{13}\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O VI}$$

$$\mathbf{B}_{208} = -z_{13}\,\mathbf{a}_1+\left(\tfrac{1}{2}+x_{13}\right)\mathbf{a}_2+\left(\tfrac{1}{2}-y_{13}\right)\mathbf{a}_3 = -z_{13}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_{13}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-y_{13}\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O VI}$$

$$\mathbf{B}_{209} = y_{13}\,\mathbf{a}_1+z_{13}\,\mathbf{a}_2+x_{13}\,\mathbf{a}_3 = y_{13}a\,\hat{\mathbf{x}}+z_{13}a\,\hat{\mathbf{y}}+x_{13}a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O VI}$$

$$\mathbf{B}_{210} = -y_{13}\,\mathbf{a}_1+\left(\tfrac{1}{2}+z_{13}\right)\mathbf{a}_2+\left(\tfrac{1}{2}-x_{13}\right)\mathbf{a}_3 = -y_{13}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+z_{13}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-x_{13}\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O VI}$$

$$\mathbf{B}_{211} = \left(\tfrac{1}{2}+y_{13}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-z_{13}\right)\mathbf{a}_2-x_{13}\,\mathbf{a}_3 = \left(\tfrac{1}{2}+y_{13}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-z_{13}\right)a\,\hat{\mathbf{y}}-x_{13}a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O VI}$$

$$\mathbf{B}_{212} = \left(\tfrac{1}{2}-y_{13}\right)\mathbf{a}_1-z_{13}\,\mathbf{a}_2+\left(\tfrac{1}{2}+x_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-y_{13}\right)a\,\hat{\mathbf{x}}-z_{13}a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+x_{13}\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O VI}$$

$$\mathbf{B}_{213} = -x_{13}\,\mathbf{a}_1-y_{13}\,\mathbf{a}_2-z_{13}\,\mathbf{a}_3 = -x_{13}a\,\hat{\mathbf{x}}-y_{13}a\,\hat{\mathbf{y}}-z_{13}a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O VI}$$

$$\mathbf{B}_{214} = \left(\tfrac{1}{2}+x_{13}\right)\mathbf{a}_1+y_{13}\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{13}\right)a\,\hat{\mathbf{x}}+y_{13}a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{13}\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O VI}$$

$$\mathbf{B}_{215} = x_{13}\,\mathbf{a}_1+\left(\tfrac{1}{2}-y_{13}\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_{13}\right)\mathbf{a}_3 = x_{13}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{13}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{13}\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O VI}$$

$$\mathbf{B}_{216} = \left(\tfrac{1}{2}-x_{13}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{13}\right)\mathbf{a}_2+z_{13}\,\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{13}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{13}\right)a\,\hat{\mathbf{y}}+z_{13}a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O VI}$$

$$\mathbf{B}_{217} = -z_{13}\,\mathbf{a}_1-x_{13}\,\mathbf{a}_2-y_{13}\,\mathbf{a}_3 = -z_{13}a\,\hat{\mathbf{x}}-x_{13}a\,\hat{\mathbf{y}}-y_{13}a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O VI}$$

$$\mathbf{B}_{218} = \left(\tfrac{1}{2}-z_{13}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_{13}\right)\mathbf{a}_2+y_{13}\,\mathbf{a}_3 = \left(\tfrac{1}{2}-z_{13}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_{13}\right)a\,\hat{\mathbf{y}}+y_{13}a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O VI}$$

$$\mathbf{B}_{219} = \left(\tfrac{1}{2}+z_{13}\right)\mathbf{a}_1+x_{13}\,\mathbf{a}_2+\left(\tfrac{1}{2}-y_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+z_{13}\right)a\,\hat{\mathbf{x}}+x_{13}a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-y_{13}\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O VI}$$

$$\mathbf{B}_{220} = z_{13}\,\mathbf{a}_1+\left(\tfrac{1}{2}-x_{13}\right)\mathbf{a}_2+\left(\tfrac{1}{2}+y_{13}\right)\mathbf{a}_3 = z_{13}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_{13}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+y_{13}\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O VI}$$

$$\mathbf{B}_{221} = -y_{13}\,\mathbf{a}_1-z_{13}\,\mathbf{a}_2-x_{13}\,\mathbf{a}_3 = -y_{13}a\,\hat{\mathbf{x}}-z_{13}a\,\hat{\mathbf{y}}-x_{13}a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O VI}$$

$$\mathbf{B}_{222} = y_{13}\,\mathbf{a}_1+\left(\tfrac{1}{2}-z_{13}\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_{13}\right)\mathbf{a}_3 = y_{13}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-z_{13}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+x_{13}\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O VI}$$

$$\mathbf{B}_{223} = \left(\tfrac{1}{2}-y_{13}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+z_{13}\right)\mathbf{a}_2+x_{13}\,\mathbf{a}_3 = \left(\tfrac{1}{2}-y_{13}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+z_{13}\right)a\,\hat{\mathbf{y}}+x_{13}a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O VI}$$

$$\mathbf{B}_{224} = \left(\tfrac{1}{2}+y_{13}\right)\mathbf{a}_1+z_{13}\,\mathbf{a}_2+\left(\tfrac{1}{2}-x_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+y_{13}\right)a\,\hat{\mathbf{x}}+z_{13}a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-x_{13}\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O VI}$$

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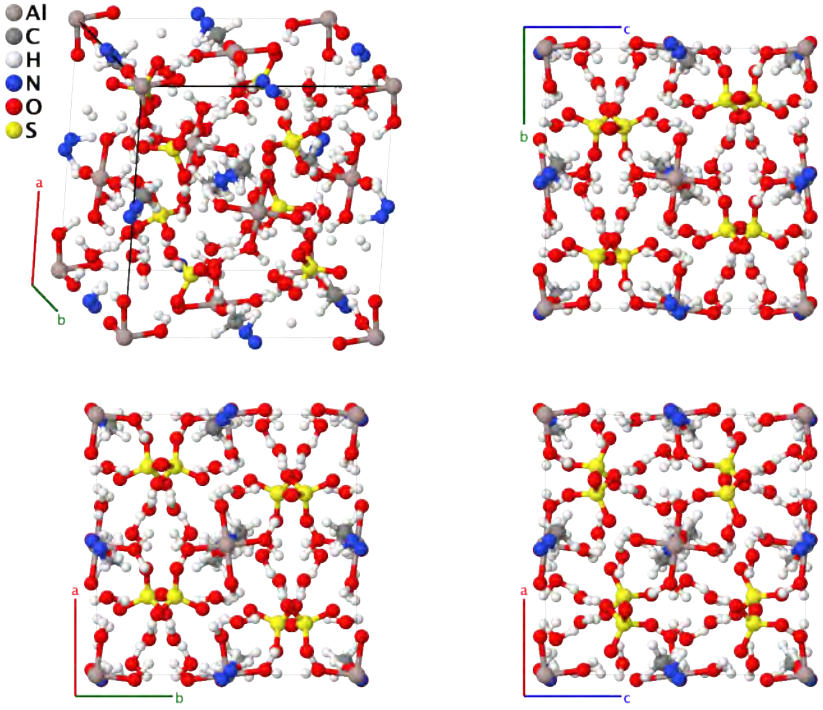
Geometry files:

- CIF: pp. 1785
- POSCAR: pp. 1785

β-Alum [$Al(NH_3CH_3)_2(SO_4)_2{\cdot}12H_2O$, $H4_{14}$] Structure: AB2C36D2E20F2_cP252_205_a_c_6d_c_c3d_c

http://aflow.org/prototype-encyclopedia/AB2C36D2E20F2_cP252_205_a_c_6d_c_c3d_c



Prototype	:	$AlC_2H_{36}N_2O_{20}S_2$
AFLOW prototype label	:	AB2C36D2E20F2_cP252_205_a_c_6d_c_c3d_c
***Strukturbericht* designation**	:	$H4_{14}$
Pearson symbol	:	cP252
Space group number	:	205
Space group symbol	:	$Pa\bar{3}$
AFLOW prototype command	:	`aflow --proto=AB2C36D2E20F2_cP252_205_a_c_6d_c_c3d_c --params=`$a, x_2, x_3, x_4, x_5, x_6, y_6, z_6, x_7, y_7, z_7, x_8, y_8, z_8, x_9, y_9, z_9, x_{10}, y_{10}, z_{10}, x_{11}, y_{11}, z_{11}, x_{12}, y_{12}, z_{12}, x_{13}, y_{13}, z_{13}, x_{14}, y_{14}, z_{14}$

Other compounds with this structure

- $AlCs(SO_4)_2{\cdot}12H_2O$

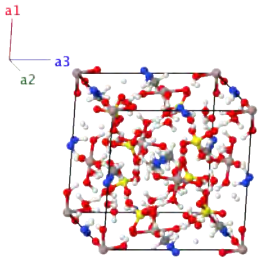
- The alums have the general formula $AB(XO_4)_2{\cdot}12H_2O$, where A is a monovalent ion, B is a trivalent ion, and X is a chalcogen. In most cases atom B is aluminum and atom X is sulfur, leading to the name alum.
- All alums have their room-temperature form in space group $Pa\bar{3}$ #205, but the bonding between the A and B ions and the XO_4 complex can be quite different.
- (Lipson, 1935ab) described three general forms of alum based on the sizes of the monovalent ions. Each of these forms was given a *Strukturbericht* designation by (Gottfried, 1937):
  - α-alum, with intermediate sized ions, prototype $KAl(SO_4)_2{\cdot}12H_2O$, $H4_{13}$,

  - β-alum, with large ions, prototype $(NH_3CH_3)Al(SO_4)_2{\cdot}12H_2O$, $H4_{14}$ (this structure), and
  - γ-alum, with small ions, prototype $NaAl(SO_4)_2{\cdot}12H_2O$, $H4_{15}$.

- This classification scheme is not complete, *e.g.*, (Ledsham, 1968) points out that $NaCr(SO_4)_2{\cdot}12H_2O$ does not fit into any of these categories, and that the actual structure depends on the combination of monovalent and trivalent ions.
- As noted above, the $Pa\overline{3}$ structures of alum are the room temperature form. As the temperature decreases the alum structure may transform. For example, in the temperature range 150-170 K the β-alum $(NH_3CH_3)Al(SO_4)_2{\cdot}12H_2O$ transforms into an orthorhombic structure with fully ordered NH_3CH_3 ions.
- This structure was originally determined by (Lipsom, 1935c), who could only determine that the NH_3CH_3 ion occupied the (4*b*) Wyckoff position. (Abdeen, 1981) showed that the ion was statistically distributed at two possible sites. The C-N bond distance is 1.4 Å, slightly smaller than the 1.51 Å distance observed in the low temperature structure. At any site, one of the two nitrogen positions is occupied, along with the carbon position 1.4 Å away. Six hydrogen positions from the (H-I) and (H-II) sites are then occupied.

Simple Cubic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= a\,\hat{\mathbf{x}}\\ \mathbf{a}_2 &= a\,\hat{\mathbf{y}}\\ \mathbf{a}_3 &= a\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(4*a*)	Al
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(4*a*)	Al
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(4*a*)	Al
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$	(4*a*)	Al
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$x_2a\,\hat{\mathbf{x}} + x_2a\,\hat{\mathbf{y}} + x_2a\,\hat{\mathbf{z}}$	(8*c*)	C
$\mathbf{B}_6$	=	$\left(\frac{1}{2}-x_2\right)\mathbf{a}_1 - x_2\,\mathbf{a}_2 + \left(\frac{1}{2}+x_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_2\right)a\,\hat{\mathbf{x}} - x_2a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+x_2\right)a\,\hat{\mathbf{z}}$	(8*c*)	C
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2}+x_2\right)\mathbf{a}_2 + \left(\frac{1}{2}-x_2\right)\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_2\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-x_2\right)a\,\hat{\mathbf{z}}$	(8*c*)	C
$\mathbf{B}_8$	=	$\left(\frac{1}{2}+x_2\right)\mathbf{a}_1 + \left(\frac{1}{2}-x_2\right)\mathbf{a}_2 - x_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_2\right)a\,\hat{\mathbf{y}} - x_2a\,\hat{\mathbf{z}}$	(8*c*)	C
$\mathbf{B}_9$	=	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 - x_2\,\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}} - x_2a\,\hat{\mathbf{y}} - x_2a\,\hat{\mathbf{z}}$	(8*c*)	C
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2}+x_2\right)\mathbf{a}_1 + x_2\,\mathbf{a}_2 + \left(\frac{1}{2}-x_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_2\right)a\,\hat{\mathbf{x}} + x_2a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-x_2\right)a\,\hat{\mathbf{z}}$	(8*c*)	C
$\mathbf{B}_{11}$	=	$x_2\,\mathbf{a}_1 + \left(\frac{1}{2}-x_2\right)\mathbf{a}_2 + \left(\frac{1}{2}+x_2\right)\mathbf{a}_3$	=	$x_2a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_2\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+x_2\right)a\,\hat{\mathbf{z}}$	(8*c*)	C
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2}-x_2\right)\mathbf{a}_1 + \left(\frac{1}{2}+x_2\right)\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_2\right)a\,\hat{\mathbf{y}} + x_2a\,\hat{\mathbf{z}}$	(8*c*)	C
$\mathbf{B}_{13}$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(8*c*)	N
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2}-x_3\right)\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\frac{1}{2}+x_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{z}}$	(8*c*)	N
$\mathbf{B}_{15}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2}+x_3\right)\mathbf{a}_2 + \left(\frac{1}{2}-x_3\right)\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{z}}$	(8*c*)	N
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2}+x_3\right)\mathbf{a}_1 + \left(\frac{1}{2}-x_3\right)\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(8*c*)	N
$\mathbf{B}_{17}$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(8*c*)	N

$$
\begin{array}{lcccccc}
\mathbf{B}_{18} & = & \left(\frac{1}{2}+x_3\right)\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \left(\frac{1}{2}-x_3\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{z}} & (8c) & \mathrm{N} \\
\mathbf{B}_{19} & = & x_3\,\mathbf{a}_1 + \left(\frac{1}{2}-x_3\right)\mathbf{a}_2 + \left(\frac{1}{2}+x_3\right)\mathbf{a}_3 & = & x_3 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{z}} & (8c) & \mathrm{N} \\
\mathbf{B}_{20} & = & \left(\frac{1}{2}-x_3\right)\mathbf{a}_1 + \left(\frac{1}{2}+x_3\right)\mathbf{a}_2 + x_3\,\mathbf{a}_3 & = & \left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{y}} + x_3 a\,\hat{\mathbf{z}} & (8c) & \mathrm{N} \\
\mathbf{B}_{21} & = & x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3 & = & x_4 a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}} + x_4 a\,\hat{\mathbf{z}} & (8c) & \mathrm{O\,I} \\
\mathbf{B}_{22} & = & \left(\frac{1}{2}-x_4\right)\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \left(\frac{1}{2}+x_4\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{z}} & (8c) & \mathrm{O\,I} \\
\mathbf{B}_{23} & = & -x_4\,\mathbf{a}_1 + \left(\frac{1}{2}+x_4\right)\mathbf{a}_2 + \left(\frac{1}{2}-x_4\right)\mathbf{a}_3 & = & -x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{z}} & (8c) & \mathrm{O\,I} \\
\mathbf{B}_{24} & = & \left(\frac{1}{2}+x_4\right)\mathbf{a}_1 + \left(\frac{1}{2}-x_4\right)\mathbf{a}_2 - x_4\,\mathbf{a}_3 & = & \left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{y}} - x_4 a\,\hat{\mathbf{z}} & (8c) & \mathrm{O\,I} \\
\mathbf{B}_{25} & = & -x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3 & = & -x_4 a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}} - x_4 a\,\hat{\mathbf{z}} & (8c) & \mathrm{O\,I} \\
\mathbf{B}_{26} & = & \left(\frac{1}{2}+x_4\right)\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \left(\frac{1}{2}-x_4\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{z}} & (8c) & \mathrm{O\,I} \\
\mathbf{B}_{27} & = & x_4\,\mathbf{a}_1 + \left(\frac{1}{2}-x_4\right)\mathbf{a}_2 + \left(\frac{1}{2}+x_4\right)\mathbf{a}_3 & = & x_4 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{z}} & (8c) & \mathrm{O\,I} \\
\mathbf{B}_{28} & = & \left(\frac{1}{2}-x_4\right)\mathbf{a}_1 + \left(\frac{1}{2}+x_4\right)\mathbf{a}_2 + x_4\,\mathbf{a}_3 & = & \left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{y}} + x_4 a\,\hat{\mathbf{z}} & (8c) & \mathrm{O\,I} \\
\mathbf{B}_{29} & = & x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3 & = & x_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}} & (8c) & \mathrm{S} \\
\mathbf{B}_{30} & = & \left(\frac{1}{2}-x_5\right)\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \left(\frac{1}{2}+x_5\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{z}} & (8c) & \mathrm{S} \\
\mathbf{B}_{31} & = & -x_5\,\mathbf{a}_1 + \left(\frac{1}{2}+x_5\right)\mathbf{a}_2 + \left(\frac{1}{2}-x_5\right)\mathbf{a}_3 & = & -x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{z}} & (8c) & \mathrm{S} \\
\mathbf{B}_{32} & = & \left(\frac{1}{2}+x_5\right)\mathbf{a}_1 + \left(\frac{1}{2}-x_5\right)\mathbf{a}_2 - x_5\,\mathbf{a}_3 & = & \left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}} & (8c) & \mathrm{S} \\
\mathbf{B}_{33} & = & -x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - x_5\,\mathbf{a}_3 & = & -x_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}} & (8c) & \mathrm{S} \\
\mathbf{B}_{34} & = & \left(\frac{1}{2}+x_5\right)\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \left(\frac{1}{2}-x_5\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{z}} & (8c) & \mathrm{S} \\
\mathbf{B}_{35} & = & x_5\,\mathbf{a}_1 + \left(\frac{1}{2}-x_5\right)\mathbf{a}_2 + \left(\frac{1}{2}+x_5\right)\mathbf{a}_3 & = & x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{z}} & (8c) & \mathrm{S} \\
\mathbf{B}_{36} & = & \left(\frac{1}{2}-x_5\right)\mathbf{a}_1 + \left(\frac{1}{2}+x_5\right)\mathbf{a}_2 + x_5\,\mathbf{a}_3 & = & \left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}} & (8c) & \mathrm{S} \\
\mathbf{B}_{37} & = & x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 & = & x_6 a\,\hat{\mathbf{x}} + y_6 a\,\hat{\mathbf{y}} + z_6 a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\,I} \\
\mathbf{B}_{38} & = & \left(\frac{1}{2}-x_6\right)\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\frac{1}{2}+z_6\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{x}} - y_6 a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_6\right)a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\,I} \\
\mathbf{B}_{39} & = & -x_6\,\mathbf{a}_1 + \left(\frac{1}{2}+y_6\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_6\right)\mathbf{a}_3 & = & -x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_6\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_6\right)a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\,I} \\
\mathbf{B}_{40} & = & \left(\frac{1}{2}+x_6\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3 & = & \left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_6\right)a\,\hat{\mathbf{y}} - z_6 a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\,I} \\
\mathbf{B}_{41} & = & z_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + y_6\,\mathbf{a}_3 & = & z_6 a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} + y_6 a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\,I} \\
\mathbf{B}_{42} & = & \left(\frac{1}{2}+z_6\right)\mathbf{a}_1 + \left(\frac{1}{2}-x_6\right)\mathbf{a}_2 - y_6\,\mathbf{a}_3 & = & \left(\frac{1}{2}+z_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{y}} - y_6 a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\,I} \\
\mathbf{B}_{43} & = & \left(\frac{1}{2}-z_6\right)\mathbf{a}_1 - x_6\,\mathbf{a}_2 + \left(\frac{1}{2}+y_6\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-z_6\right)a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+y_6\right)a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\,I} \\
\mathbf{B}_{44} & = & -z_6\,\mathbf{a}_1 + \left(\frac{1}{2}+x_6\right)\mathbf{a}_2 + \left(\frac{1}{2}-y_6\right)\mathbf{a}_3 & = & -z_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-y_6\right)a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\,I} \\
\mathbf{B}_{45} & = & y_6\,\mathbf{a}_1 + z_6\,\mathbf{a}_2 + x_6\,\mathbf{a}_3 & = & y_6 a\,\hat{\mathbf{x}} + z_6 a\,\hat{\mathbf{y}} + x_6 a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\,I} \\
\mathbf{B}_{46} & = & -y_6\,\mathbf{a}_1 + \left(\frac{1}{2}+z_6\right)\mathbf{a}_2 + \left(\frac{1}{2}-x_6\right)\mathbf{a}_3 & = & -y_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+z_6\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\,I} \\
\mathbf{B}_{47} & = & \left(\frac{1}{2}+y_6\right)\mathbf{a}_1 + \left(\frac{1}{2}-z_6\right)\mathbf{a}_2 - x_6\,\mathbf{a}_3 & = & \left(\frac{1}{2}+y_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-z_6\right)a\,\hat{\mathbf{y}} - x_6 a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\,I} \\
\mathbf{B}_{48} & = & \left(\frac{1}{2}-y_6\right)\mathbf{a}_1 - z_6\,\mathbf{a}_2 + \left(\frac{1}{2}+x_6\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-y_6\right)a\,\hat{\mathbf{x}} - z_6 a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\,I} \\
\mathbf{B}_{49} & = & -x_6\,\mathbf{a}_1 - y_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3 & = & -x_6 a\,\hat{\mathbf{x}} - y_6 a\,\hat{\mathbf{y}} - z_6 a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\,I} \\
\mathbf{B}_{50} & = & \left(\frac{1}{2}+x_6\right)\mathbf{a}_1 + y_6\,\mathbf{a}_2 + \left(\frac{1}{2}-z_6\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{x}} + y_6 a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_6\right)a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\,I} \\
\mathbf{B}_{51} & = & x_6\,\mathbf{a}_1 + \left(\frac{1}{2}-y_6\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_6\right)\mathbf{a}_3 & = & x_6 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_6\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_6\right)a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\,I} \\
\mathbf{B}_{52} & = & \left(\frac{1}{2}-x_6\right)\mathbf{a}_1 + \left(\frac{1}{2}+y_6\right)\mathbf{a}_2 + z_6\,\mathbf{a}_3 & = & \left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_6\right)a\,\hat{\mathbf{y}} + z_6 a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\,I} \\
\mathbf{B}_{53} & = & -z_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 - y_6\,\mathbf{a}_3 & = & -z_6 a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} - y_6 a\,\hat{\mathbf{z}} & (24d) & \mathrm{H\,I}
\end{array}
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\begin{aligned}
\mathbf{B}_{54} &= \left(\tfrac{1}{2}-z_6\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_6\right)\mathbf{a}_2+y_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-z_6\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{y}}+y_6a\,\hat{\mathbf{z}} && (24d) && \text{H I}\\
\mathbf{B}_{55} &= \left(\tfrac{1}{2}+z_6\right)\mathbf{a}_1+x_6\,\mathbf{a}_2+\left(\tfrac{1}{2}-y_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+z_6\right)a\,\hat{\mathbf{x}}+x_6a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-y_6\right)a\,\hat{\mathbf{z}} && (24d) && \text{H I}\\
\mathbf{B}_{56} &= z_6\,\mathbf{a}_1+\left(\tfrac{1}{2}-x_6\right)\mathbf{a}_2+\left(\tfrac{1}{2}+y_6\right)\mathbf{a}_3 &&= z_6a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+y_6\right)a\,\hat{\mathbf{z}} && (24d) && \text{H I}\\
\mathbf{B}_{57} &= -y_6\,\mathbf{a}_1-z_6\,\mathbf{a}_2-x_6\,\mathbf{a}_3 &&= -y_6a\,\hat{\mathbf{x}}-z_6a\,\hat{\mathbf{y}}-x_6a\,\hat{\mathbf{z}} && (24d) && \text{H I}\\
\mathbf{B}_{58} &= y_6\,\mathbf{a}_1+\left(\tfrac{1}{2}-z_6\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_6\right)\mathbf{a}_3 &&= y_6a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-z_6\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{z}} && (24d) && \text{H I}\\
\mathbf{B}_{59} &= \left(\tfrac{1}{2}-y_6\right)\mathbf{a}_1+\left(\tfrac{1}{2}+z_6\right)\mathbf{a}_2+x_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-y_6\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+z_6\right)a\,\hat{\mathbf{y}}+x_6a\,\hat{\mathbf{z}} && (24d) && \text{H I}\\
\mathbf{B}_{60} &= \left(\tfrac{1}{2}+y_6\right)\mathbf{a}_1+z_6\,\mathbf{a}_2+\left(\tfrac{1}{2}-x_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+y_6\right)a\,\hat{\mathbf{x}}+z_6a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{z}} && (24d) && \text{H I}\\
\mathbf{B}_{61} &= x_7\,\mathbf{a}_1+y_7\,\mathbf{a}_2+z_7\,\mathbf{a}_3 &&= x_7a\,\hat{\mathbf{x}}+y_7a\,\hat{\mathbf{y}}+z_7a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{62} &= \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_1-y_7\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{x}}-y_7a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{63} &= -x_7\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_7\right)\mathbf{a}_3 &&= -x_7a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_7\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_7\right)a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{64} &= \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_7\right)\mathbf{a}_2-z_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_7\right)a\,\hat{\mathbf{y}}-z_7a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{65} &= z_7\,\mathbf{a}_1+x_7\,\mathbf{a}_2+y_7\,\mathbf{a}_3 &&= z_7a\,\hat{\mathbf{x}}+x_7a\,\hat{\mathbf{y}}+y_7a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{66} &= \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}-x_7\right)\mathbf{a}_2-y_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{y}}-y_7a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{67} &= \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_1-x_7\,\mathbf{a}_2+\left(\tfrac{1}{2}+y_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-z_7\right)a\,\hat{\mathbf{x}}-x_7a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+y_7\right)a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{68} &= -z_7\,\mathbf{a}_1+\left(\tfrac{1}{2}+x_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}-y_7\right)\mathbf{a}_3 &&= -z_7a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-y_7\right)a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{69} &= y_7\,\mathbf{a}_1+z_7\,\mathbf{a}_2+x_7\,\mathbf{a}_3 &&= y_7a\,\hat{\mathbf{x}}+z_7a\,\hat{\mathbf{y}}+x_7a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{70} &= -y_7\,\mathbf{a}_1+\left(\tfrac{1}{2}+z_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}-x_7\right)\mathbf{a}_3 &&= -y_7a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{71} &= \left(\tfrac{1}{2}+y_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}-z_7\right)\mathbf{a}_2-x_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+y_7\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-z_7\right)a\,\hat{\mathbf{y}}-x_7a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{72} &= \left(\tfrac{1}{2}-y_7\right)\mathbf{a}_1-z_7\,\mathbf{a}_2+\left(\tfrac{1}{2}+x_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-y_7\right)a\,\hat{\mathbf{x}}-z_7a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{73} &= -x_7\,\mathbf{a}_1-y_7\,\mathbf{a}_2-z_7\,\mathbf{a}_3 &&= -x_7a\,\hat{\mathbf{x}}-y_7a\,\hat{\mathbf{y}}-z_7a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{74} &= \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1+y_7\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{x}}+y_7a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_7\right)a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{75} &= x_7\,\mathbf{a}_1+\left(\tfrac{1}{2}-y_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 &&= x_7a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_7\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{76} &= \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_7\right)\mathbf{a}_2+z_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_7\right)a\,\hat{\mathbf{y}}+z_7a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{77} &= -z_7\,\mathbf{a}_1-x_7\,\mathbf{a}_2-y_7\,\mathbf{a}_3 &&= -z_7a\,\hat{\mathbf{x}}-x_7a\,\hat{\mathbf{y}}-y_7a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{78} &= \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_7\right)\mathbf{a}_2+y_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-z_7\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{y}}+y_7a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{79} &= \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_1+x_7\,\mathbf{a}_2+\left(\tfrac{1}{2}-y_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{x}}+x_7a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-y_7\right)a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{80} &= z_7\,\mathbf{a}_1+\left(\tfrac{1}{2}-x_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}+y_7\right)\mathbf{a}_3 &&= z_7a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+y_7\right)a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{81} &= -y_7\,\mathbf{a}_1-z_7\,\mathbf{a}_2-x_7\,\mathbf{a}_3 &&= -y_7a\,\hat{\mathbf{x}}-z_7a\,\hat{\mathbf{y}}-x_7a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{82} &= y_7\,\mathbf{a}_1+\left(\tfrac{1}{2}-z_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_7\right)\mathbf{a}_3 &&= y_7a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-z_7\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{83} &= \left(\tfrac{1}{2}-y_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}+z_7\right)\mathbf{a}_2+x_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-y_7\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{y}}+x_7a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{84} &= \left(\tfrac{1}{2}+y_7\right)\mathbf{a}_1+z_7\,\mathbf{a}_2+\left(\tfrac{1}{2}-x_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+y_7\right)a\,\hat{\mathbf{x}}+z_7a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{z}} && (24d) && \text{H II}\\
\mathbf{B}_{85} &= x_8\,\mathbf{a}_1+y_8\,\mathbf{a}_2+z_8\,\mathbf{a}_3 &&= x_8a\,\hat{\mathbf{x}}+y_8a\,\hat{\mathbf{y}}+z_8a\,\hat{\mathbf{z}} && (24d) && \text{H III}\\
\mathbf{B}_{86} &= \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_1-y_8\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_8\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{x}}-y_8a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_8\right)a\,\hat{\mathbf{z}} && (24d) && \text{H III}\\
\mathbf{B}_{87} &= -x_8\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_8\right)\mathbf{a}_3 &&= -x_8a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_8\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_8\right)a\,\hat{\mathbf{z}} && (24d) && \text{H III}\\
\mathbf{B}_{88} &= \left(\tfrac{1}{2}+x_8\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_8\right)\mathbf{a}_2-z_8\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_8\right)a\,\hat{\mathbf{y}}-z_8a\,\hat{\mathbf{z}} && (24d) && \text{H III}\\
\mathbf{B}_{89} &= z_8\,\mathbf{a}_1+x_8\,\mathbf{a}_2+y_8\,\mathbf{a}_3 &&= z_8a\,\hat{\mathbf{x}}+x_8a\,\hat{\mathbf{y}}+y_8a\,\hat{\mathbf{z}} && (24d) && \text{H III}
\end{aligned}
$$

$$\mathbf{B}_{90} = \left(\tfrac{1}{2}+z_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_2 - y_8\,\mathbf{a}_3 = \left(\tfrac{1}{2}+z_8\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{y}} - y_8 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H III}$$

$$\mathbf{B}_{91} = \left(\tfrac{1}{2}-z_8\right)\mathbf{a}_1 - x_8\,\mathbf{a}_2 + \left(\tfrac{1}{2}+y_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-z_8\right)a\,\hat{\mathbf{x}} - x_8 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+y_8\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H III}$$

$$\mathbf{B}_{92} = -z_8\,\mathbf{a}_1 + \left(\tfrac{1}{2}+x_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-y_8\right)\mathbf{a}_3 = -z_8 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-y_8\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H III}$$

$$\mathbf{B}_{93} = y_8\,\mathbf{a}_1 + z_8\,\mathbf{a}_2 + x_8\,\mathbf{a}_3 = y_8 a\,\hat{\mathbf{x}} + z_8 a\,\hat{\mathbf{y}} + x_8 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H III}$$

$$\mathbf{B}_{94} = -y_8\,\mathbf{a}_1 + \left(\tfrac{1}{2}+z_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_3 = -y_8 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+z_8\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H III}$$

$$\mathbf{B}_{95} = \left(\tfrac{1}{2}+y_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-z_8\right)\mathbf{a}_2 - x_8\,\mathbf{a}_3 = \left(\tfrac{1}{2}+y_8\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-z_8\right)a\,\hat{\mathbf{y}} - x_8 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H III}$$

$$\mathbf{B}_{96} = \left(\tfrac{1}{2}-y_8\right)\mathbf{a}_1 - z_8\,\mathbf{a}_2 + \left(\tfrac{1}{2}+x_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-y_8\right)a\,\hat{\mathbf{x}} - z_8 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H III}$$

$$\mathbf{B}_{97} = -x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3 = -x_8 a\,\hat{\mathbf{x}} - y_8 a\,\hat{\mathbf{y}} - z_8 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H III}$$

$$\mathbf{B}_{98} = \left(\tfrac{1}{2}+x_8\right)\mathbf{a}_1 + y_8\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{x}} + y_8 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_8\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H III}$$

$$\mathbf{B}_{99} = x_8\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_8\right)\mathbf{a}_3 = x_8 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_8\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_8\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H III}$$

$$\mathbf{B}_{100} = \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_8\right)\mathbf{a}_2 + z_8\,\mathbf{a}_3 = \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_8\right)a\,\hat{\mathbf{y}} + z_8 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H III}$$

$$\mathbf{B}_{101} = -z_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 - y_8\,\mathbf{a}_3 = -z_8 a\,\hat{\mathbf{x}} - x_8 a\,\hat{\mathbf{y}} - y_8 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H III}$$

$$\mathbf{B}_{102} = \left(\tfrac{1}{2}-z_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+x_8\right)\mathbf{a}_2 + y_8\,\mathbf{a}_3 = \left(\tfrac{1}{2}-z_8\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{y}} + y_8 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H III}$$

$$\mathbf{B}_{103} = \left(\tfrac{1}{2}+z_8\right)\mathbf{a}_1 + x_8\,\mathbf{a}_2 + \left(\tfrac{1}{2}-y_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+z_8\right)a\,\hat{\mathbf{x}} + x_8 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-y_8\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H III}$$

$$\mathbf{B}_{104} = z_8\,\mathbf{a}_1 + \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+y_8\right)\mathbf{a}_3 = z_8 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+y_8\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H III}$$

$$\mathbf{B}_{105} = -y_8\,\mathbf{a}_1 - z_8\,\mathbf{a}_2 - x_8\,\mathbf{a}_3 = -y_8 a\,\hat{\mathbf{x}} - z_8 a\,\hat{\mathbf{y}} - x_8 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H III}$$

$$\mathbf{B}_{106} = y_8\,\mathbf{a}_1 + \left(\tfrac{1}{2}-z_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+x_8\right)\mathbf{a}_3 = y_8 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-z_8\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H III}$$

$$\mathbf{B}_{107} = \left(\tfrac{1}{2}-y_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+z_8\right)\mathbf{a}_2 + x_8\,\mathbf{a}_3 = \left(\tfrac{1}{2}-y_8\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+z_8\right)a\,\hat{\mathbf{y}} + x_8 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H III}$$

$$\mathbf{B}_{108} = \left(\tfrac{1}{2}+y_8\right)\mathbf{a}_1 + z_8\,\mathbf{a}_2 + \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+y_8\right)a\,\hat{\mathbf{x}} + z_8 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H III}$$

$$\mathbf{B}_{109} = x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 = x_9 a\,\hat{\mathbf{x}} + y_9 a\,\hat{\mathbf{y}} + z_9 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$\mathbf{B}_{110} = \left(\tfrac{1}{2}-x_9\right)\mathbf{a}_1 - y_9\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_9\right)a\,\hat{\mathbf{x}} - y_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_9\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$\mathbf{B}_{111} = -x_9\,\mathbf{a}_1 + \left(\tfrac{1}{2}+y_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_9\right)\mathbf{a}_3 = -x_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_9\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$\mathbf{B}_{112} = \left(\tfrac{1}{2}+x_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_9\right)\mathbf{a}_2 - z_9\,\mathbf{a}_3 = \left(\tfrac{1}{2}+x_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_9\right)a\,\hat{\mathbf{y}} - z_9 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$\mathbf{B}_{113} = z_9\,\mathbf{a}_1 + x_9\,\mathbf{a}_2 + y_9\,\mathbf{a}_3 = z_9 a\,\hat{\mathbf{x}} + x_9 a\,\hat{\mathbf{y}} + y_9 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$\mathbf{B}_{114} = \left(\tfrac{1}{2}+z_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-x_9\right)\mathbf{a}_2 - y_9\,\mathbf{a}_3 = \left(\tfrac{1}{2}+z_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_9\right)a\,\hat{\mathbf{y}} - y_9 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$\mathbf{B}_{115} = \left(\tfrac{1}{2}-z_9\right)\mathbf{a}_1 - x_9\,\mathbf{a}_2 + \left(\tfrac{1}{2}+y_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-z_9\right)a\,\hat{\mathbf{x}} - x_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+y_9\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$\mathbf{B}_{116} = -z_9\,\mathbf{a}_1 + \left(\tfrac{1}{2}+x_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-y_9\right)\mathbf{a}_3 = -z_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-y_9\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$\mathbf{B}_{117} = y_9\,\mathbf{a}_1 + z_9\,\mathbf{a}_2 + x_9\,\mathbf{a}_3 = y_9 a\,\hat{\mathbf{x}} + z_9 a\,\hat{\mathbf{y}} + x_9 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$\mathbf{B}_{118} = -y_9\,\mathbf{a}_1 + \left(\tfrac{1}{2}+z_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-x_9\right)\mathbf{a}_3 = -y_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+z_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-x_9\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$\mathbf{B}_{119} = \left(\tfrac{1}{2}+y_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-z_9\right)\mathbf{a}_2 - x_9\,\mathbf{a}_3 = \left(\tfrac{1}{2}+y_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-z_9\right)a\,\hat{\mathbf{y}} - x_9 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$\mathbf{B}_{120} = \left(\tfrac{1}{2}-y_9\right)\mathbf{a}_1 - z_9\,\mathbf{a}_2 + \left(\tfrac{1}{2}+x_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-y_9\right)a\,\hat{\mathbf{x}} - z_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_9\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$\mathbf{B}_{121} = -x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3 = -x_9 a\,\hat{\mathbf{x}} - y_9 a\,\hat{\mathbf{y}} - z_9 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$\mathbf{B}_{122} = \left(\tfrac{1}{2}+x_9\right)\mathbf{a}_1 + y_9\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_9\right)a\,\hat{\mathbf{x}} + y_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_9\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$\mathbf{B}_{123} = x_9\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_9\right)\mathbf{a}_3 = x_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_9\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$\mathbf{B}_{124} = \left(\tfrac{1}{2}-x_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_9\right)\mathbf{a}_2 + z_9\,\mathbf{a}_3 = \left(\tfrac{1}{2}-x_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+y_9\right)a\,\hat{\mathbf{y}} + z_9 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$\mathbf{B}_{125} = -z_9\,\mathbf{a}_1 - x_9\,\mathbf{a}_2 - y_9\,\mathbf{a}_3 = -z_9 a\,\hat{\mathbf{x}} - x_9 a\,\hat{\mathbf{y}} - y_9 a\,\hat{\mathbf{z}} \quad (24d) \quad \text{H IV}$$

$$
\begin{array}{lcccccc}
\mathbf{B}_{126} & = & \left(\frac{1}{2}-z_9\right)\mathbf{a}_1+\left(\frac{1}{2}+x_9\right)\mathbf{a}_2+y_9\,\mathbf{a}_3 & = & \left(\frac{1}{2}-z_9\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_9\right)a\,\hat{\mathbf{y}}+y_9a\,\hat{\mathbf{z}} & (24d) & \text{H IV}\\
\mathbf{B}_{127} & = & \left(\frac{1}{2}+z_9\right)\mathbf{a}_1+x_9\,\mathbf{a}_2+\left(\frac{1}{2}-y_9\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+z_9\right)a\,\hat{\mathbf{x}}+x_9a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-y_9\right)a\,\hat{\mathbf{z}} & (24d) & \text{H IV}\\
\mathbf{B}_{128} & = & z_9\,\mathbf{a}_1+\left(\frac{1}{2}-x_9\right)\mathbf{a}_2+\left(\frac{1}{2}+y_9\right)\mathbf{a}_3 & = & z_9a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_9\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+y_9\right)a\,\hat{\mathbf{z}} & (24d) & \text{H IV}\\
\mathbf{B}_{129} & = & -y_9\,\mathbf{a}_1-z_9\,\mathbf{a}_2-x_9\,\mathbf{a}_3 & = & -y_9a\,\hat{\mathbf{x}}-z_9a\,\hat{\mathbf{y}}-x_9a\,\hat{\mathbf{z}} & (24d) & \text{H IV}\\
\mathbf{B}_{130} & = & y_9\,\mathbf{a}_1+\left(\frac{1}{2}-z_9\right)\mathbf{a}_2+\left(\frac{1}{2}+x_9\right)\mathbf{a}_3 & = & y_9a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-z_9\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+x_9\right)a\,\hat{\mathbf{z}} & (24d) & \text{H IV}\\
\mathbf{B}_{131} & = & \left(\frac{1}{2}-y_9\right)\mathbf{a}_1+\left(\frac{1}{2}+z_9\right)\mathbf{a}_2+x_9\,\mathbf{a}_3 & = & \left(\frac{1}{2}-y_9\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+z_9\right)a\,\hat{\mathbf{y}}+x_9a\,\hat{\mathbf{z}} & (24d) & \text{H IV}\\
\mathbf{B}_{132} & = & \left(\frac{1}{2}+y_9\right)\mathbf{a}_1+z_9\,\mathbf{a}_2+\left(\frac{1}{2}-x_9\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+y_9\right)a\,\hat{\mathbf{x}}+z_9a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-x_9\right)a\,\hat{\mathbf{z}} & (24d) & \text{H IV}\\
\mathbf{B}_{133} & = & x_{10}\,\mathbf{a}_1+y_{10}\,\mathbf{a}_2+z_{10}\,\mathbf{a}_3 & = & x_{10}a\,\hat{\mathbf{x}}+y_{10}a\,\hat{\mathbf{y}}+z_{10}a\,\hat{\mathbf{z}} & (24d) & \text{H V}\\
\mathbf{B}_{134} & = & \left(\frac{1}{2}-x_{10}\right)\mathbf{a}_1-y_{10}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{10}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{x}}-y_{10}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{10}\right)a\,\hat{\mathbf{z}} & (24d) & \text{H V}\\
\mathbf{B}_{135} & = & -x_{10}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{10}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{10}\right)\mathbf{a}_3 & = & -x_{10}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{10}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{10}\right)a\,\hat{\mathbf{z}} & (24d) & \text{H V}\\
\mathbf{B}_{136} & = & \left(\frac{1}{2}+x_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{10}\right)\mathbf{a}_2-z_{10}\,\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{10}\right)a\,\hat{\mathbf{y}}-z_{10}a\,\hat{\mathbf{z}} & (24d) & \text{H V}\\
\mathbf{B}_{137} & = & z_{10}\,\mathbf{a}_1+x_{10}\,\mathbf{a}_2+y_{10}\,\mathbf{a}_3 & = & z_{10}a\,\hat{\mathbf{x}}+x_{10}a\,\hat{\mathbf{y}}+y_{10}a\,\hat{\mathbf{z}} & (24d) & \text{H V}\\
\mathbf{B}_{138} & = & \left(\frac{1}{2}+z_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}-x_{10}\right)\mathbf{a}_2-y_{10}\,\mathbf{a}_3 & = & \left(\frac{1}{2}+z_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{y}}-y_{10}a\,\hat{\mathbf{z}} & (24d) & \text{H V}\\
\mathbf{B}_{139} & = & \left(\frac{1}{2}-z_{10}\right)\mathbf{a}_1-x_{10}\,\mathbf{a}_2+\left(\frac{1}{2}+y_{10}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-z_{10}\right)a\,\hat{\mathbf{x}}-x_{10}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+y_{10}\right)a\,\hat{\mathbf{z}} & (24d) & \text{H V}\\
\mathbf{B}_{140} & = & -z_{10}\,\mathbf{a}_1+\left(\frac{1}{2}+x_{10}\right)\mathbf{a}_2+\left(\frac{1}{2}-y_{10}\right)\mathbf{a}_3 & = & -z_{10}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-y_{10}\right)a\,\hat{\mathbf{z}} & (24d) & \text{H V}\\
\mathbf{B}_{141} & = & y_{10}\,\mathbf{a}_1+z_{10}\,\mathbf{a}_2+x_{10}\,\mathbf{a}_3 & = & y_{10}a\,\hat{\mathbf{x}}+z_{10}a\,\hat{\mathbf{y}}+x_{10}a\,\hat{\mathbf{z}} & (24d) & \text{H V}\\
\mathbf{B}_{142} & = & -y_{10}\,\mathbf{a}_1+\left(\frac{1}{2}+z_{10}\right)\mathbf{a}_2+\left(\frac{1}{2}-x_{10}\right)\mathbf{a}_3 & = & -y_{10}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+z_{10}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{z}} & (24d) & \text{H V}\\
\mathbf{B}_{143} & = & \left(\frac{1}{2}+y_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}-z_{10}\right)\mathbf{a}_2-x_{10}\,\mathbf{a}_3 & = & \left(\frac{1}{2}+y_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-z_{10}\right)a\,\hat{\mathbf{y}}-x_{10}a\,\hat{\mathbf{z}} & (24d) & \text{H V}\\
\mathbf{B}_{144} & = & \left(\frac{1}{2}-y_{10}\right)\mathbf{a}_1-z_{10}\,\mathbf{a}_2+\left(\frac{1}{2}+x_{10}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-y_{10}\right)a\,\hat{\mathbf{x}}-z_{10}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{z}} & (24d) & \text{H V}\\
\mathbf{B}_{145} & = & -x_{10}\,\mathbf{a}_1-y_{10}\,\mathbf{a}_2-z_{10}\,\mathbf{a}_3 & = & -x_{10}a\,\hat{\mathbf{x}}-y_{10}a\,\hat{\mathbf{y}}-z_{10}a\,\hat{\mathbf{z}} & (24d) & \text{H V}\\
\mathbf{B}_{146} & = & \left(\frac{1}{2}+x_{10}\right)\mathbf{a}_1+y_{10}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{10}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{x}}+y_{10}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{10}\right)a\,\hat{\mathbf{z}} & (24d) & \text{H V}\\
\mathbf{B}_{147} & = & x_{10}\,\mathbf{a}_1+\left(\frac{1}{2}-y_{10}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{10}\right)\mathbf{a}_3 & = & x_{10}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{10}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{10}\right)a\,\hat{\mathbf{z}} & (24d) & \text{H V}\\
\mathbf{B}_{148} & = & \left(\frac{1}{2}-x_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{10}\right)\mathbf{a}_2+z_{10}\,\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{10}\right)a\,\hat{\mathbf{y}}+z_{10}a\,\hat{\mathbf{z}} & (24d) & \text{H V}\\
\mathbf{B}_{149} & = & -z_{10}\,\mathbf{a}_1-x_{10}\,\mathbf{a}_2-y_{10}\,\mathbf{a}_3 & = & -z_{10}a\,\hat{\mathbf{x}}-x_{10}a\,\hat{\mathbf{y}}-y_{10}a\,\hat{\mathbf{z}} & (24d) & \text{H V}\\
\mathbf{B}_{150} & = & \left(\frac{1}{2}-z_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}+x_{10}\right)\mathbf{a}_2+y_{10}\,\mathbf{a}_3 & = & \left(\frac{1}{2}-z_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{y}}+y_{10}a\,\hat{\mathbf{z}} & (24d) & \text{H V}\\
\mathbf{B}_{151} & = & \left(\frac{1}{2}+z_{10}\right)\mathbf{a}_1+x_{10}\,\mathbf{a}_2+\left(\frac{1}{2}-y_{10}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+z_{10}\right)a\,\hat{\mathbf{x}}+x_{10}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-y_{10}\right)a\,\hat{\mathbf{z}} & (24d) & \text{H V}\\
\mathbf{B}_{152} & = & z_{10}\,\mathbf{a}_1+\left(\frac{1}{2}-x_{10}\right)\mathbf{a}_2+\left(\frac{1}{2}+y_{10}\right)\mathbf{a}_3 & = & z_{10}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+y_{10}\right)a\,\hat{\mathbf{z}} & (24d) & \text{H V}\\
\mathbf{B}_{153} & = & -y_{10}\,\mathbf{a}_1-z_{10}\,\mathbf{a}_2-x_{10}\,\mathbf{a}_3 & = & -y_{10}a\,\hat{\mathbf{x}}-z_{10}a\,\hat{\mathbf{y}}-x_{10}a\,\hat{\mathbf{z}} & (24d) & \text{H V}\\
\mathbf{B}_{154} & = & y_{10}\,\mathbf{a}_1+\left(\frac{1}{2}-z_{10}\right)\mathbf{a}_2+\left(\frac{1}{2}+x_{10}\right)\mathbf{a}_3 & = & y_{10}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-z_{10}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+x_{10}\right)a\,\hat{\mathbf{z}} & (24d) & \text{H V}\\
\mathbf{B}_{155} & = & \left(\frac{1}{2}-y_{10}\right)\mathbf{a}_1+\left(\frac{1}{2}+z_{10}\right)\mathbf{a}_2+x_{10}\,\mathbf{a}_3 & = & \left(\frac{1}{2}-y_{10}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+z_{10}\right)a\,\hat{\mathbf{y}}+x_{10}a\,\hat{\mathbf{z}} & (24d) & \text{H V}\\
\mathbf{B}_{156} & = & \left(\frac{1}{2}+y_{10}\right)\mathbf{a}_1+z_{10}\,\mathbf{a}_2+\left(\frac{1}{2}-x_{10}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+y_{10}\right)a\,\hat{\mathbf{x}}+z_{10}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-x_{10}\right)a\,\hat{\mathbf{z}} & (24d) & \text{H V}\\
\mathbf{B}_{157} & = & x_{11}\,\mathbf{a}_1+y_{11}\,\mathbf{a}_2+z_{11}\,\mathbf{a}_3 & = & x_{11}a\,\hat{\mathbf{x}}+y_{11}a\,\hat{\mathbf{y}}+z_{11}a\,\hat{\mathbf{z}} & (24d) & \text{H VI}\\
\mathbf{B}_{158} & = & \left(\frac{1}{2}-x_{11}\right)\mathbf{a}_1-y_{11}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{11}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{11}\right)a\,\hat{\mathbf{x}}-y_{11}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{11}\right)a\,\hat{\mathbf{z}} & (24d) & \text{H VI}
\end{array}
$$

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\begin{array}{lcccll}
\mathbf{B}_{159} & = & -x_{11}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{11}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{11}\right)\mathbf{a}_3 & = & -x_{11}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{11}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{11}\right)a\,\hat{\mathbf{z}} & (24d)\ \text{H VI}\\
\mathbf{B}_{160} & = & \left(\frac{1}{2}+x_{11}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{11}\right)\mathbf{a}_2-z_{11}\,\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{11}\right)a\,\hat{\mathbf{y}}-z_{11}a\,\hat{\mathbf{z}} & (24d)\ \text{H VI}\\
\mathbf{B}_{161} & = & z_{11}\,\mathbf{a}_1+x_{11}\,\mathbf{a}_2+y_{11}\,\mathbf{a}_3 & = & z_{11}a\,\hat{\mathbf{x}}+x_{11}a\,\hat{\mathbf{y}}+y_{11}a\,\hat{\mathbf{z}} & (24d)\ \text{H VI}\\
\mathbf{B}_{162} & = & \left(\frac{1}{2}+z_{11}\right)\mathbf{a}_1+\left(\frac{1}{2}-x_{11}\right)\mathbf{a}_2-y_{11}\,\mathbf{a}_3 & = & \left(\frac{1}{2}+z_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_{11}\right)a\,\hat{\mathbf{y}}-y_{11}a\,\hat{\mathbf{z}} & (24d)\ \text{H VI}\\
\mathbf{B}_{163} & = & \left(\frac{1}{2}-z_{11}\right)\mathbf{a}_1-x_{11}\,\mathbf{a}_2+\left(\frac{1}{2}+y_{11}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-z_{11}\right)a\,\hat{\mathbf{x}}-x_{11}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+y_{11}\right)a\,\hat{\mathbf{z}} & (24d)\ \text{H VI}\\
\mathbf{B}_{164} & = & -z_{11}\,\mathbf{a}_1+\left(\frac{1}{2}+x_{11}\right)\mathbf{a}_2+\left(\frac{1}{2}-y_{11}\right)\mathbf{a}_3 & = & -z_{11}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_{11}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-y_{11}\right)a\,\hat{\mathbf{z}} & (24d)\ \text{H VI}\\
\mathbf{B}_{165} & = & y_{11}\,\mathbf{a}_1+z_{11}\,\mathbf{a}_2+x_{11}\,\mathbf{a}_3 & = & y_{11}a\,\hat{\mathbf{x}}+z_{11}a\,\hat{\mathbf{y}}+x_{11}a\,\hat{\mathbf{z}} & (24d)\ \text{H VI}\\
\mathbf{B}_{166} & = & -y_{11}\,\mathbf{a}_1+\left(\frac{1}{2}+z_{11}\right)\mathbf{a}_2+\left(\frac{1}{2}-x_{11}\right)\mathbf{a}_3 & = & -y_{11}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+z_{11}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-x_{11}\right)a\,\hat{\mathbf{z}} & (24d)\ \text{H VI}\\
\mathbf{B}_{167} & = & \left(\frac{1}{2}+y_{11}\right)\mathbf{a}_1+\left(\frac{1}{2}-z_{11}\right)\mathbf{a}_2-x_{11}\,\mathbf{a}_3 & = & \left(\frac{1}{2}+y_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-z_{11}\right)a\,\hat{\mathbf{y}}-x_{11}a\,\hat{\mathbf{z}} & (24d)\ \text{H VI}\\
\mathbf{B}_{168} & = & \left(\frac{1}{2}-y_{11}\right)\mathbf{a}_1-z_{11}\,\mathbf{a}_2+\left(\frac{1}{2}+x_{11}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-y_{11}\right)a\,\hat{\mathbf{x}}-z_{11}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+x_{11}\right)a\,\hat{\mathbf{z}} & (24d)\ \text{H VI}\\
\mathbf{B}_{169} & = & -x_{11}\,\mathbf{a}_1-y_{11}\,\mathbf{a}_2-z_{11}\,\mathbf{a}_3 & = & -x_{11}a\,\hat{\mathbf{x}}-y_{11}a\,\hat{\mathbf{y}}-z_{11}a\,\hat{\mathbf{z}} & (24d)\ \text{H VI}\\
\mathbf{B}_{170} & = & \left(\frac{1}{2}+x_{11}\right)\mathbf{a}_1+y_{11}\,\mathbf{a}_2+\left(\frac{1}{2}-z_{11}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{11}\right)a\,\hat{\mathbf{x}}+y_{11}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{11}\right)a\,\hat{\mathbf{z}} & (24d)\ \text{H VI}\\
\mathbf{B}_{171} & = & x_{11}\,\mathbf{a}_1+\left(\frac{1}{2}-y_{11}\right)\mathbf{a}_2+\left(\frac{1}{2}+z_{11}\right)\mathbf{a}_3 & = & x_{11}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{11}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{11}\right)a\,\hat{\mathbf{z}} & (24d)\ \text{H VI}\\
\mathbf{B}_{172} & = & \left(\frac{1}{2}-x_{11}\right)\mathbf{a}_1+\left(\frac{1}{2}+y_{11}\right)\mathbf{a}_2+z_{11}\,\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{11}\right)a\,\hat{\mathbf{y}}+z_{11}a\,\hat{\mathbf{z}} & (24d)\ \text{H VI}\\
\mathbf{B}_{173} & = & -z_{11}\,\mathbf{a}_1-x_{11}\,\mathbf{a}_2-y_{11}\,\mathbf{a}_3 & = & -z_{11}a\,\hat{\mathbf{x}}-x_{11}a\,\hat{\mathbf{y}}-y_{11}a\,\hat{\mathbf{z}} & (24d)\ \text{H VI}\\
\mathbf{B}_{174} & = & \left(\frac{1}{2}-z_{11}\right)\mathbf{a}_1+\left(\frac{1}{2}+x_{11}\right)\mathbf{a}_2+y_{11}\,\mathbf{a}_3 & = & \left(\frac{1}{2}-z_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_{11}\right)a\,\hat{\mathbf{y}}+y_{11}a\,\hat{\mathbf{z}} & (24d)\ \text{H VI}\\
\mathbf{B}_{175} & = & \left(\frac{1}{2}+z_{11}\right)\mathbf{a}_1+x_{11}\,\mathbf{a}_2+\left(\frac{1}{2}-y_{11}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+z_{11}\right)a\,\hat{\mathbf{x}}+x_{11}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-y_{11}\right)a\,\hat{\mathbf{z}} & (24d)\ \text{H VI}\\
\mathbf{B}_{176} & = & z_{11}\,\mathbf{a}_1+\left(\frac{1}{2}-x_{11}\right)\mathbf{a}_2+\left(\frac{1}{2}+y_{11}\right)\mathbf{a}_3 & = & z_{11}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_{11}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+y_{11}\right)a\,\hat{\mathbf{z}} & (24d)\ \text{H VI}\\
\mathbf{B}_{177} & = & -y_{11}\,\mathbf{a}_1-z_{11}\,\mathbf{a}_2-x_{11}\,\mathbf{a}_3 & = & -y_{11}a\,\hat{\mathbf{x}}-z_{11}a\,\hat{\mathbf{y}}-x_{11}a\,\hat{\mathbf{z}} & (24d)\ \text{H VI}\\
\mathbf{B}_{178} & = & y_{11}\,\mathbf{a}_1+\left(\frac{1}{2}-z_{11}\right)\mathbf{a}_2+\left(\frac{1}{2}+x_{11}\right)\mathbf{a}_3 & = & y_{11}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-z_{11}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+x_{11}\right)a\,\hat{\mathbf{z}} & (24d)\ \text{H VI}\\
\mathbf{B}_{179} & = & \left(\frac{1}{2}-y_{11}\right)\mathbf{a}_1+\left(\frac{1}{2}+z_{11}\right)\mathbf{a}_2+x_{11}\,\mathbf{a}_3 & = & \left(\frac{1}{2}-y_{11}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+z_{11}\right)a\,\hat{\mathbf{y}}+x_{11}a\,\hat{\mathbf{z}} & (24d)\ \text{H VI}\\
\mathbf{B}_{180} & = & \left(\frac{1}{2}+y_{11}\right)\mathbf{a}_1+z_{11}\,\mathbf{a}_2+\left(\frac{1}{2}-x_{11}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+y_{11}\right)a\,\hat{\mathbf{x}}+z_{11}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-x_{11}\right)a\,\hat{\mathbf{z}} & (24d)\ \text{H VI}\\
\mathbf{B}_{181} & = & x_{12}\,\mathbf{a}_1+y_{12}\,\mathbf{a}_2+z_{12}\,\mathbf{a}_3 & = & x_{12}a\,\hat{\mathbf{x}}+y_{12}a\,\hat{\mathbf{y}}+z_{12}a\,\hat{\mathbf{z}} & (24d)\ \text{O II}\\
\mathbf{B}_{182} & = & \left(\frac{1}{2}-x_{12}\right)\mathbf{a}_1-y_{12}\,\mathbf{a}_2+\left(\frac{1}{2}+z_{12}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_{12}\right)a\,\hat{\mathbf{x}}-y_{12}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+z_{12}\right)a\,\hat{\mathbf{z}} & (24d)\ \text{O II}\\
\mathbf{B}_{183} & = & -x_{12}\,\mathbf{a}_1+\left(\frac{1}{2}+y_{12}\right)\mathbf{a}_2+\left(\frac{1}{2}-z_{12}\right)\mathbf{a}_3 & = & -x_{12}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_{12}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_{12}\right)a\,\hat{\mathbf{z}} & (24d)\ \text{O II}\\
\mathbf{B}_{184} & = & \left(\frac{1}{2}+x_{12}\right)\mathbf{a}_1+\left(\frac{1}{2}-y_{12}\right)\mathbf{a}_2-z_{12}\,\mathbf{a}_3 & = & \left(\frac{1}{2}+x_{12}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_{12}\right)a\,\hat{\mathbf{y}}-z_{12}a\,\hat{\mathbf{z}} & (24d)\ \text{O II}\\
\mathbf{B}_{185} & = & z_{12}\,\mathbf{a}_1+x_{12}\,\mathbf{a}_2+y_{12}\,\mathbf{a}_3 & = & z_{12}a\,\hat{\mathbf{x}}+x_{12}a\,\hat{\mathbf{y}}+y_{12}a\,\hat{\mathbf{z}} & (24d)\ \text{O II}\\
\mathbf{B}_{186} & = & \left(\frac{1}{2}+z_{12}\right)\mathbf{a}_1+\left(\frac{1}{2}-x_{12}\right)\mathbf{a}_2-y_{12}\,\mathbf{a}_3 & = & \left(\frac{1}{2}+z_{12}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_{12}\right)a\,\hat{\mathbf{y}}-y_{12}a\,\hat{\mathbf{z}} & (24d)\ \text{O II}\\
\mathbf{B}_{187} & = & \left(\frac{1}{2}-z_{12}\right)\mathbf{a}_1-x_{12}\,\mathbf{a}_2+\left(\frac{1}{2}+y_{12}\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-z_{12}\right)a\,\hat{\mathbf{x}}-x_{12}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+y_{12}\right)a\,\hat{\mathbf{z}} & (24d)\ \text{O II}\\
\mathbf{B}_{188} & = & -z_{12}\,\mathbf{a}_1+\left(\frac{1}{2}+x_{12}\right)\mathbf{a}_2+\left(\frac{1}{2}-y_{12}\right)\mathbf{a}_3 & = & -z_{12}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_{12}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-y_{12}\right)a\,\hat{\mathbf{z}} & (24d)\ \text{O II}\\
\mathbf{B}_{189} & = & y_{12}\,\mathbf{a}_1+z_{12}\,\mathbf{a}_2+x_{12}\,\mathbf{a}_3 & = & y_{12}a\,\hat{\mathbf{x}}+z_{12}a\,\hat{\mathbf{y}}+x_{12}a\,\hat{\mathbf{z}} & (24d)\ \text{O II}\\
\mathbf{B}_{190} & = & -y_{12}\,\mathbf{a}_1+\left(\frac{1}{2}+z_{12}\right)\mathbf{a}_2+\left(\frac{1}{2}-x_{12}\right)\mathbf{a}_3 & = & -y_{12}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+z_{12}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-x_{12}\right)a\,\hat{\mathbf{z}} & (24d)\ \text{O II}
\end{array}
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$$\mathbf{B}_{191} = \left(\tfrac{1}{2}+y_{12}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-z_{12}\right)\mathbf{a}_2-x_{12}\,\mathbf{a}_3 = \left(\tfrac{1}{2}+y_{12}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-z_{12}\right)a\,\hat{\mathbf{y}}-x_{12}a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O II}$$

$$\mathbf{B}_{192} = \left(\tfrac{1}{2}-y_{12}\right)\mathbf{a}_1-z_{12}\,\mathbf{a}_2+\left(\tfrac{1}{2}+x_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-y_{12}\right)a\,\hat{\mathbf{x}}-z_{12}a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+x_{12}\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O II}$$

$$\mathbf{B}_{193} = -x_{12}\,\mathbf{a}_1-y_{12}\,\mathbf{a}_2-z_{12}\,\mathbf{a}_3 = -x_{12}a\,\hat{\mathbf{x}}-y_{12}a\,\hat{\mathbf{y}}-z_{12}a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O II}$$

$$\mathbf{B}_{194} = \left(\tfrac{1}{2}+x_{12}\right)\mathbf{a}_1+y_{12}\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{12}\right)a\,\hat{\mathbf{x}}+y_{12}a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{12}\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O II}$$

$$\mathbf{B}_{195} = x_{12}\,\mathbf{a}_1+\left(\tfrac{1}{2}-y_{12}\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_{12}\right)\mathbf{a}_3 = x_{12}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{12}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{12}\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O II}$$

$$\mathbf{B}_{196} = \left(\tfrac{1}{2}-x_{12}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{12}\right)\mathbf{a}_2+z_{12}\,\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{12}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{12}\right)a\,\hat{\mathbf{y}}+z_{12}a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O II}$$

$$\mathbf{B}_{197} = -z_{12}\,\mathbf{a}_1-x_{12}\,\mathbf{a}_2-y_{12}\,\mathbf{a}_3 = -z_{12}a\,\hat{\mathbf{x}}-x_{12}a\,\hat{\mathbf{y}}-y_{12}a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O II}$$

$$\mathbf{B}_{198} = \left(\tfrac{1}{2}-z_{12}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_{12}\right)\mathbf{a}_2+y_{12}\,\mathbf{a}_3 = \left(\tfrac{1}{2}-z_{12}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_{12}\right)a\,\hat{\mathbf{y}}+y_{12}a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O II}$$

$$\mathbf{B}_{199} = \left(\tfrac{1}{2}+z_{12}\right)\mathbf{a}_1+x_{12}\,\mathbf{a}_2+\left(\tfrac{1}{2}-y_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+z_{12}\right)a\,\hat{\mathbf{x}}+x_{12}a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-y_{12}\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O II}$$

$$\mathbf{B}_{200} = z_{12}\,\mathbf{a}_1+\left(\tfrac{1}{2}-x_{12}\right)\mathbf{a}_2+\left(\tfrac{1}{2}+y_{12}\right)\mathbf{a}_3 = z_{12}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_{12}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+y_{12}\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O II}$$

$$\mathbf{B}_{201} = -y_{12}\,\mathbf{a}_1-z_{12}\,\mathbf{a}_2-x_{12}\,\mathbf{a}_3 = -y_{12}a\,\hat{\mathbf{x}}-z_{12}a\,\hat{\mathbf{y}}-x_{12}a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O II}$$

$$\mathbf{B}_{202} = y_{12}\,\mathbf{a}_1+\left(\tfrac{1}{2}-z_{12}\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_{12}\right)\mathbf{a}_3 = y_{12}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-z_{12}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+x_{12}\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O II}$$

$$\mathbf{B}_{203} = \left(\tfrac{1}{2}-y_{12}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+z_{12}\right)\mathbf{a}_2+x_{12}\,\mathbf{a}_3 = \left(\tfrac{1}{2}-y_{12}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+z_{12}\right)a\,\hat{\mathbf{y}}+x_{12}a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O II}$$

$$\mathbf{B}_{204} = \left(\tfrac{1}{2}+y_{12}\right)\mathbf{a}_1+z_{12}\,\mathbf{a}_2+\left(\tfrac{1}{2}-x_{12}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+y_{12}\right)a\,\hat{\mathbf{x}}+z_{12}a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-x_{12}\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O II}$$

$$\mathbf{B}_{205} = x_{13}\,\mathbf{a}_1+y_{13}\,\mathbf{a}_2+z_{13}\,\mathbf{a}_3 = x_{13}a\,\hat{\mathbf{x}}+y_{13}a\,\hat{\mathbf{y}}+z_{13}a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O III}$$

$$\mathbf{B}_{206} = \left(\tfrac{1}{2}-x_{13}\right)\mathbf{a}_1-y_{13}\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{13}\right)a\,\hat{\mathbf{x}}-y_{13}a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{13}\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O III}$$

$$\mathbf{B}_{207} = -x_{13}\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_{13}\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_{13}\right)\mathbf{a}_3 = -x_{13}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{13}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{13}\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O III}$$

$$\mathbf{B}_{208} = \left(\tfrac{1}{2}+x_{13}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_{13}\right)\mathbf{a}_2-z_{13}\,\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{13}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{13}\right)a\,\hat{\mathbf{y}}-z_{13}a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O III}$$

$$\mathbf{B}_{209} = z_{13}\,\mathbf{a}_1+x_{13}\,\mathbf{a}_2+y_{13}\,\mathbf{a}_3 = z_{13}a\,\hat{\mathbf{x}}+x_{13}a\,\hat{\mathbf{y}}+y_{13}a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O III}$$

$$\mathbf{B}_{210} = \left(\tfrac{1}{2}+z_{13}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-x_{13}\right)\mathbf{a}_2-y_{13}\,\mathbf{a}_3 = \left(\tfrac{1}{2}+z_{13}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_{13}\right)a\,\hat{\mathbf{y}}-y_{13}a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O III}$$

$$\mathbf{B}_{211} = \left(\tfrac{1}{2}-z_{13}\right)\mathbf{a}_1-x_{13}\,\mathbf{a}_2+\left(\tfrac{1}{2}+y_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-z_{13}\right)a\,\hat{\mathbf{x}}-x_{13}a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+y_{13}\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O III}$$

$$\mathbf{B}_{212} = -z_{13}\,\mathbf{a}_1+\left(\tfrac{1}{2}+x_{13}\right)\mathbf{a}_2+\left(\tfrac{1}{2}-y_{13}\right)\mathbf{a}_3 = -z_{13}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_{13}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-y_{13}\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O III}$$

$$\mathbf{B}_{213} = y_{13}\,\mathbf{a}_1+z_{13}\,\mathbf{a}_2+x_{13}\,\mathbf{a}_3 = y_{13}a\,\hat{\mathbf{x}}+z_{13}a\,\hat{\mathbf{y}}+x_{13}a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O III}$$

$$\mathbf{B}_{214} = -y_{13}\,\mathbf{a}_1+\left(\tfrac{1}{2}+z_{13}\right)\mathbf{a}_2+\left(\tfrac{1}{2}-x_{13}\right)\mathbf{a}_3 = -y_{13}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+z_{13}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-x_{13}\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O III}$$

$$\mathbf{B}_{215} = \left(\tfrac{1}{2}+y_{13}\right)\mathbf{a}_1+\left(\tfrac{1}{2}-z_{13}\right)\mathbf{a}_2-x_{13}\,\mathbf{a}_3 = \left(\tfrac{1}{2}+y_{13}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-z_{13}\right)a\,\hat{\mathbf{y}}-x_{13}a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O III}$$

$$\mathbf{B}_{216} = \left(\tfrac{1}{2}-y_{13}\right)\mathbf{a}_1-z_{13}\,\mathbf{a}_2+\left(\tfrac{1}{2}+x_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-y_{13}\right)a\,\hat{\mathbf{x}}-z_{13}a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+x_{13}\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O III}$$

$$\mathbf{B}_{217} = -x_{13}\,\mathbf{a}_1-y_{13}\,\mathbf{a}_2-z_{13}\,\mathbf{a}_3 = -x_{13}a\,\hat{\mathbf{x}}-y_{13}a\,\hat{\mathbf{y}}-z_{13}a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O III}$$

$$\mathbf{B}_{218} = \left(\tfrac{1}{2}+x_{13}\right)\mathbf{a}_1+y_{13}\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_{13}\right)a\,\hat{\mathbf{x}}+y_{13}a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_{13}\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O III}$$

$$\mathbf{B}_{219} = x_{13}\,\mathbf{a}_1+\left(\tfrac{1}{2}-y_{13}\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_{13}\right)\mathbf{a}_3 = x_{13}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_{13}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_{13}\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O III}$$

$$\mathbf{B}_{220} = \left(\tfrac{1}{2}-x_{13}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_{13}\right)\mathbf{a}_2+z_{13}\,\mathbf{a}_3 = \left(\tfrac{1}{2}-x_{13}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_{13}\right)a\,\hat{\mathbf{y}}+z_{13}a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O III}$$

$$\mathbf{B}_{221} = -z_{13}\,\mathbf{a}_1-x_{13}\,\mathbf{a}_2-y_{13}\,\mathbf{a}_3 = -z_{13}a\,\hat{\mathbf{x}}-x_{13}a\,\hat{\mathbf{y}}-y_{13}a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O III}$$

$$\mathbf{B}_{222} = \left(\tfrac{1}{2}-z_{13}\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_{13}\right)\mathbf{a}_2+y_{13}\,\mathbf{a}_3 = \left(\tfrac{1}{2}-z_{13}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_{13}\right)a\,\hat{\mathbf{y}}+y_{13}a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O III}$$

$$\mathbf{B}_{223} = \left(\tfrac{1}{2}+z_{13}\right)\mathbf{a}_1+x_{13}\,\mathbf{a}_2+\left(\tfrac{1}{2}-y_{13}\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+z_{13}\right)a\,\hat{\mathbf{x}}+x_{13}a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-y_{13}\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O III}$$

$$\mathbf{B}_{224} = z_{13}\,\mathbf{a}_1+\left(\tfrac{1}{2}-x_{13}\right)\mathbf{a}_2+\left(\tfrac{1}{2}+y_{13}\right)\mathbf{a}_3 = z_{13}a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_{13}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+y_{13}\right)a\,\hat{\mathbf{z}} \quad (24d) \quad \text{O III}$$

$\mathbf{B}_{225}$	=	$-y_{13}\,\mathbf{a}_1 - z_{13}\,\mathbf{a}_2 - x_{13}\,\mathbf{a}_3$	=	$-y_{13}a\,\hat{\mathbf{x}} - z_{13}a\,\hat{\mathbf{y}} - x_{13}a\,\hat{\mathbf{z}}$	(24*d*)	O III
$\mathbf{B}_{226}$	=	$y_{13}\,\mathbf{a}_1 + \left(\frac{1}{2} - z_{13}\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_{13}\right)\mathbf{a}_3$	=	$y_{13}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - z_{13}\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_{13}\right)a\,\hat{\mathbf{z}}$	(24*d*)	O III
$\mathbf{B}_{227}$	=	$\left(\frac{1}{2} - y_{13}\right)\mathbf{a}_1 + \left(\frac{1}{2} + z_{13}\right)\mathbf{a}_2 + x_{13}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_{13}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + z_{13}\right)a\,\hat{\mathbf{y}} + x_{13}a\,\hat{\mathbf{z}}$	(24*d*)	O III
$\mathbf{B}_{228}$	=	$\left(\frac{1}{2} + y_{13}\right)\mathbf{a}_1 + z_{13}\,\mathbf{a}_2 + \left(\frac{1}{2} - x_{13}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_{13}\right)a\,\hat{\mathbf{x}} + z_{13}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_{13}\right)a\,\hat{\mathbf{z}}$	(24*d*)	O III
$\mathbf{B}_{229}$	=	$x_{14}\,\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3$	=	$x_{14}a\,\hat{\mathbf{x}} + y_{14}a\,\hat{\mathbf{y}} + z_{14}a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{230}$	=	$\left(\frac{1}{2} - x_{14}\right)\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 + \left(\frac{1}{2} + z_{14}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{14}\right)a\,\hat{\mathbf{x}} - y_{14}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{14}\right)a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{231}$	=	$-x_{14}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_{14}\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_{14}\right)\mathbf{a}_3$	=	$-x_{14}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{14}\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{14}\right)a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{232}$	=	$\left(\frac{1}{2} + x_{14}\right)\mathbf{a}_1 + \left(\frac{1}{2} - y_{14}\right)\mathbf{a}_2 - z_{14}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{14}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{14}\right)a\,\hat{\mathbf{y}} - z_{14}a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{233}$	=	$z_{14}\,\mathbf{a}_1 + x_{14}\,\mathbf{a}_2 + y_{14}\,\mathbf{a}_3$	=	$z_{14}a\,\hat{\mathbf{x}} + x_{14}a\,\hat{\mathbf{y}} + y_{14}a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{234}$	=	$\left(\frac{1}{2} + z_{14}\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_{14}\right)\mathbf{a}_2 - y_{14}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + z_{14}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_{14}\right)a\,\hat{\mathbf{y}} - y_{14}a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{235}$	=	$\left(\frac{1}{2} - z_{14}\right)\mathbf{a}_1 - x_{14}\,\mathbf{a}_2 + \left(\frac{1}{2} + y_{14}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - z_{14}\right)a\,\hat{\mathbf{x}} - x_{14}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + y_{14}\right)a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{236}$	=	$-z_{14}\,\mathbf{a}_1 + \left(\frac{1}{2} + x_{14}\right)\mathbf{a}_2 + \left(\frac{1}{2} - y_{14}\right)\mathbf{a}_3$	=	$-z_{14}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_{14}\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - y_{14}\right)a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{237}$	=	$y_{14}\,\mathbf{a}_1 + z_{14}\,\mathbf{a}_2 + x_{14}\,\mathbf{a}_3$	=	$y_{14}a\,\hat{\mathbf{x}} + z_{14}a\,\hat{\mathbf{y}} + x_{14}a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{238}$	=	$-y_{14}\,\mathbf{a}_1 + \left(\frac{1}{2} + z_{14}\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_{14}\right)\mathbf{a}_3$	=	$-y_{14}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + z_{14}\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_{14}\right)a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{239}$	=	$\left(\frac{1}{2} + y_{14}\right)\mathbf{a}_1 + \left(\frac{1}{2} - z_{14}\right)\mathbf{a}_2 - x_{14}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_{14}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - z_{14}\right)a\,\hat{\mathbf{y}} - x_{14}a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{240}$	=	$\left(\frac{1}{2} - y_{14}\right)\mathbf{a}_1 - z_{14}\,\mathbf{a}_2 + \left(\frac{1}{2} + x_{14}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_{14}\right)a\,\hat{\mathbf{x}} - z_{14}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_{14}\right)a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{241}$	=	$-x_{14}\,\mathbf{a}_1 - y_{14}\,\mathbf{a}_2 - z_{14}\,\mathbf{a}_3$	=	$-x_{14}a\,\hat{\mathbf{x}} - y_{14}a\,\hat{\mathbf{y}} - z_{14}a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{242}$	=	$\left(\frac{1}{2} + x_{14}\right)\mathbf{a}_1 + y_{14}\,\mathbf{a}_2 + \left(\frac{1}{2} - z_{14}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_{14}\right)a\,\hat{\mathbf{x}} + y_{14}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_{14}\right)a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{243}$	=	$x_{14}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_{14}\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_{14}\right)\mathbf{a}_3$	=	$x_{14}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_{14}\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_{14}\right)a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{244}$	=	$\left(\frac{1}{2} - x_{14}\right)\mathbf{a}_1 + \left(\frac{1}{2} + y_{14}\right)\mathbf{a}_2 + z_{14}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{14}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_{14}\right)a\,\hat{\mathbf{y}} + z_{14}a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{245}$	=	$-z_{14}\,\mathbf{a}_1 - x_{14}\,\mathbf{a}_2 - y_{14}\,\mathbf{a}_3$	=	$-z_{14}a\,\hat{\mathbf{x}} - x_{14}a\,\hat{\mathbf{y}} - y_{14}a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{246}$	=	$\left(\frac{1}{2} - z_{14}\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_{14}\right)\mathbf{a}_2 + y_{14}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - z_{14}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_{14}\right)a\,\hat{\mathbf{y}} + y_{14}a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{247}$	=	$\left(\frac{1}{2} + z_{14}\right)\mathbf{a}_1 + x_{14}\,\mathbf{a}_2 + \left(\frac{1}{2} - y_{14}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + z_{14}\right)a\,\hat{\mathbf{x}} + x_{14}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - y_{14}\right)a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{248}$	=	$z_{14}\,\mathbf{a}_1 + \left(\frac{1}{2} - x_{14}\right)\mathbf{a}_2 + \left(\frac{1}{2} + y_{14}\right)\mathbf{a}_3$	=	$z_{14}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_{14}\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + y_{14}\right)a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{249}$	=	$-y_{14}\,\mathbf{a}_1 - z_{14}\,\mathbf{a}_2 - x_{14}\,\mathbf{a}_3$	=	$-y_{14}a\,\hat{\mathbf{x}} - z_{14}a\,\hat{\mathbf{y}} - x_{14}a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{250}$	=	$y_{14}\,\mathbf{a}_1 + \left(\frac{1}{2} - z_{14}\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_{14}\right)\mathbf{a}_3$	=	$y_{14}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - z_{14}\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_{14}\right)a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{251}$	=	$\left(\frac{1}{2} - y_{14}\right)\mathbf{a}_1 + \left(\frac{1}{2} + z_{14}\right)\mathbf{a}_2 + x_{14}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_{14}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + z_{14}\right)a\,\hat{\mathbf{y}} + x_{14}a\,\hat{\mathbf{z}}$	(24*d*)	O IV
$\mathbf{B}_{252}$	=	$\left(\frac{1}{2} + y_{14}\right)\mathbf{a}_1 + z_{14}\,\mathbf{a}_2 + \left(\frac{1}{2} - x_{14}\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_{14}\right)a\,\hat{\mathbf{x}} + z_{14}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_{14}\right)a\,\hat{\mathbf{z}}$	(24*d*)	O IV

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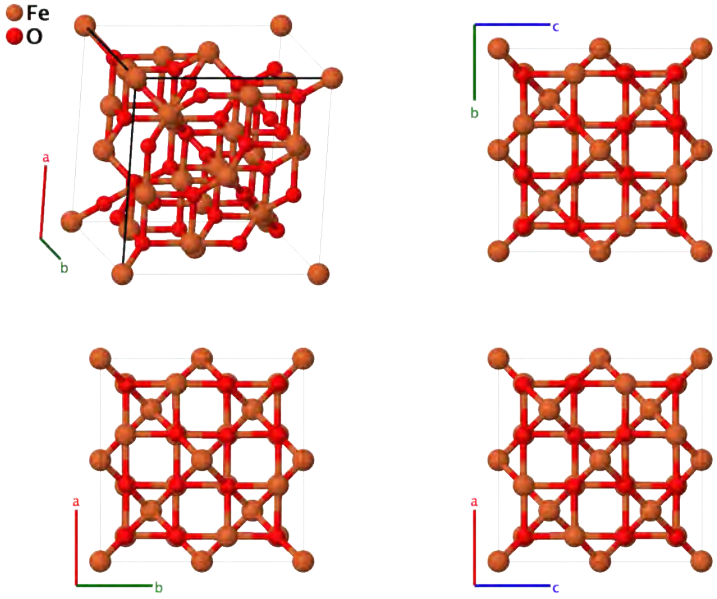
Geometry files:
- CIF: pp. 1787
- POSCAR: pp. 1787

Maghemite (γ-Fe_2O_3, $D5_7$) Structure: A2B3_cP60_212_bcd_ace

http://aflow.org/prototype-encyclopedia/A2B3_cP60_212_bcd_ace



Prototype	:	Fe_2O_3
AFLOW prototype label	:	A2B3_cP60_212_bcd_ace
***Strukturbericht* designation**	:	$D5_7$
Pearson symbol	:	cP60
Space group number	:	212
Space group symbol	:	$P4_332$
AFLOW prototype command	:	`aflow --proto=A2B3_cP60_212_bcd_ace --params=`$a, x_3, x_4, y_5, x_6, y_6, z_6$

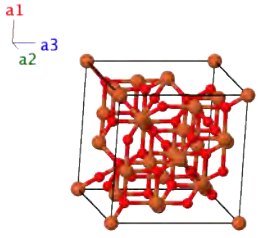
Other compounds with this structure

- γ-Al_2O_3 (γ-corundum)

- (Hermann, 1937) gives γ-Al_2O_3 as the prototype for *Strukturbericht* $D5_7$, but states that the data for γ-Fe_2O_3 is more reliable and presents the data for the latter compound, which we use as the prototype. This is a rock-salt ($B1$) structure with defects. This structure can also be expressed in the enantiomorphic space group $P4_132$ #213.

Simple Cubic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = a\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{8}\,\mathbf{a}_1 + \frac{1}{8}\,\mathbf{a}_2 + \frac{1}{8}\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_2$	=	$\frac{3}{8}\,\mathbf{a}_1 + \frac{7}{8}\,\mathbf{a}_2 + \frac{5}{8}\,\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} + \frac{7}{8}a\,\hat{\mathbf{y}} + \frac{5}{8}a\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_3$	=	$\frac{7}{8}\,\mathbf{a}_1 + \frac{5}{8}\,\mathbf{a}_2 + \frac{3}{8}\,\mathbf{a}_3$	=	$\frac{7}{8}a\,\hat{\mathbf{x}} + \frac{5}{8}a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_4$	=	$\frac{5}{8}\,\mathbf{a}_1 + \frac{3}{8}\,\mathbf{a}_2 + \frac{7}{8}\,\mathbf{a}_3$	=	$\frac{5}{8}a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} + \frac{7}{8}a\,\hat{\mathbf{z}}$	(4*a*)	O I
$\mathbf{B}_5$	=	$\frac{5}{8}\,\mathbf{a}_1 + \frac{5}{8}\,\mathbf{a}_2 + \frac{5}{8}\,\mathbf{a}_3$	=	$\frac{5}{8}a\,\hat{\mathbf{x}} + \frac{5}{8}a\,\hat{\mathbf{y}} + \frac{5}{8}a\,\hat{\mathbf{z}}$	(4*b*)	Fe I
$\mathbf{B}_6$	=	$\frac{7}{8}\,\mathbf{a}_1 + \frac{3}{8}\,\mathbf{a}_2 + \frac{1}{8}\,\mathbf{a}_3$	=	$\frac{7}{8}a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(4*b*)	Fe I
$\mathbf{B}_7$	=	$\frac{3}{8}\,\mathbf{a}_1 + \frac{1}{8}\,\mathbf{a}_2 + \frac{7}{8}\,\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \frac{7}{8}a\,\hat{\mathbf{z}}$	(4*b*)	Fe I
$\mathbf{B}_8$	=	$\frac{1}{8}\,\mathbf{a}_1 + \frac{7}{8}\,\mathbf{a}_2 + \frac{3}{8}\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \frac{7}{8}a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}}$	(4*b*)	Fe I
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(8*c*)	Fe II
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{z}}$	(8*c*)	Fe II
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{z}}$	(8*c*)	Fe II
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(8*c*)	Fe II
$\mathbf{B}_{13}$	=	$\left(\frac{1}{4} + x_3\right)\mathbf{a}_1 + \left(\frac{3}{4} + x_3\right)\mathbf{a}_2 + \left(\frac{3}{4} - x_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{3}{4} + x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{3}{4} - x_3\right)a\,\hat{\mathbf{z}}$	(8*c*)	Fe II
$\mathbf{B}_{14}$	=	$\left(\frac{1}{4} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{4} - x_3\right)\mathbf{a}_2 + \left(\frac{1}{4} - x_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{z}}$	(8*c*)	Fe II
$\mathbf{B}_{15}$	=	$\left(\frac{3}{4} + x_3\right)\mathbf{a}_1 + \left(\frac{3}{4} - x_3\right)\mathbf{a}_2 + \left(\frac{1}{4} + x_3\right)\mathbf{a}_3$	=	$\left(\frac{3}{4} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{3}{4} - x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{z}}$	(8*c*)	Fe II
$\mathbf{B}_{16}$	=	$\left(\frac{3}{4} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{4} + x_3\right)\mathbf{a}_2 + \left(\frac{3}{4} + x_3\right)\mathbf{a}_3$	=	$\left(\frac{3}{4} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{3}{4} + x_3\right)a\,\hat{\mathbf{z}}$	(8*c*)	Fe II
$\mathbf{B}_{17}$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	(8*c*)	O II
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \left(\frac{1}{2} + x_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{z}}$	(8*c*)	O II
$\mathbf{B}_{19}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + x_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{z}}$	(8*c*)	O II
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{y}} - x_4a\,\hat{\mathbf{z}}$	(8*c*)	O II
$\mathbf{B}_{21}$	=	$\left(\frac{1}{4} + x_4\right)\mathbf{a}_1 + \left(\frac{3}{4} + x_4\right)\mathbf{a}_2 + \left(\frac{3}{4} - x_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{3}{4} + x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{3}{4} - x_4\right)a\,\hat{\mathbf{z}}$	(8*c*)	O II
$\mathbf{B}_{22}$	=	$\left(\frac{1}{4} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{4} - x_4\right)\mathbf{a}_2 + \left(\frac{1}{4} - x_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{z}}$	(8*c*)	O II
$\mathbf{B}_{23}$	=	$\left(\frac{3}{4} + x_4\right)\mathbf{a}_1 + \left(\frac{3}{4} - x_4\right)\mathbf{a}_2 + \left(\frac{1}{4} + x_4\right)\mathbf{a}_3$	=	$\left(\frac{3}{4} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{3}{4} - x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + x_4\right)a\,\hat{\mathbf{z}}$	(8*c*)	O II
$\mathbf{B}_{24}$	=	$\left(\frac{3}{4} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{4} + x_4\right)\mathbf{a}_2 + \left(\frac{3}{4} + x_4\right)\mathbf{a}_3$	=	$\left(\frac{3}{4} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{3}{4} + x_4\right)a\,\hat{\mathbf{z}}$	(8*c*)	O II
$\mathbf{B}_{25}$	=	$\frac{1}{8}\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + \left(\frac{1}{4} - y_5\right)\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + y_5a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - y_5\right)a\,\hat{\mathbf{z}}$	(12*d*)	Fe III
$\mathbf{B}_{26}$	=	$\frac{3}{8}\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + \left(\frac{3}{4} - y_5\right)\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} - y_5a\,\hat{\mathbf{y}} + \left(\frac{3}{4} - y_5\right)a\,\hat{\mathbf{z}}$	(12*d*)	Fe III
$\mathbf{B}_{27}$	=	$\frac{7}{8}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_5\right)\mathbf{a}_2 + \left(\frac{1}{4} + y_5\right)\mathbf{a}_3$	=	$\frac{7}{8}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + y_5\right)a\,\hat{\mathbf{z}}$	(12*d*)	Fe III

$$\mathbf{B}_{28} = \tfrac{5}{8}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_5\right)\mathbf{a}_2 + \left(\tfrac{3}{4} + y_5\right)\mathbf{a}_3 = \tfrac{5}{8}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{3}{4} + y_5\right)a\,\hat{\mathbf{z}} \quad (12d) \quad \text{Fe III}$$

$$\mathbf{B}_{29} = \left(\tfrac{1}{4} - y_5\right)\mathbf{a}_1 + \tfrac{1}{8}\,\mathbf{a}_2 + y_5\,\mathbf{a}_3 = \left(\tfrac{1}{4} - y_5\right)a\,\hat{\mathbf{x}} + \tfrac{1}{8}a\,\hat{\mathbf{y}} + y_5 a\,\hat{\mathbf{z}} \quad (12d) \quad \text{Fe III}$$

$$\mathbf{B}_{30} = \left(\tfrac{3}{4} - y_5\right)\mathbf{a}_1 + \tfrac{3}{8}\,\mathbf{a}_2 - y_5\,\mathbf{a}_3 = \left(\tfrac{3}{4} - y_5\right)a\,\hat{\mathbf{x}} + \tfrac{3}{8}a\,\hat{\mathbf{y}} - y_5 a\,\hat{\mathbf{z}} \quad (12d) \quad \text{Fe III}$$

$$\mathbf{B}_{31} = \left(\tfrac{1}{4} + y_5\right)\mathbf{a}_1 + \tfrac{7}{8}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + y_5\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + y_5\right)a\,\hat{\mathbf{x}} + \tfrac{7}{8}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + y_5\right)a\,\hat{\mathbf{z}} \quad (12d) \quad \text{Fe III}$$

$$\mathbf{B}_{32} = \left(\tfrac{3}{4} + y_5\right)\mathbf{a}_1 + \tfrac{5}{8}\,\mathbf{a}_2 + \left(\tfrac{1}{2} - y_5\right)\mathbf{a}_3 = \left(\tfrac{3}{4} + y_5\right)a\,\hat{\mathbf{x}} + \tfrac{5}{8}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - y_5\right)a\,\hat{\mathbf{z}} \quad (12d) \quad \text{Fe III}$$

$$\mathbf{B}_{33} = y_5\,\mathbf{a}_1 + \left(\tfrac{1}{4} - y_5\right)\mathbf{a}_2 + \tfrac{1}{8}\,\mathbf{a}_3 = y_5 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - y_5\right)a\,\hat{\mathbf{y}} + \tfrac{1}{8}a\,\hat{\mathbf{z}} \quad (12d) \quad \text{Fe III}$$

$$\mathbf{B}_{34} = -y_5\,\mathbf{a}_1 + \left(\tfrac{3}{4} - y_5\right)\mathbf{a}_2 + \tfrac{3}{8}\,\mathbf{a}_3 = -y_5 a\,\hat{\mathbf{x}} + \left(\tfrac{3}{4} - y_5\right)a\,\hat{\mathbf{y}} + \tfrac{3}{8}a\,\hat{\mathbf{z}} \quad (12d) \quad \text{Fe III}$$

$$\mathbf{B}_{35} = \left(\tfrac{1}{2} + y_5\right)\mathbf{a}_1 + \left(\tfrac{1}{4} + y_5\right)\mathbf{a}_2 + \tfrac{7}{8}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + y_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + y_5\right)a\,\hat{\mathbf{y}} + \tfrac{7}{8}a\,\hat{\mathbf{z}} \quad (12d) \quad \text{Fe III}$$

$$\mathbf{B}_{36} = \left(\tfrac{1}{2} - y_5\right)\mathbf{a}_1 + \left(\tfrac{3}{4} + y_5\right)\mathbf{a}_2 + \tfrac{5}{8}\,\mathbf{a}_3 = \left(\tfrac{1}{2} - y_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{3}{4} + y_5\right)a\,\hat{\mathbf{y}} + \tfrac{5}{8}a\,\hat{\mathbf{z}} \quad (12d) \quad \text{Fe III}$$

$$\mathbf{B}_{37} = x_6\,\mathbf{a}_1 + y_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 = x_6 a\,\hat{\mathbf{x}} + y_6 a\,\hat{\mathbf{y}} + z_6 a\,\hat{\mathbf{z}} \quad (24e) \quad \text{O III}$$

$$\mathbf{B}_{38} = \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_1 - y_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_6\right)a\,\hat{\mathbf{x}} - y_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right)a\,\hat{\mathbf{z}} \quad (24e) \quad \text{O III}$$

$$\mathbf{B}_{39} = -x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_3 = -x_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_6\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_6\right)a\,\hat{\mathbf{z}} \quad (24e) \quad \text{O III}$$

$$\mathbf{B}_{40} = \left(\tfrac{1}{2} + x_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_6\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_6\right)a\,\hat{\mathbf{y}} - z_6 a\,\hat{\mathbf{z}} \quad (24e) \quad \text{O III}$$

$$\mathbf{B}_{41} = z_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + y_6\,\mathbf{a}_3 = z_6 a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} + y_6 a\,\hat{\mathbf{z}} \quad (24e) \quad \text{O III}$$

$$\mathbf{B}_{42} = \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_2 - y_6\,\mathbf{a}_3 = \left(\tfrac{1}{2} + z_6\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_6\right)a\,\hat{\mathbf{y}} - y_6 a\,\hat{\mathbf{z}} \quad (24e) \quad \text{O III}$$

$$\mathbf{B}_{43} = \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_1 - x_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} + y_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - z_6\right)a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + y_6\right)a\,\hat{\mathbf{z}} \quad (24e) \quad \text{O III}$$

$$\mathbf{B}_{44} = -z_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - y_6\right)\mathbf{a}_3 = -z_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_6\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - y_6\right)a\,\hat{\mathbf{z}} \quad (24e) \quad \text{O III}$$

$$\mathbf{B}_{45} = y_6\,\mathbf{a}_1 + z_6\,\mathbf{a}_2 + x_6\,\mathbf{a}_3 = y_6 a\,\hat{\mathbf{x}} + z_6 a\,\hat{\mathbf{y}} + x_6 a\,\hat{\mathbf{z}} \quad (24e) \quad \text{O III}$$

$$\mathbf{B}_{46} = -y_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_3 = -y_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + z_6\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_6\right)a\,\hat{\mathbf{z}} \quad (24e) \quad \text{O III}$$

$$\mathbf{B}_{47} = \left(\tfrac{1}{2} + y_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_2 - x_6\,\mathbf{a}_3 = \left(\tfrac{1}{2} + y_6\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - z_6\right)a\,\hat{\mathbf{y}} - x_6 a\,\hat{\mathbf{z}} \quad (24e) \quad \text{O III}$$

$$\mathbf{B}_{48} = \left(\tfrac{1}{2} - y_6\right)\mathbf{a}_1 - z_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - y_6\right)a\,\hat{\mathbf{x}} - z_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_6\right)a\,\hat{\mathbf{z}} \quad (24e) \quad \text{O III}$$

$$\mathbf{B}_{49} = \left(\tfrac{1}{4} + y_6\right)\mathbf{a}_1 + \left(\tfrac{3}{4} + x_6\right)\mathbf{a}_2 + \left(\tfrac{3}{4} - z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + y_6\right)a\,\hat{\mathbf{x}} + \left(\tfrac{3}{4} + x_6\right)a\,\hat{\mathbf{y}} + \left(\tfrac{3}{4} - z_6\right)a\,\hat{\mathbf{z}} \quad (24e) \quad \text{O III}$$

$$\mathbf{B}_{50} = \left(\tfrac{1}{4} - y_6\right)\mathbf{a}_1 + \left(\tfrac{1}{4} - x_6\right)\mathbf{a}_2 + \left(\tfrac{1}{4} - z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - y_6\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_6\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - z_6\right)a\,\hat{\mathbf{z}} \quad (24e) \quad \text{O III}$$

$$\mathbf{B}_{51} = \left(\tfrac{3}{4} + y_6\right)\mathbf{a}_1 + \left(\tfrac{3}{4} - x_6\right)\mathbf{a}_2 + \left(\tfrac{1}{4} + z_6\right)\mathbf{a}_3 = \left(\tfrac{3}{4} + y_6\right)a\,\hat{\mathbf{x}} + \left(\tfrac{3}{4} - x_6\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_6\right)a\,\hat{\mathbf{z}} \quad (24e) \quad \text{O III}$$

$$\mathbf{B}_{52} = \left(\tfrac{3}{4} - y_6\right)\mathbf{a}_1 + \left(\tfrac{1}{4} + x_6\right)\mathbf{a}_2 + \left(\tfrac{3}{4} + z_6\right)\mathbf{a}_3 = \left(\tfrac{3}{4} - y_6\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_6\right)a\,\hat{\mathbf{y}} + \left(\tfrac{3}{4} + z_6\right)a\,\hat{\mathbf{z}} \quad (24e) \quad \text{O III}$$

$$\mathbf{B}_{53} = \left(\tfrac{1}{4} + x_6\right)\mathbf{a}_1 + \left(\tfrac{3}{4} + z_6\right)\mathbf{a}_2 + \left(\tfrac{3}{4} - y_6\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + x_6\right)a\,\hat{\mathbf{x}} + \left(\tfrac{3}{4} + z_6\right)a\,\hat{\mathbf{y}} + \left(\tfrac{3}{4} - y_6\right)a\,\hat{\mathbf{z}} \quad (24e) \quad \text{O III}$$

$$\mathbf{B}_{54} = \left(\tfrac{3}{4} - x_6\right)\mathbf{a}_1 + \left(\tfrac{1}{4} + z_6\right)\mathbf{a}_2 + \left(\tfrac{3}{4} + y_6\right)\mathbf{a}_3 = \left(\tfrac{3}{4} - x_6\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + z_6\right)a\,\hat{\mathbf{y}} + \left(\tfrac{3}{4} + y_6\right)a\,\hat{\mathbf{z}} \quad (24e) \quad \text{O III}$$

$$\mathbf{B}_{55} = \left(\tfrac{1}{4} - x_6\right)\mathbf{a}_1 + \left(\tfrac{1}{4} - z_6\right)\mathbf{a}_2 + \left(\tfrac{1}{4} - y_6\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - x_6\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - z_6\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - y_6\right)a\,\hat{\mathbf{z}} \quad (24e) \quad \text{O III}$$

$$\mathbf{B}_{56} = \left(\tfrac{3}{4} + x_6\right)\mathbf{a}_1 + \left(\tfrac{3}{4} - z_6\right)\mathbf{a}_2 + \left(\tfrac{1}{4} + y_6\right)\mathbf{a}_3 = \left(\tfrac{3}{4} + x_6\right)a\,\hat{\mathbf{x}} + \left(\tfrac{3}{4} - z_6\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + y_6\right)a\,\hat{\mathbf{z}} \quad (24e) \quad \text{O III}$$

$\mathbf{B}_{57}$	=	$\left(\frac{1}{4}+z_6\right)\mathbf{a}_1+\left(\frac{3}{4}+y_6\right)\mathbf{a}_2+\left(\frac{3}{4}-x_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}+z_6\right)a\,\hat{\mathbf{x}}+\left(\frac{3}{4}+y_6\right)a\,\hat{\mathbf{y}}+\left(\frac{3}{4}-x_6\right)a\,\hat{\mathbf{z}}$	(24*e*)	O III
$\mathbf{B}_{58}$	=	$\left(\frac{3}{4}+z_6\right)\mathbf{a}_1+\left(\frac{3}{4}-y_6\right)\mathbf{a}_2+\left(\frac{1}{4}+x_6\right)\mathbf{a}_3$	=	$\left(\frac{3}{4}+z_6\right)a\,\hat{\mathbf{x}}+\left(\frac{3}{4}-y_6\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+x_6\right)a\,\hat{\mathbf{z}}$	(24*e*)	O III
$\mathbf{B}_{59}$	=	$\left(\frac{3}{4}-z_6\right)\mathbf{a}_1+\left(\frac{1}{4}+y_6\right)\mathbf{a}_2+\left(\frac{3}{4}+x_6\right)\mathbf{a}_3$	=	$\left(\frac{3}{4}-z_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+y_6\right)a\,\hat{\mathbf{y}}+\left(\frac{3}{4}+x_6\right)a\,\hat{\mathbf{z}}$	(24*e*)	O III
$\mathbf{B}_{60}$	=	$\left(\frac{1}{4}-z_6\right)\mathbf{a}_1+\left(\frac{1}{4}-y_6\right)\mathbf{a}_2+\left(\frac{1}{4}-x_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}-z_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-y_6\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-x_6\right)a\,\hat{\mathbf{z}}$	(24*e*)	O III

References:

- J. Thewlis, *The structure of ferromagnetic ferric oxide*, Philos. Mag. **12**, 1089–1106 (1931), doi:10.1080/14786443109461890.

Found in:

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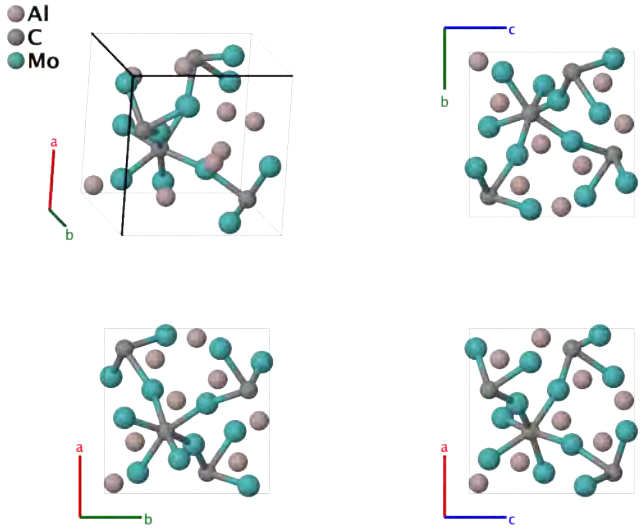
Geometry files:

- CIF: pp. 1788
- POSCAR: pp. 1789

Al_2Mo_3C Structure: A2BC3_cP24_213_c_a_d

http://aflow.org/prototype-encyclopedia/A2BC3_cP24_213_c_a_d



Prototype	:	Al_2CMo_3
AFLOW prototype label	:	A2BC3_cP24_213_c_a_d
***Strukturbericht* designation**	:	None
Pearson symbol	:	cP24
Space group number	:	213
Space group symbol	:	$P4_132$
AFLOW prototype command	:	`aflow --proto=A2BC3_cP24_213_c_a_d --params=`a, x_2, y_3

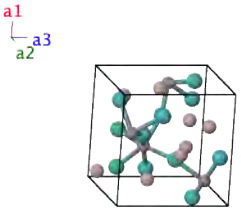
Other compounds with this structure

- Ag_2Pd_3Sn, Al_2Nb_3C, Al_2Nb_3N, Al_2Ta_3C, Cr_2Re_3B, Li_2Pd_3B, Li_2Pt_3B, Mn_2Rh_3P, Ni_2W_3N, Rh_2Mo_3N, and $(Fe_{2-x}Rh_x)Mo_3N$

- This is a "filled" β-Mn ($A13$) structure, with the aluminum and molybdenum atoms almost exactly on the sites of the manganese atoms in $A13$.

Simple Cubic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = a\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{3}{8}\,\mathbf{a}_1 + \frac{3}{8}\,\mathbf{a}_2 + \frac{3}{8}\,\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}}$	(4*a*)	C
$\mathbf{B}_2$	=	$\frac{1}{8}\,\mathbf{a}_1 + \frac{5}{8}\,\mathbf{a}_2 + \frac{7}{8}\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \frac{5}{8}a\,\hat{\mathbf{y}} + \frac{7}{8}a\,\hat{\mathbf{z}}$	(4*a*)	C
$\mathbf{B}_3$	=	$\frac{5}{8}\,\mathbf{a}_1 + \frac{7}{8}\,\mathbf{a}_2 + \frac{1}{8}\,\mathbf{a}_3$	=	$\frac{5}{8}a\,\hat{\mathbf{x}} + \frac{7}{8}a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(4*a*)	C
$\mathbf{B}_4$	=	$\frac{7}{8}\,\mathbf{a}_1 + \frac{1}{8}\,\mathbf{a}_2 + \frac{5}{8}\,\mathbf{a}_3$	=	$\frac{7}{8}a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \frac{5}{8}a\,\hat{\mathbf{z}}$	(4*a*)	C
$\mathbf{B}_5$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$x_2a\,\hat{\mathbf{x}} + x_2a\,\hat{\mathbf{y}} + x_2a\,\hat{\mathbf{z}}$	(8*c*)	Al
$\mathbf{B}_6$	=	$\left(\frac{1}{2} - x_2\right)\mathbf{a}_1 - x_2\,\mathbf{a}_2 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} - x_2a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{z}}$	(8*c*)	Al
$\mathbf{B}_7$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{1}{2} + x_2\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{z}}$	(8*c*)	Al
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_2\right)\mathbf{a}_2 - x_2\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{y}} - x_2a\,\hat{\mathbf{z}}$	(8*c*)	Al
$\mathbf{B}_9$	=	$\left(\frac{3}{4} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{4} + x_2\right)\mathbf{a}_2 + \left(\frac{1}{4} - x_2\right)\mathbf{a}_3$	=	$\left(\frac{3}{4} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_2\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_2\right)a\,\hat{\mathbf{z}}$	(8*c*)	Al
$\mathbf{B}_{10}$	=	$\left(\frac{3}{4} - x_2\right)\mathbf{a}_1 + \left(\frac{3}{4} - x_2\right)\mathbf{a}_2 + \left(\frac{3}{4} - x_2\right)\mathbf{a}_3$	=	$\left(\frac{3}{4} - x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{3}{4} - x_2\right)a\,\hat{\mathbf{y}} + \left(\frac{3}{4} - x_2\right)a\,\hat{\mathbf{z}}$	(8*c*)	Al
$\mathbf{B}_{11}$	=	$\left(\frac{1}{4} + x_2\right)\mathbf{a}_1 + \left(\frac{1}{4} - x_2\right)\mathbf{a}_2 + \left(\frac{3}{4} + x_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_2\right)a\,\hat{\mathbf{y}} + \left(\frac{3}{4} + x_2\right)a\,\hat{\mathbf{z}}$	(8*c*)	Al
$\mathbf{B}_{12}$	=	$\left(\frac{1}{4} - x_2\right)\mathbf{a}_1 + \left(\frac{3}{4} + x_2\right)\mathbf{a}_2 + \left(\frac{1}{4} + x_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{3}{4} + x_2\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + x_2\right)a\,\hat{\mathbf{z}}$	(8*c*)	Al
$\mathbf{B}_{13}$	=	$\frac{1}{8}\,\mathbf{a}_1 + y_3\,\mathbf{a}_2 + \left(\frac{1}{4} + y_3\right)\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + y_3a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + y_3\right)a\,\hat{\mathbf{z}}$	(12*d*)	Mo
$\mathbf{B}_{14}$	=	$\frac{3}{8}\,\mathbf{a}_1 - y_3\,\mathbf{a}_2 + \left(\frac{3}{4} + y_3\right)\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} - y_3a\,\hat{\mathbf{y}} + \left(\frac{3}{4} + y_3\right)a\,\hat{\mathbf{z}}$	(12*d*)	Mo
$\mathbf{B}_{15}$	=	$\frac{7}{8}\,\mathbf{a}_1 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_2 + \left(\frac{1}{4} - y_3\right)\mathbf{a}_3$	=	$\frac{7}{8}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - y_3\right)a\,\hat{\mathbf{z}}$	(12*d*)	Mo
$\mathbf{B}_{16}$	=	$\frac{5}{8}\,\mathbf{a}_1 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_2 + \left(\frac{3}{4} - y_3\right)\mathbf{a}_3$	=	$\frac{5}{8}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{y}} + \left(\frac{3}{4} - y_3\right)a\,\hat{\mathbf{z}}$	(12*d*)	Mo
$\mathbf{B}_{17}$	=	$\left(\frac{1}{4} + y_3\right)\mathbf{a}_1 + \frac{1}{8}\,\mathbf{a}_2 + y_3\,\mathbf{a}_3$	=	$\left(\frac{1}{4} + y_3\right)a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + y_3a\,\hat{\mathbf{z}}$	(12*d*)	Mo
$\mathbf{B}_{18}$	=	$\left(\frac{3}{4} + y_3\right)\mathbf{a}_1 + \frac{3}{8}\,\mathbf{a}_2 - y_3\,\mathbf{a}_3$	=	$\left(\frac{3}{4} + y_3\right)a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} - y_3a\,\hat{\mathbf{z}}$	(12*d*)	Mo
$\mathbf{B}_{19}$	=	$\left(\frac{1}{4} - y_3\right)\mathbf{a}_1 + \frac{7}{8}\,\mathbf{a}_2 + \left(\frac{1}{2} + y_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - y_3\right)a\,\hat{\mathbf{x}} + \frac{7}{8}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{z}}$	(12*d*)	Mo
$\mathbf{B}_{20}$	=	$\left(\frac{3}{4} - y_3\right)\mathbf{a}_1 + \frac{5}{8}\,\mathbf{a}_2 + \left(\frac{1}{2} - y_3\right)\mathbf{a}_3$	=	$\left(\frac{3}{4} - y_3\right)a\,\hat{\mathbf{x}} + \frac{5}{8}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{z}}$	(12*d*)	Mo
$\mathbf{B}_{21}$	=	$y_3\,\mathbf{a}_1 + \left(\frac{1}{4} + y_3\right)\mathbf{a}_2 + \frac{1}{8}\,\mathbf{a}_3$	=	$y_3a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + y_3\right)a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(12*d*)	Mo
$\mathbf{B}_{22}$	=	$-y_3\,\mathbf{a}_1 + \left(\frac{3}{4} + y_3\right)\mathbf{a}_2 + \frac{3}{8}\,\mathbf{a}_3$	=	$-y_3a\,\hat{\mathbf{x}} + \left(\frac{3}{4} + y_3\right)a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}}$	(12*d*)	Mo
$\mathbf{B}_{23}$	=	$\left(\frac{1}{2} + y_3\right)\mathbf{a}_1 + \left(\frac{1}{4} - y_3\right)\mathbf{a}_2 + \frac{7}{8}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + y_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - y_3\right)a\,\hat{\mathbf{y}} + \frac{7}{8}a\,\hat{\mathbf{z}}$	(12*d*)	Mo
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} - y_3\right)\mathbf{a}_1 + \left(\frac{3}{4} - y_3\right)\mathbf{a}_2 + \frac{5}{8}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{x}} + \left(\frac{3}{4} - y_3\right)a\,\hat{\mathbf{y}} + \frac{5}{8}a\,\hat{\mathbf{z}}$	(12*d*)	Mo

References:

- W. Jeitschko, H. Nowotny, and F. Benesovsky, *Ein Beitrag zum Dreistoff: Molybdän-Aluminium-Kohlenstoff*, Monatsh. Chem. Verw. Teile Anderer Wiss. **94**, 247–251 (1963), doi:10.1007/BF00900244.

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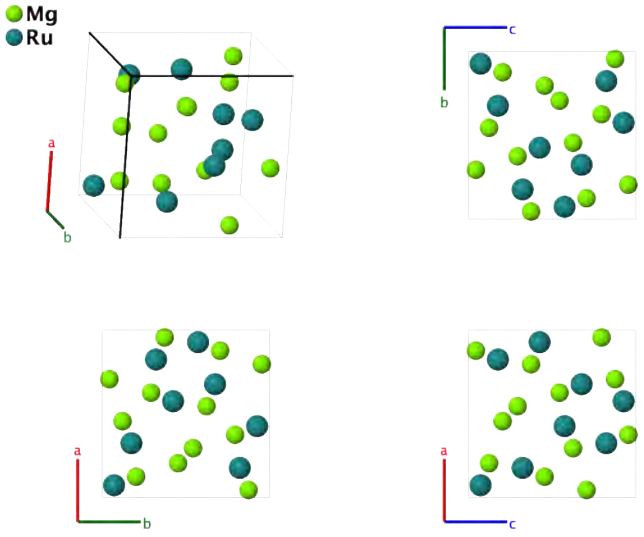
- A. Iyo, I. Hase, H. Fujihisa, Y. Gotoh, N. Takeshita, S. Ishida, H. Ninomiya, Y. Yoshida, H. Eisaki, and K. Kawashima, *Superconductivity induced by Mg deficiency in noncentrosymmetric phosphide Mg_2Rh_3P*, Phys. Rev. Mater. **3**, 124802 (2019), doi:10.1103/PhysRevMaterials.3.124802.
- J. Johnston, L. Toth, K. Kennedy, and E. R. Parker, *Superconductivity of Mo_3Al_2C*, Solid State Commun. **2**, 123 (1964), doi:10.1016/0038-1098(64)90251-0.

Geometry files:
- CIF: pp. 1789
- POSCAR: pp. 1790

Mg_3Ru_2 Structure: A3B2_cP20_213_d_c

http://aflow.org/prototype-encyclopedia/A3B2_cP20_213_d_c

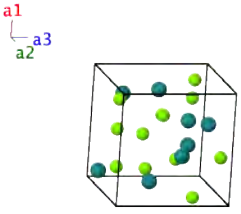


Prototype	:	Mg_3Ru_2
AFLOW prototype label	:	A3B2_cP20_213_d_c
***Strukturbericht* designation**	:	None
Pearson symbol	:	cP20
Space group number	:	213
Space group symbol	:	$P4_132$
AFLOW prototype command	:	`aflow --proto=A3B2_cP20_213_d_c --params=`a, x_1, y_2

- This is the binary form of the β-Mn ($A13$) structure.

Simple Cubic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = a\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$= x_1\,\mathbf{a}_1 + x_1\,\mathbf{a}_2 + x_1\,\mathbf{a}_3$	$=$	$x_1 a\,\hat{\mathbf{x}} + x_1 a\,\hat{\mathbf{y}} + x_1 a\,\hat{\mathbf{z}}$	($8c$)	Ru
$\mathbf{B}_2$	$= \left(\frac{1}{2} - x_1\right)\mathbf{a}_1 - x_1\,\mathbf{a}_2 + \left(\frac{1}{2} + x_1\right)\mathbf{a}_3$	$=$	$\left(\frac{1}{2} - x_1\right)a\,\hat{\mathbf{x}} - x_1 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_1\right)a\,\hat{\mathbf{z}}$	($8c$)	Ru
$\mathbf{B}_3$	$= -x_1\,\mathbf{a}_1 + \left(\frac{1}{2} + x_1\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_1\right)\mathbf{a}_3$	$=$	$-x_1 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_1\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_1\right)a\,\hat{\mathbf{z}}$	($8c$)	Ru

$\mathbf{B}_4$	=	$\left(\frac{1}{2}+x_1\right)\mathbf{a}_1+\left(\frac{1}{2}-x_1\right)\mathbf{a}_2-x_1\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_1\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_1\right)a\,\hat{\mathbf{y}}-x_1a\,\hat{\mathbf{z}}$	(8*c*)	Ru
$\mathbf{B}_5$	=	$\left(\frac{3}{4}+x_1\right)\mathbf{a}_1+\left(\frac{1}{4}+x_1\right)\mathbf{a}_2+\left(\frac{1}{4}-x_1\right)\mathbf{a}_3$	=	$\left(\frac{3}{4}+x_1\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+x_1\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-x_1\right)a\,\hat{\mathbf{z}}$	(8*c*)	Ru
$\mathbf{B}_6$	=	$\left(\frac{3}{4}-x_1\right)\mathbf{a}_1+\left(\frac{3}{4}-x_1\right)\mathbf{a}_2+\left(\frac{3}{4}-x_1\right)\mathbf{a}_3$	=	$\left(\frac{3}{4}-x_1\right)a\,\hat{\mathbf{x}}+\left(\frac{3}{4}-x_1\right)a\,\hat{\mathbf{y}}+\left(\frac{3}{4}-x_1\right)a\,\hat{\mathbf{z}}$	(8*c*)	Ru
$\mathbf{B}_7$	=	$\left(\frac{1}{4}+x_1\right)\mathbf{a}_1+\left(\frac{1}{4}-x_1\right)\mathbf{a}_2+\left(\frac{3}{4}+x_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}+x_1\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-x_1\right)a\,\hat{\mathbf{y}}+\left(\frac{3}{4}+x_1\right)a\,\hat{\mathbf{z}}$	(8*c*)	Ru
$\mathbf{B}_8$	=	$\left(\frac{1}{4}-x_1\right)\mathbf{a}_1+\left(\frac{3}{4}+x_1\right)\mathbf{a}_2+\left(\frac{1}{4}+x_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}-x_1\right)a\,\hat{\mathbf{x}}+\left(\frac{3}{4}+x_1\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+x_1\right)a\,\hat{\mathbf{z}}$	(8*c*)	Ru
$\mathbf{B}_9$	=	$\frac{1}{8}\mathbf{a}_1+y_2\mathbf{a}_2+\left(\frac{1}{4}+y_2\right)\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}}+y_2a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+y_2\right)a\,\hat{\mathbf{z}}$	(12*d*)	Mg
$\mathbf{B}_{10}$	=	$\frac{3}{8}\mathbf{a}_1-y_2\mathbf{a}_2+\left(\frac{3}{4}+y_2\right)\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}}-y_2a\,\hat{\mathbf{y}}+\left(\frac{3}{4}+y_2\right)a\,\hat{\mathbf{z}}$	(12*d*)	Mg
$\mathbf{B}_{11}$	=	$\frac{7}{8}\mathbf{a}_1+\left(\frac{1}{2}+y_2\right)\mathbf{a}_2+\left(\frac{1}{4}-y_2\right)\mathbf{a}_3$	=	$\frac{7}{8}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+y_2\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-y_2\right)a\,\hat{\mathbf{z}}$	(12*d*)	Mg
$\mathbf{B}_{12}$	=	$\frac{5}{8}\mathbf{a}_1+\left(\frac{1}{2}-y_2\right)\mathbf{a}_2+\left(\frac{3}{4}-y_2\right)\mathbf{a}_3$	=	$\frac{5}{8}a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_2\right)a\,\hat{\mathbf{y}}+\left(\frac{3}{4}-y_2\right)a\,\hat{\mathbf{z}}$	(12*d*)	Mg
$\mathbf{B}_{13}$	=	$\left(\frac{1}{4}+y_2\right)\mathbf{a}_1+\frac{1}{8}\mathbf{a}_2+y_2\mathbf{a}_3$	=	$\left(\frac{1}{4}+y_2\right)a\,\hat{\mathbf{x}}+\frac{1}{8}a\,\hat{\mathbf{y}}+y_2a\,\hat{\mathbf{z}}$	(12*d*)	Mg
$\mathbf{B}_{14}$	=	$\left(\frac{3}{4}+y_2\right)\mathbf{a}_1+\frac{3}{8}\mathbf{a}_2-y_2\mathbf{a}_3$	=	$\left(\frac{3}{4}+y_2\right)a\,\hat{\mathbf{x}}+\frac{3}{8}a\,\hat{\mathbf{y}}-y_2a\,\hat{\mathbf{z}}$	(12*d*)	Mg
$\mathbf{B}_{15}$	=	$\left(\frac{1}{4}-y_2\right)\mathbf{a}_1+\frac{7}{8}\mathbf{a}_2+\left(\frac{1}{2}+y_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}-y_2\right)a\,\hat{\mathbf{x}}+\frac{7}{8}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+y_2\right)a\,\hat{\mathbf{z}}$	(12*d*)	Mg
$\mathbf{B}_{16}$	=	$\left(\frac{3}{4}-y_2\right)\mathbf{a}_1+\frac{5}{8}\mathbf{a}_2+\left(\frac{1}{2}-y_2\right)\mathbf{a}_3$	=	$\left(\frac{3}{4}-y_2\right)a\,\hat{\mathbf{x}}+\frac{5}{8}a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-y_2\right)a\,\hat{\mathbf{z}}$	(12*d*)	Mg
$\mathbf{B}_{17}$	=	$y_2\mathbf{a}_1+\left(\frac{1}{4}+y_2\right)\mathbf{a}_2+\frac{1}{8}\mathbf{a}_3$	=	$y_2a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+y_2\right)a\,\hat{\mathbf{y}}+\frac{1}{8}a\,\hat{\mathbf{z}}$	(12*d*)	Mg
$\mathbf{B}_{18}$	=	$-y_2\mathbf{a}_1+\left(\frac{3}{4}+y_2\right)\mathbf{a}_2+\frac{3}{8}\mathbf{a}_3$	=	$-y_2a\,\hat{\mathbf{x}}+\left(\frac{3}{4}+y_2\right)a\,\hat{\mathbf{y}}+\frac{3}{8}a\,\hat{\mathbf{z}}$	(12*d*)	Mg
$\mathbf{B}_{19}$	=	$\left(\frac{1}{2}+y_2\right)\mathbf{a}_1+\left(\frac{1}{4}-y_2\right)\mathbf{a}_2+\frac{7}{8}\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_2\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-y_2\right)a\,\hat{\mathbf{y}}+\frac{7}{8}a\,\hat{\mathbf{z}}$	(12*d*)	Mg
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2}-y_2\right)\mathbf{a}_1+\left(\frac{3}{4}-y_2\right)\mathbf{a}_2+\frac{5}{8}\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_2\right)a\,\hat{\mathbf{x}}+\left(\frac{3}{4}-y_2\right)a\,\hat{\mathbf{y}}+\frac{5}{8}a\,\hat{\mathbf{z}}$	(12*d*)	Mg

References:

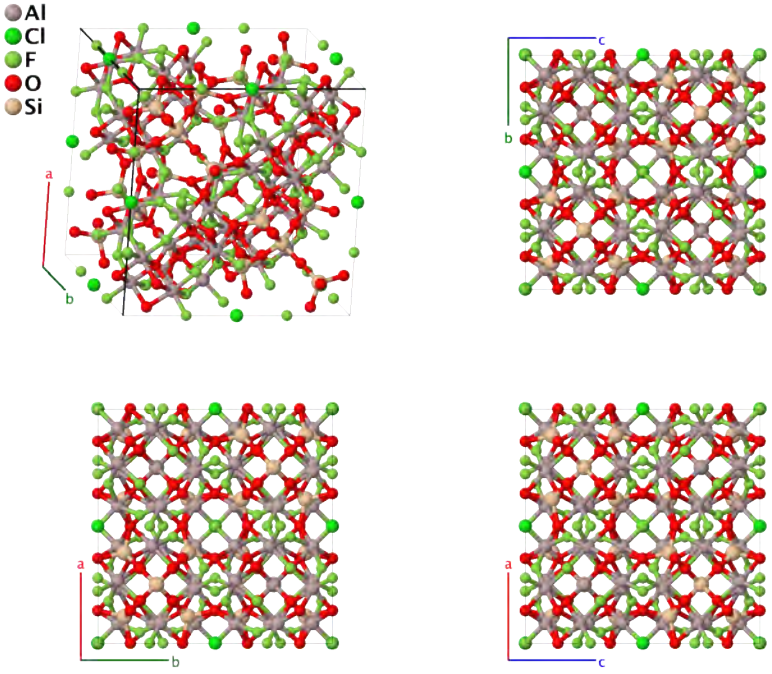
- R. Pöttgen, V. Hlukhyy, A. Baranov, and Y. Grin, *Crystal Structure and Chemical Bonding of Mg_3Ru_2*, Inorg. Chem. **47**, 6051–6055 (2008), doi:10.1021/ic800387a.

Geometry files:

- CIF: pp. 1790
- POSCAR: pp. 1790

Zunyite [$Al_{13}(OH,F)_{18}Si_5O_{20}Cl$, $S0_8$] Structure: A13BC18D20E5_cF228_216_dh_b_fh_2eh_ce

http://aflow.org/prototype-encyclopedia/A13BC18D20E5_cF228_216_dh_b_fh_2eh_ce

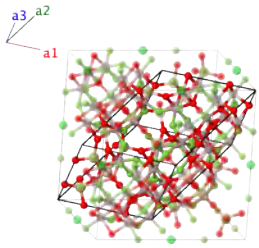


Prototype	:	$Al_{13}ClF_{18}O_{20}Si_5$
AFLOW prototype label	:	A13BC18D20E5_cF228_216_dh_b_fh_2eh_ce
***Strukturbericht* designation**	:	$S0_8$
Pearson symbol	:	cF228
Space group number	:	216
Space group symbol	:	$F\bar{4}3m$
AFLOW prototype command	:	`aflow --proto=A13BC18D20E5_cF228_216_dh_b_fh_2eh_ce --params=`$a, x_4, x_5, x_6, x_7, x_8, z_8, x_9, z_9, x_{10}, z_{10}$

- We use the structure proposed by (Kamb, 1960), a refinement of the original (Pauling, 1933) $S0_8$ structure. The only major difference is the y coordinate of the OH/F-I site, which is now at a more reasonable distance from the chlorine atom.
- For easy of visualization, we have used fluorine atoms to represent all of the $(OH,F)_{18}$ positions, but in reality the system is dominated by OH, not F. Kamb argues that the physics of hydrogen bonding makes it likely that the actual structure has composition $(OH)_{16}F_2$, with the fluorine atoms substituting for OH on the second ($48h$) site.

Face-centered Cubic primitive vectors:

$$\mathbf{a}_1 = \tfrac{1}{2}\,a\,\hat{\mathbf{y}} + \tfrac{1}{2}\,a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{y}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(4*b*)	Cl
$\mathbf{B}_2$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(4*c*)	Si I
$\mathbf{B}_3$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(4*d*)	Al I
$\mathbf{B}_4$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	(16*e*)	O I
$\mathbf{B}_5$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 - 3x_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	(16*e*)	O I
$\mathbf{B}_6$	=	$x_4\,\mathbf{a}_1 - 3x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} - x_4a\,\hat{\mathbf{z}}$	(16*e*)	O I
$\mathbf{B}_7$	=	$-3x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} - x_4a\,\hat{\mathbf{z}}$	(16*e*)	O I
$\mathbf{B}_8$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}} + x_5a\,\hat{\mathbf{y}} + x_5a\,\hat{\mathbf{z}}$	(16*e*)	O II
$\mathbf{B}_9$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 - 3x_5\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}} - x_5a\,\hat{\mathbf{y}} + x_5a\,\hat{\mathbf{z}}$	(16*e*)	O II
$\mathbf{B}_{10}$	=	$x_5\,\mathbf{a}_1 - 3x_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}} + x_5a\,\hat{\mathbf{y}} - x_5a\,\hat{\mathbf{z}}$	(16*e*)	O II
$\mathbf{B}_{11}$	=	$-3x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}} - x_5a\,\hat{\mathbf{y}} - x_5a\,\hat{\mathbf{z}}$	(16*e*)	O II
$\mathbf{B}_{12}$	=	$x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + x_6\,\mathbf{a}_3$	=	$x_6a\,\hat{\mathbf{x}} + x_6a\,\hat{\mathbf{y}} + x_6a\,\hat{\mathbf{z}}$	(16*e*)	Si II
$\mathbf{B}_{13}$	=	$x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 - 3x_6\,\mathbf{a}_3$	=	$-x_6a\,\hat{\mathbf{x}} - x_6a\,\hat{\mathbf{y}} + x_6a\,\hat{\mathbf{z}}$	(16*e*)	Si II
$\mathbf{B}_{14}$	=	$x_6\,\mathbf{a}_1 - 3x_6\,\mathbf{a}_2 + x_6\,\mathbf{a}_3$	=	$-x_6a\,\hat{\mathbf{x}} + x_6a\,\hat{\mathbf{y}} - x_6a\,\hat{\mathbf{z}}$	(16*e*)	Si II
$\mathbf{B}_{15}$	=	$-3x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + x_6\,\mathbf{a}_3$	=	$x_6a\,\hat{\mathbf{x}} - x_6a\,\hat{\mathbf{y}} - x_6a\,\hat{\mathbf{z}}$	(16*e*)	Si II
$\mathbf{B}_{16}$	=	$-x_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + x_7\,\mathbf{a}_3$	=	$x_7a\,\hat{\mathbf{x}}$	(24*f*)	F I
$\mathbf{B}_{17}$	=	$x_7\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 - x_7\,\mathbf{a}_3$	=	$-x_7a\,\hat{\mathbf{x}}$	(24*f*)	F I
$\mathbf{B}_{18}$	=	$x_7\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 + x_7\,\mathbf{a}_3$	=	$x_7a\,\hat{\mathbf{y}}$	(24*f*)	F I
$\mathbf{B}_{19}$	=	$-x_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 - x_7\,\mathbf{a}_3$	=	$-x_7a\,\hat{\mathbf{y}}$	(24*f*)	F I
$\mathbf{B}_{20}$	=	$x_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 - x_7\,\mathbf{a}_3$	=	$x_7a\,\hat{\mathbf{z}}$	(24*f*)	F I
$\mathbf{B}_{21}$	=	$-x_7\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 + x_7\,\mathbf{a}_3$	=	$-x_7a\,\hat{\mathbf{z}}$	(24*f*)	F I
$\mathbf{B}_{22}$	=	$z_8\,\mathbf{a}_1 + z_8\,\mathbf{a}_2 + (2x_8 - z_8)\,\mathbf{a}_3$	=	$x_8a\,\hat{\mathbf{x}} + x_8a\,\hat{\mathbf{y}} + z_8a\,\hat{\mathbf{z}}$	(48*h*)	Al II
$\mathbf{B}_{23}$	=	$z_8\,\mathbf{a}_1 + z_8\,\mathbf{a}_2 + (-2x_8 - z_8)\,\mathbf{a}_3$	=	$-x_8a\,\hat{\mathbf{x}} - x_8a\,\hat{\mathbf{y}} + z_8a\,\hat{\mathbf{z}}$	(48*h*)	Al II
$\mathbf{B}_{24}$	=	$(2x_8 - z_8)\,\mathbf{a}_1 + (-2x_8 - z_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$-x_8a\,\hat{\mathbf{x}} + x_8a\,\hat{\mathbf{y}} - z_8a\,\hat{\mathbf{z}}$	(48*h*)	Al II
$\mathbf{B}_{25}$	=	$(-2x_8 - z_8)\,\mathbf{a}_1 + (2x_8 - z_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$x_8a\,\hat{\mathbf{x}} - x_8a\,\hat{\mathbf{y}} - z_8a\,\hat{\mathbf{z}}$	(48*h*)	Al II
$\mathbf{B}_{26}$	=	$(2x_8 - z_8)\,\mathbf{a}_1 + z_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$z_8a\,\hat{\mathbf{x}} + x_8a\,\hat{\mathbf{y}} + x_8a\,\hat{\mathbf{z}}$	(48*h*)	Al II
$\mathbf{B}_{27}$	=	$(-2x_8 - z_8)\,\mathbf{a}_1 + z_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$z_8a\,\hat{\mathbf{x}} - x_8a\,\hat{\mathbf{y}} - x_8a\,\hat{\mathbf{z}}$	(48*h*)	Al II

$\mathbf{B}_{28}$	=	$z_8\,\mathbf{a}_1 + (2x_8 - z_8)\,\mathbf{a}_2 + (-2x_8 - z_8)\,\mathbf{a}_3$	=	$-z_8a\,\hat{\mathbf{x}} - x_8a\,\hat{\mathbf{y}} + x_8a\,\hat{\mathbf{z}}$	(48*h*)	Al II
$\mathbf{B}_{29}$	=	$z_8\,\mathbf{a}_1 + (-2x_8 - z_8)\,\mathbf{a}_2 + (2x_8 - z_8)\,\mathbf{a}_3$	=	$-z_8a\,\hat{\mathbf{x}} + x_8a\,\hat{\mathbf{y}} - x_8a\,\hat{\mathbf{z}}$	(48*h*)	Al II
$\mathbf{B}_{30}$	=	$z_8\,\mathbf{a}_1 + (2x_8 - z_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$x_8a\,\hat{\mathbf{x}} + z_8a\,\hat{\mathbf{y}} + x_8a\,\hat{\mathbf{z}}$	(48*h*)	Al II
$\mathbf{B}_{31}$	=	$z_8\,\mathbf{a}_1 + (-2x_8 - z_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$-x_8a\,\hat{\mathbf{x}} + z_8a\,\hat{\mathbf{y}} - x_8a\,\hat{\mathbf{z}}$	(48*h*)	Al II
$\mathbf{B}_{32}$	=	$(-2x_8 - z_8)\,\mathbf{a}_1 + z_8\,\mathbf{a}_2 + (2x_8 - z_8)\,\mathbf{a}_3$	=	$x_8a\,\hat{\mathbf{x}} - z_8a\,\hat{\mathbf{y}} - x_8a\,\hat{\mathbf{z}}$	(48*h*)	Al II
$\mathbf{B}_{33}$	=	$(2x_8 - z_8)\,\mathbf{a}_1 + z_8\,\mathbf{a}_2 + (-2x_8 - z_8)\,\mathbf{a}_3$	=	$-x_8a\,\hat{\mathbf{x}} - z_8a\,\hat{\mathbf{y}} + x_8a\,\hat{\mathbf{z}}$	(48*h*)	Al II
$\mathbf{B}_{34}$	=	$z_9\,\mathbf{a}_1 + z_9\,\mathbf{a}_2 + (2x_9 - z_9)\,\mathbf{a}_3$	=	$x_9a\,\hat{\mathbf{x}} + x_9a\,\hat{\mathbf{y}} + z_9a\,\hat{\mathbf{z}}$	(48*h*)	F II
$\mathbf{B}_{35}$	=	$z_9\,\mathbf{a}_1 + z_9\,\mathbf{a}_2 + (-2x_9 - z_9)\,\mathbf{a}_3$	=	$-x_9a\,\hat{\mathbf{x}} - x_9a\,\hat{\mathbf{y}} + z_9a\,\hat{\mathbf{z}}$	(48*h*)	F II
$\mathbf{B}_{36}$	=	$(2x_9 - z_9)\,\mathbf{a}_1 + (-2x_9 - z_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$-x_9a\,\hat{\mathbf{x}} + x_9a\,\hat{\mathbf{y}} - z_9a\,\hat{\mathbf{z}}$	(48*h*)	F II
$\mathbf{B}_{37}$	=	$(-2x_9 - z_9)\,\mathbf{a}_1 + (2x_9 - z_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$x_9a\,\hat{\mathbf{x}} - x_9a\,\hat{\mathbf{y}} - z_9a\,\hat{\mathbf{z}}$	(48*h*)	F II
$\mathbf{B}_{38}$	=	$(2x_9 - z_9)\,\mathbf{a}_1 + z_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$z_9a\,\hat{\mathbf{x}} + x_9a\,\hat{\mathbf{y}} + x_9a\,\hat{\mathbf{z}}$	(48*h*)	F II
$\mathbf{B}_{39}$	=	$(-2x_9 - z_9)\,\mathbf{a}_1 + z_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$z_9a\,\hat{\mathbf{x}} - x_9a\,\hat{\mathbf{y}} - x_9a\,\hat{\mathbf{z}}$	(48*h*)	F II
$\mathbf{B}_{40}$	=	$z_9\,\mathbf{a}_1 + (2x_9 - z_9)\,\mathbf{a}_2 + (-2x_9 - z_9)\,\mathbf{a}_3$	=	$-z_9a\,\hat{\mathbf{x}} - x_9a\,\hat{\mathbf{y}} + x_9a\,\hat{\mathbf{z}}$	(48*h*)	F II
$\mathbf{B}_{41}$	=	$z_9\,\mathbf{a}_1 + (-2x_9 - z_9)\,\mathbf{a}_2 + (2x_9 - z_9)\,\mathbf{a}_3$	=	$-z_9a\,\hat{\mathbf{x}} + x_9a\,\hat{\mathbf{y}} - x_9a\,\hat{\mathbf{z}}$	(48*h*)	F II
$\mathbf{B}_{42}$	=	$z_9\,\mathbf{a}_1 + (2x_9 - z_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$x_9a\,\hat{\mathbf{x}} + z_9a\,\hat{\mathbf{y}} + x_9a\,\hat{\mathbf{z}}$	(48*h*)	F II
$\mathbf{B}_{43}$	=	$z_9\,\mathbf{a}_1 + (-2x_9 - z_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3$	=	$-x_9a\,\hat{\mathbf{x}} + z_9a\,\hat{\mathbf{y}} - x_9a\,\hat{\mathbf{z}}$	(48*h*)	F II
$\mathbf{B}_{44}$	=	$(-2x_9 - z_9)\,\mathbf{a}_1 + z_9\,\mathbf{a}_2 + (2x_9 - z_9)\,\mathbf{a}_3$	=	$x_9a\,\hat{\mathbf{x}} - z_9a\,\hat{\mathbf{y}} - x_9a\,\hat{\mathbf{z}}$	(48*h*)	F II
$\mathbf{B}_{45}$	=	$(2x_9 - z_9)\,\mathbf{a}_1 + z_9\,\mathbf{a}_2 + (-2x_9 - z_9)\,\mathbf{a}_3$	=	$-x_9a\,\hat{\mathbf{x}} - z_9a\,\hat{\mathbf{y}} + x_9a\,\hat{\mathbf{z}}$	(48*h*)	F II
$\mathbf{B}_{46}$	=	$z_{10}\,\mathbf{a}_1 + z_{10}\,\mathbf{a}_2 + (2x_{10} - z_{10})\,\mathbf{a}_3$	=	$x_{10}a\,\hat{\mathbf{x}} + x_{10}a\,\hat{\mathbf{y}} + z_{10}a\,\hat{\mathbf{z}}$	(48*h*)	O III
$\mathbf{B}_{47}$	=	$z_{10}\,\mathbf{a}_1 + z_{10}\,\mathbf{a}_2 + (-2x_{10} - z_{10})\,\mathbf{a}_3$	=	$-x_{10}a\,\hat{\mathbf{x}} - x_{10}a\,\hat{\mathbf{y}} + z_{10}a\,\hat{\mathbf{z}}$	(48*h*)	O III
$\mathbf{B}_{48}$	=	$(2x_{10} - z_{10})\,\mathbf{a}_1 + (-2x_{10} - z_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$-x_{10}a\,\hat{\mathbf{x}} + x_{10}a\,\hat{\mathbf{y}} - z_{10}a\,\hat{\mathbf{z}}$	(48*h*)	O III
$\mathbf{B}_{49}$	=	$(-2x_{10} - z_{10})\,\mathbf{a}_1 + (2x_{10} - z_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$x_{10}a\,\hat{\mathbf{x}} - x_{10}a\,\hat{\mathbf{y}} - z_{10}a\,\hat{\mathbf{z}}$	(48*h*)	O III
$\mathbf{B}_{50}$	=	$(2x_{10} - z_{10})\,\mathbf{a}_1 + z_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$z_{10}a\,\hat{\mathbf{x}} + x_{10}a\,\hat{\mathbf{y}} + x_{10}a\,\hat{\mathbf{z}}$	(48*h*)	O III
$\mathbf{B}_{51}$	=	$(-2x_{10} - z_{10})\,\mathbf{a}_1 + z_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$z_{10}a\,\hat{\mathbf{x}} - x_{10}a\,\hat{\mathbf{y}} - x_{10}a\,\hat{\mathbf{z}}$	(48*h*)	O III
$\mathbf{B}_{52}$	=	$z_{10}\,\mathbf{a}_1 + (2x_{10} - z_{10})\,\mathbf{a}_2 + (-2x_{10} - z_{10})\,\mathbf{a}_3$	=	$-z_{10}a\,\hat{\mathbf{x}} - x_{10}a\,\hat{\mathbf{y}} + x_{10}a\,\hat{\mathbf{z}}$	(48*h*)	O III
$\mathbf{B}_{53}$	=	$z_{10}\,\mathbf{a}_1 + (-2x_{10} - z_{10})\,\mathbf{a}_2 + (2x_{10} - z_{10})\,\mathbf{a}_3$	=	$-z_{10}a\,\hat{\mathbf{x}} + x_{10}a\,\hat{\mathbf{y}} - x_{10}a\,\hat{\mathbf{z}}$	(48*h*)	O III
$\mathbf{B}_{54}$	=	$z_{10}\,\mathbf{a}_1 + (2x_{10} - z_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$x_{10}a\,\hat{\mathbf{x}} + z_{10}a\,\hat{\mathbf{y}} + x_{10}a\,\hat{\mathbf{z}}$	(48*h*)	O III
$\mathbf{B}_{55}$	=	$z_{10}\,\mathbf{a}_1 + (-2x_{10} - z_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$-x_{10}a\,\hat{\mathbf{x}} + z_{10}a\,\hat{\mathbf{y}} - x_{10}a\,\hat{\mathbf{z}}$	(48*h*)	O III
$\mathbf{B}_{56}$	=	$(-2x_{10} - z_{10})\,\mathbf{a}_1 + z_{10}\,\mathbf{a}_2 + (2x_{10} - z_{10})\,\mathbf{a}_3$	=	$x_{10}a\,\hat{\mathbf{x}} - z_{10}a\,\hat{\mathbf{y}} - x_{10}a\,\hat{\mathbf{z}}$	(48*h*)	O III
$\mathbf{B}_{57}$	=	$(2x_{10} - z_{10})\,\mathbf{a}_1 + z_{10}\,\mathbf{a}_2 + (-2x_{10} - z_{10})\,\mathbf{a}_3$	=	$-x_{10}a\,\hat{\mathbf{x}} - z_{10}a\,\hat{\mathbf{y}} + x_{10}a\,\hat{\mathbf{z}}$	(48*h*)	O III

References:

- W. B. Kamb, *The Crystal Structure of Zunyite*, Acta Cryst. **13**, 15–24 (1960), doi:10.1107/S0365110X60000030.
- L. Pauling, *The Crystal Structure of Zunyite, $Al_{13}Si_5O_{20}(OH,F)_{18}Cl$*, Zeitschrift für Kristallographie - Crystalline Materials **84**, 442–452 (1933), doi:10.1524/zkri.1933.84.1.442.

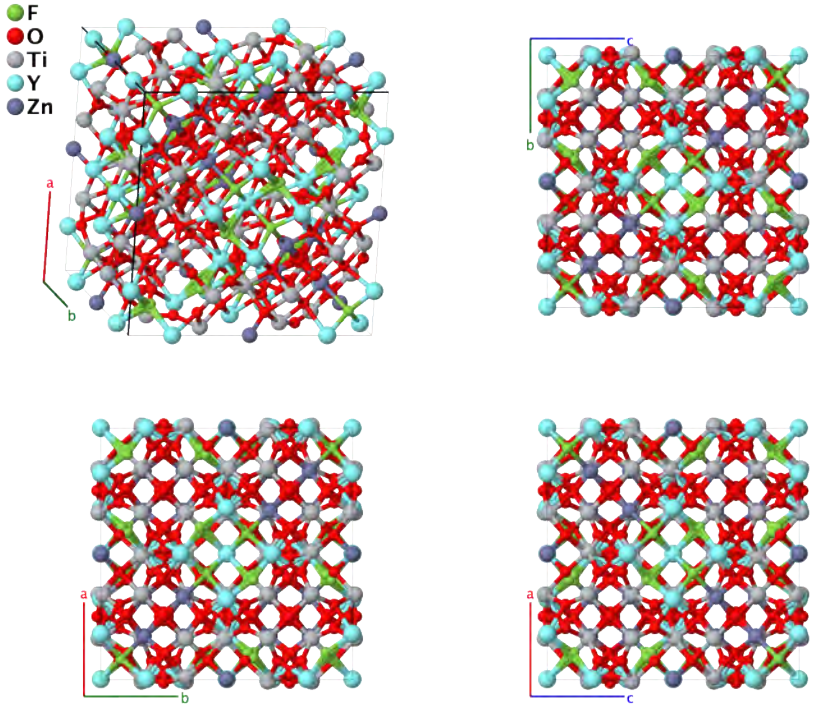
Geometry files:

- CIF: pp. 1790

- POSCAR: pp. 1791

Murataite $[(Y,Na)_6(Zn,Fe)_5Ti_{12}O_{29}(O,F)_{10}F_4]$ Structure: A16B40C12D6E5_cF316_216_eh_e2g2h_h_f_be

http://aflow.org/prototype-encyclopedia/A16B40C12D6E5_cF316_216_eh_e2g2h_h_f_be



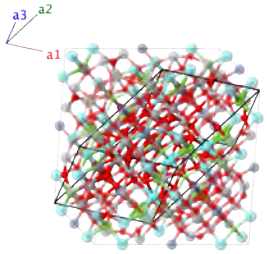
Prototype	:	$F_{16}O_{40}Ti_{12}Y_6Zn_5$
AFLOW prototype label	:	A16B40C12D6E5_cF316_216_eh_e2g2h_h_f_be
***Strukturbericht* designation**	:	None
Pearson symbol	:	cF316
Space group number	:	216
Space group symbol	:	$F\bar{4}3m$
AFLOW prototype command	:	`aflow --proto=A16B40C12D6E5_cF316_216_eh_e2g2h_h_f_be --params=`$a, x_2, x_3, x_4, x_5, x_6, x_7, x_8, z_8, x_9, z_9, x_{10}, z_{10}, x_{11}, z_{11}$

- Most of the sites in this structure are somewhat disordered. The "nominal" composition is given as $F_{16}O_{40}Ti_{12}Y_6Zn_5$ by (Ercit, 1995), but as the CIF in (Downs, 2003) shows, even these labels are not quite correct. In our listing we label each Wyckoff position by the type of atom that has the largest concentration on that site. Following (Downs, 2003):
  - Site Zn-I has the composition $Zn_{0.89}Si_{0.11}$.
  - Site F-I has the composition $F_{0.55}O_{0.45}$.
  - Site O-I is pure oxygen.
  - Site Zn-II has the composition $Zn_{0.48}Fe_{0.25}Na_{0.16}Ti_{0.11}$.
  - Site Y has the composition $Y_{0.37}Na_{0.35}Mn_{0.03}HREE_{0.25}$, where "HREE" is a mixture of heavy Rare Earth elements.
  - Site O-II is pure oxygen, but only 8.3333% of the sites are occupied.
  - Site O-III has the composition $O_{0.7}F_{0.3}$.

- Site O-IV is pure oxygen.
- Site O-V is pure oxygen, but only 87% of the sites are occupied.
- Site F-II is pure fluorine, but only 33.333% of the sites are occupied.
- Site Ti has the composition $Ti_{0.76}Nb_{0.13}Na_{0.11}$.

Face-centered Cubic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{y}} + \tfrac{1}{2}\,a\,\hat{\mathbf{z}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{z}}\\ \mathbf{a}_3 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{y}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(4*b*)	Zn I
$\mathbf{B}_2$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$x_2a\,\hat{\mathbf{x}} + x_2a\,\hat{\mathbf{y}} + x_2a\,\hat{\mathbf{z}}$	(16*e*)	F I
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 - 3x_2\,\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}} - x_2a\,\hat{\mathbf{y}} + x_2a\,\hat{\mathbf{z}}$	(16*e*)	F I
$\mathbf{B}_4$	=	$x_2\,\mathbf{a}_1 - 3x_2\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}} + x_2a\,\hat{\mathbf{y}} - x_2a\,\hat{\mathbf{z}}$	(16*e*)	F I
$\mathbf{B}_5$	=	$-3x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$x_2a\,\hat{\mathbf{x}} - x_2a\,\hat{\mathbf{y}} - x_2a\,\hat{\mathbf{z}}$	(16*e*)	F I
$\mathbf{B}_6$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(16*e*)	O I
$\mathbf{B}_7$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 - 3x_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(16*e*)	O I
$\mathbf{B}_8$	=	$x_3\,\mathbf{a}_1 - 3x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(16*e*)	O I
$\mathbf{B}_9$	=	$-3x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(16*e*)	O I
$\mathbf{B}_{10}$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	(16*e*)	Zn II
$\mathbf{B}_{11}$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 - 3x_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	(16*e*)	Zn II
$\mathbf{B}_{12}$	=	$x_4\,\mathbf{a}_1 - 3x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} - x_4a\,\hat{\mathbf{z}}$	(16*e*)	Zn II
$\mathbf{B}_{13}$	=	$-3x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} - x_4a\,\hat{\mathbf{z}}$	(16*e*)	Zn II
$\mathbf{B}_{14}$	=	$-x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}}$	(24*f*)	Y
$\mathbf{B}_{15}$	=	$x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - x_5\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}}$	(24*f*)	Y
$\mathbf{B}_{16}$	=	$x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{y}}$	(24*f*)	Y
$\mathbf{B}_{17}$	=	$-x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 - x_5\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{y}}$	(24*f*)	Y
$\mathbf{B}_{18}$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 - x_5\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{z}}$	(24*f*)	Y
$\mathbf{B}_{19}$	=	$-x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{z}}$	(24*f*)	Y
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} - x_6\right)\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + x_6\,\mathbf{a}_3$	=	$x_6a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(24*g*)	O II
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1 + \left(\frac{1}{2} - x_6\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - x_6\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(24*g*)	O II
$\mathbf{B}_{22}$	=	$x_6\,\mathbf{a}_1 + \left(\frac{1}{2} - x_6\right)\,\mathbf{a}_2 + x_6\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + x_6a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(24*g*)	O II
$\mathbf{B}_{23}$	=	$\left(\frac{1}{2} - x_6\right)\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + \left(\frac{1}{2} - x_6\right)\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_6\right)a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(24*g*)	O II

$$\mathbf{B}_{24} = x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_3 = \tfrac{1}{4}a\,\hat{\mathbf{x}} + \tfrac{1}{4}a\,\hat{\mathbf{y}} + x_6 a\,\hat{\mathbf{z}} \quad (24g) \quad \text{O II}$$

$$\mathbf{B}_{25} = \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_2 + x_6\,\mathbf{a}_3 = \tfrac{1}{4}a\,\hat{\mathbf{x}} + \tfrac{1}{4}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_6\right)a\,\hat{\mathbf{z}} \quad (24g) \quad \text{O II}$$

$$\mathbf{B}_{26} = \left(\tfrac{1}{2} - x_7\right)\mathbf{a}_1 + x_7\,\mathbf{a}_2 + x_7\,\mathbf{a}_3 = x_7 a\,\hat{\mathbf{x}} + \tfrac{1}{4}a\,\hat{\mathbf{y}} + \tfrac{1}{4}a\,\hat{\mathbf{z}} \quad (24g) \quad \text{O III}$$

$$\mathbf{B}_{27} = x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_7\right)a\,\hat{\mathbf{x}} + \tfrac{1}{4}a\,\hat{\mathbf{y}} + \tfrac{1}{4}a\,\hat{\mathbf{z}} \quad (24g) \quad \text{O III}$$

$$\mathbf{B}_{28} = x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_7\right)\mathbf{a}_2 + x_7\,\mathbf{a}_3 = \tfrac{1}{4}a\,\hat{\mathbf{x}} + x_7 a\,\hat{\mathbf{y}} + \tfrac{1}{4}a\,\hat{\mathbf{z}} \quad (24g) \quad \text{O III}$$

$$\mathbf{B}_{29} = \left(\tfrac{1}{2} - x_7\right)\mathbf{a}_1 + x_7\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_7\right)\mathbf{a}_3 = \tfrac{1}{4}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_7\right)a\,\hat{\mathbf{y}} + \tfrac{1}{4}a\,\hat{\mathbf{z}} \quad (24g) \quad \text{O III}$$

$$\mathbf{B}_{30} = x_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_7\right)\mathbf{a}_3 = \tfrac{1}{4}a\,\hat{\mathbf{x}} + \tfrac{1}{4}a\,\hat{\mathbf{y}} + x_7 a\,\hat{\mathbf{z}} \quad (24g) \quad \text{O III}$$

$$\mathbf{B}_{31} = \left(\tfrac{1}{2} - x_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_7\right)\mathbf{a}_2 + x_7\,\mathbf{a}_3 = \tfrac{1}{4}a\,\hat{\mathbf{x}} + \tfrac{1}{4}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_7\right)a\,\hat{\mathbf{z}} \quad (24g) \quad \text{O III}$$

$$\mathbf{B}_{32} = z_8\,\mathbf{a}_1 + z_8\,\mathbf{a}_2 + (2x_8 - z_8)\,\mathbf{a}_3 = x_8 a\,\hat{\mathbf{x}} + x_8 a\,\hat{\mathbf{y}} + z_8 a\,\hat{\mathbf{z}} \quad (48h) \quad \text{F II}$$

$$\mathbf{B}_{33} = z_8\,\mathbf{a}_1 + z_8\,\mathbf{a}_2 + (-2x_8 - z_8)\,\mathbf{a}_3 = -x_8 a\,\hat{\mathbf{x}} - x_8 a\,\hat{\mathbf{y}} + z_8 a\,\hat{\mathbf{z}} \quad (48h) \quad \text{F II}$$

$$\mathbf{B}_{34} = (2x_8 - z_8)\,\mathbf{a}_1 + (-2x_8 - z_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 = -x_8 a\,\hat{\mathbf{x}} + x_8 a\,\hat{\mathbf{y}} - z_8 a\,\hat{\mathbf{z}} \quad (48h) \quad \text{F II}$$

$$\mathbf{B}_{35} = (-2x_8 - z_8)\,\mathbf{a}_1 + (2x_8 - z_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 = x_8 a\,\hat{\mathbf{x}} - x_8 a\,\hat{\mathbf{y}} - z_8 a\,\hat{\mathbf{z}} \quad (48h) \quad \text{F II}$$

$$\mathbf{B}_{36} = (2x_8 - z_8)\,\mathbf{a}_1 + z_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 = z_8 a\,\hat{\mathbf{x}} + x_8 a\,\hat{\mathbf{y}} + x_8 a\,\hat{\mathbf{z}} \quad (48h) \quad \text{F II}$$

$$\mathbf{B}_{37} = (-2x_8 - z_8)\,\mathbf{a}_1 + z_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 = z_8 a\,\hat{\mathbf{x}} - x_8 a\,\hat{\mathbf{y}} - x_8 a\,\hat{\mathbf{z}} \quad (48h) \quad \text{F II}$$

$$\mathbf{B}_{38} = z_8\,\mathbf{a}_1 + (2x_8 - z_8)\,\mathbf{a}_2 + (-2x_8 - z_8)\,\mathbf{a}_3 = -z_8 a\,\hat{\mathbf{x}} - x_8 a\,\hat{\mathbf{y}} + x_8 a\,\hat{\mathbf{z}} \quad (48h) \quad \text{F II}$$

$$\mathbf{B}_{39} = z_8\,\mathbf{a}_1 + (-2x_8 - z_8)\,\mathbf{a}_2 + (2x_8 - z_8)\,\mathbf{a}_3 = -z_8 a\,\hat{\mathbf{x}} + x_8 a\,\hat{\mathbf{y}} - x_8 a\,\hat{\mathbf{z}} \quad (48h) \quad \text{F II}$$

$$\mathbf{B}_{40} = z_8\,\mathbf{a}_1 + (2x_8 - z_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 = x_8 a\,\hat{\mathbf{x}} + z_8 a\,\hat{\mathbf{y}} + x_8 a\,\hat{\mathbf{z}} \quad (48h) \quad \text{F II}$$

$$\mathbf{B}_{41} = z_8\,\mathbf{a}_1 + (-2x_8 - z_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 = -x_8 a\,\hat{\mathbf{x}} + z_8 a\,\hat{\mathbf{y}} - x_8 a\,\hat{\mathbf{z}} \quad (48h) \quad \text{F II}$$

$$\mathbf{B}_{42} = (-2x_8 - z_8)\,\mathbf{a}_1 + z_8\,\mathbf{a}_2 + (2x_8 - z_8)\,\mathbf{a}_3 = x_8 a\,\hat{\mathbf{x}} - z_8 a\,\hat{\mathbf{y}} - x_8 a\,\hat{\mathbf{z}} \quad (48h) \quad \text{F II}$$

$$\mathbf{B}_{43} = (2x_8 - z_8)\,\mathbf{a}_1 + z_8\,\mathbf{a}_2 + (-2x_8 - z_8)\,\mathbf{a}_3 = -x_8 a\,\hat{\mathbf{x}} - z_8 a\,\hat{\mathbf{y}} + x_8 a\,\hat{\mathbf{z}} \quad (48h) \quad \text{F II}$$

$$\mathbf{B}_{44} = z_9\,\mathbf{a}_1 + z_9\,\mathbf{a}_2 + (2x_9 - z_9)\,\mathbf{a}_3 = x_9 a\,\hat{\mathbf{x}} + x_9 a\,\hat{\mathbf{y}} + z_9 a\,\hat{\mathbf{z}} \quad (48h) \quad \text{O IV}$$

$$\mathbf{B}_{45} = z_9\,\mathbf{a}_1 + z_9\,\mathbf{a}_2 + (-2x_9 - z_9)\,\mathbf{a}_3 = -x_9 a\,\hat{\mathbf{x}} - x_9 a\,\hat{\mathbf{y}} + z_9 a\,\hat{\mathbf{z}} \quad (48h) \quad \text{O IV}$$

$$\mathbf{B}_{46} = (2x_9 - z_9)\,\mathbf{a}_1 + (-2x_9 - z_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 = -x_9 a\,\hat{\mathbf{x}} + x_9 a\,\hat{\mathbf{y}} - z_9 a\,\hat{\mathbf{z}} \quad (48h) \quad \text{O IV}$$

$$\mathbf{B}_{47} = (-2x_9 - z_9)\,\mathbf{a}_1 + (2x_9 - z_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 = x_9 a\,\hat{\mathbf{x}} - x_9 a\,\hat{\mathbf{y}} - z_9 a\,\hat{\mathbf{z}} \quad (48h) \quad \text{O IV}$$

$$\mathbf{B}_{48} = (2x_9 - z_9)\,\mathbf{a}_1 + z_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 = z_9 a\,\hat{\mathbf{x}} + x_9 a\,\hat{\mathbf{y}} + x_9 a\,\hat{\mathbf{z}} \quad (48h) \quad \text{O IV}$$

$$\mathbf{B}_{49} = (-2x_9 - z_9)\,\mathbf{a}_1 + z_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 = z_9 a\,\hat{\mathbf{x}} - x_9 a\,\hat{\mathbf{y}} - x_9 a\,\hat{\mathbf{z}} \quad (48h) \quad \text{O IV}$$

$$\mathbf{B}_{50} = z_9\,\mathbf{a}_1 + (2x_9 - z_9)\,\mathbf{a}_2 + (-2x_9 - z_9)\,\mathbf{a}_3 = -z_9 a\,\hat{\mathbf{x}} - x_9 a\,\hat{\mathbf{y}} + x_9 a\,\hat{\mathbf{z}} \quad (48h) \quad \text{O IV}$$

$$\mathbf{B}_{51} = z_9\,\mathbf{a}_1 + (-2x_9 - z_9)\,\mathbf{a}_2 + (2x_9 - z_9)\,\mathbf{a}_3 = -z_9 a\,\hat{\mathbf{x}} + x_9 a\,\hat{\mathbf{y}} - x_9 a\,\hat{\mathbf{z}} \quad (48h) \quad \text{O IV}$$

$$\mathbf{B}_{52} = z_9\,\mathbf{a}_1 + (2x_9 - z_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 = x_9 a\,\hat{\mathbf{x}} + z_9 a\,\hat{\mathbf{y}} + x_9 a\,\hat{\mathbf{z}} \quad (48h) \quad \text{O IV}$$

$$\mathbf{B}_{53} = z_9\,\mathbf{a}_1 + (-2x_9 - z_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 = -x_9 a\,\hat{\mathbf{x}} + z_9 a\,\hat{\mathbf{y}} - x_9 a\,\hat{\mathbf{z}} \quad (48h) \quad \text{O IV}$$

$$\mathbf{B}_{54} = (-2x_9 - z_9)\,\mathbf{a}_1 + z_9\,\mathbf{a}_2 + (2x_9 - z_9)\,\mathbf{a}_3 = x_9 a\,\hat{\mathbf{x}} - z_9 a\,\hat{\mathbf{y}} - x_9 a\,\hat{\mathbf{z}} \quad (48h) \quad \text{O IV}$$

$$\mathbf{B}_{55} = (2x_9 - z_9)\,\mathbf{a}_1 + z_9\,\mathbf{a}_2 + (-2x_9 - z_9)\,\mathbf{a}_3 = -x_9 a\,\hat{\mathbf{x}} - z_9 a\,\hat{\mathbf{y}} + x_9 a\,\hat{\mathbf{z}} \quad (48h) \quad \text{O IV}$$

$$\mathbf{B}_{56} = z_{10}\,\mathbf{a}_1 + z_{10}\,\mathbf{a}_2 + (2x_{10} - z_{10})\,\mathbf{a}_3 = x_{10} a\,\hat{\mathbf{x}} + x_{10} a\,\hat{\mathbf{y}} + z_{10} a\,\hat{\mathbf{z}} \quad (48h) \quad \text{O V}$$

$$\mathbf{B}_{57} = z_{10}\,\mathbf{a}_1 + z_{10}\,\mathbf{a}_2 + (-2x_{10} - z_{10})\,\mathbf{a}_3 = -x_{10} a\,\hat{\mathbf{x}} - x_{10} a\,\hat{\mathbf{y}} + z_{10} a\,\hat{\mathbf{z}} \quad (48h) \quad \text{O V}$$

$$\mathbf{B}_{58} = (2x_{10} - z_{10})\,\mathbf{a}_1 + (-2x_{10} - z_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 = -x_{10} a\,\hat{\mathbf{x}} + x_{10} a\,\hat{\mathbf{y}} - z_{10} a\,\hat{\mathbf{z}} \quad (48h) \quad \text{O V}$$

$$\mathbf{B}_{59} = (-2x_{10} - z_{10})\,\mathbf{a}_1 + (2x_{10} - z_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 = x_{10} a\,\hat{\mathbf{x}} - x_{10} a\,\hat{\mathbf{y}} - z_{10} a\,\hat{\mathbf{z}} \quad (48h) \quad \text{O V}$$

$\mathbf{B}_{60}$	=	$(2x_{10} - z_{10})\,\mathbf{a}_1 + z_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$z_{10}a\,\hat{\mathbf{x}} + x_{10}a\,\hat{\mathbf{y}} + x_{10}a\,\hat{\mathbf{z}}$	(48*h*)	O V
$\mathbf{B}_{61}$	=	$(-2x_{10} - z_{10})\,\mathbf{a}_1 + z_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$z_{10}a\,\hat{\mathbf{x}} - x_{10}a\,\hat{\mathbf{y}} - x_{10}a\,\hat{\mathbf{z}}$	(48*h*)	O V
$\mathbf{B}_{62}$	=	$z_{10}\,\mathbf{a}_1 + (2x_{10} - z_{10})\,\mathbf{a}_2 + (-2x_{10} - z_{10})\,\mathbf{a}_3$	=	$-z_{10}a\,\hat{\mathbf{x}} - x_{10}a\,\hat{\mathbf{y}} + x_{10}a\,\hat{\mathbf{z}}$	(48*h*)	O V
$\mathbf{B}_{63}$	=	$z_{10}\,\mathbf{a}_1 + (-2x_{10} - z_{10})\,\mathbf{a}_2 + (2x_{10} - z_{10})\,\mathbf{a}_3$	=	$-z_{10}a\,\hat{\mathbf{x}} + x_{10}a\,\hat{\mathbf{y}} - x_{10}a\,\hat{\mathbf{z}}$	(48*h*)	O V
$\mathbf{B}_{64}$	=	$z_{10}\,\mathbf{a}_1 + (2x_{10} - z_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$x_{10}a\,\hat{\mathbf{x}} + z_{10}a\,\hat{\mathbf{y}} + x_{10}a\,\hat{\mathbf{z}}$	(48*h*)	O V
$\mathbf{B}_{65}$	=	$z_{10}\,\mathbf{a}_1 + (-2x_{10} - z_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3$	=	$-x_{10}a\,\hat{\mathbf{x}} + z_{10}a\,\hat{\mathbf{y}} - x_{10}a\,\hat{\mathbf{z}}$	(48*h*)	O V
$\mathbf{B}_{66}$	=	$(-2x_{10} - z_{10})\,\mathbf{a}_1 + z_{10}\,\mathbf{a}_2 + (2x_{10} - z_{10})\,\mathbf{a}_3$	=	$x_{10}a\,\hat{\mathbf{x}} - z_{10}a\,\hat{\mathbf{y}} - x_{10}a\,\hat{\mathbf{z}}$	(48*h*)	O V
$\mathbf{B}_{67}$	=	$(2x_{10} - z_{10})\,\mathbf{a}_1 + z_{10}\,\mathbf{a}_2 + (-2x_{10} - z_{10})\,\mathbf{a}_3$	=	$-x_{10}a\,\hat{\mathbf{x}} - z_{10}a\,\hat{\mathbf{y}} + x_{10}a\,\hat{\mathbf{z}}$	(48*h*)	O V
$\mathbf{B}_{68}$	=	$z_{11}\,\mathbf{a}_1 + z_{11}\,\mathbf{a}_2 + (2x_{11} - z_{11})\,\mathbf{a}_3$	=	$x_{11}a\,\hat{\mathbf{x}} + x_{11}a\,\hat{\mathbf{y}} + z_{11}a\,\hat{\mathbf{z}}$	(48*h*)	Ti
$\mathbf{B}_{69}$	=	$z_{11}\,\mathbf{a}_1 + z_{11}\,\mathbf{a}_2 + (-2x_{11} - z_{11})\,\mathbf{a}_3$	=	$-x_{11}a\,\hat{\mathbf{x}} - x_{11}a\,\hat{\mathbf{y}} + z_{11}a\,\hat{\mathbf{z}}$	(48*h*)	Ti
$\mathbf{B}_{70}$	=	$(2x_{11} - z_{11})\,\mathbf{a}_1 + (-2x_{11} - z_{11})\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$-x_{11}a\,\hat{\mathbf{x}} + x_{11}a\,\hat{\mathbf{y}} - z_{11}a\,\hat{\mathbf{z}}$	(48*h*)	Ti
$\mathbf{B}_{71}$	=	$(-2x_{11} - z_{11})\,\mathbf{a}_1 + (2x_{11} - z_{11})\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$x_{11}a\,\hat{\mathbf{x}} - x_{11}a\,\hat{\mathbf{y}} - z_{11}a\,\hat{\mathbf{z}}$	(48*h*)	Ti
$\mathbf{B}_{72}$	=	$(2x_{11} - z_{11})\,\mathbf{a}_1 + z_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$z_{11}a\,\hat{\mathbf{x}} + x_{11}a\,\hat{\mathbf{y}} + x_{11}a\,\hat{\mathbf{z}}$	(48*h*)	Ti
$\mathbf{B}_{73}$	=	$(-2x_{11} - z_{11})\,\mathbf{a}_1 + z_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$z_{11}a\,\hat{\mathbf{x}} - x_{11}a\,\hat{\mathbf{y}} - x_{11}a\,\hat{\mathbf{z}}$	(48*h*)	Ti
$\mathbf{B}_{74}$	=	$z_{11}\,\mathbf{a}_1 + (2x_{11} - z_{11})\,\mathbf{a}_2 + (-2x_{11} - z_{11})\,\mathbf{a}_3$	=	$-z_{11}a\,\hat{\mathbf{x}} - x_{11}a\,\hat{\mathbf{y}} + x_{11}a\,\hat{\mathbf{z}}$	(48*h*)	Ti
$\mathbf{B}_{75}$	=	$z_{11}\,\mathbf{a}_1 + (-2x_{11} - z_{11})\,\mathbf{a}_2 + (2x_{11} - z_{11})\,\mathbf{a}_3$	=	$-z_{11}a\,\hat{\mathbf{x}} + x_{11}a\,\hat{\mathbf{y}} - x_{11}a\,\hat{\mathbf{z}}$	(48*h*)	Ti
$\mathbf{B}_{76}$	=	$z_{11}\,\mathbf{a}_1 + (2x_{11} - z_{11})\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$x_{11}a\,\hat{\mathbf{x}} + z_{11}a\,\hat{\mathbf{y}} + x_{11}a\,\hat{\mathbf{z}}$	(48*h*)	Ti
$\mathbf{B}_{77}$	=	$z_{11}\,\mathbf{a}_1 + (-2x_{11} - z_{11})\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3$	=	$-x_{11}a\,\hat{\mathbf{x}} + z_{11}a\,\hat{\mathbf{y}} - x_{11}a\,\hat{\mathbf{z}}$	(48*h*)	Ti
$\mathbf{B}_{78}$	=	$(-2x_{11} - z_{11})\,\mathbf{a}_1 + z_{11}\,\mathbf{a}_2 + (2x_{11} - z_{11})\,\mathbf{a}_3$	=	$x_{11}a\,\hat{\mathbf{x}} - z_{11}a\,\hat{\mathbf{y}} - x_{11}a\,\hat{\mathbf{z}}$	(48*h*)	Ti
$\mathbf{B}_{79}$	=	$(2x_{11} - z_{11})\,\mathbf{a}_1 + z_{11}\,\mathbf{a}_2 + (-2x_{11} - z_{11})\,\mathbf{a}_3$	=	$-x_{11}a\,\hat{\mathbf{x}} - z_{11}a\,\hat{\mathbf{y}} + x_{11}a\,\hat{\mathbf{z}}$	(48*h*)	Ti

References:

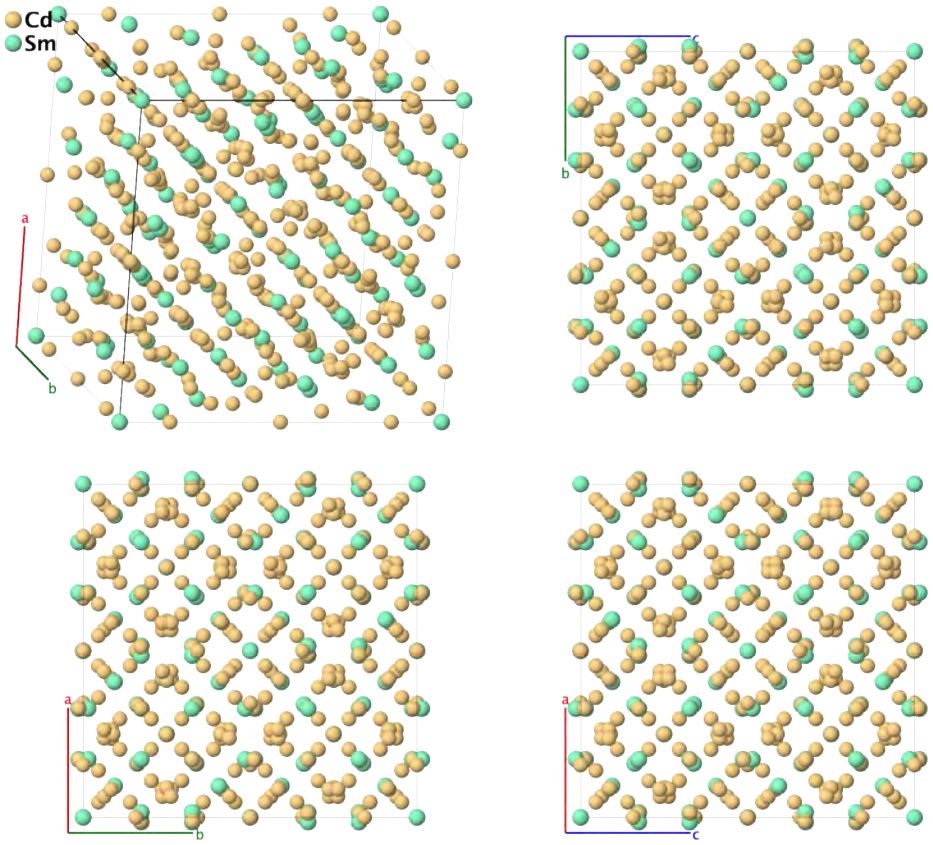
- T. S. Ercit and F. C. Hawthorne, *Murataite, A UB_{12} derivative structure with condensed Keggin molecules*, Can. Mineral. **33**, 1233–1229 (1995).
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1791
- POSCAR: pp. 1792

$Sm_{11}Cd_{45}$ Structure: A45B11_cF448_216_bd4efg5h_ac2eh

http://aflow.org/prototype-encyclopedia/A45B11_cF448_216_bd4efg5h_ac2eh



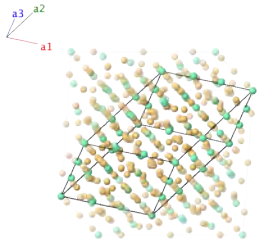
Prototype	:	$Cd_{45}Sm_{11}$
AFLOW prototype label	:	A45B11_cF448_216_bd4efg5h_ac2eh
***Strukturbericht* designation**	:	None
Pearson symbol	:	cF448
Space group number	:	216
Space group symbol	:	$F\bar{4}3m$
AFLOW prototype command	:	`aflow --proto=A45B11_cF448_216_bd4efg5h_ac2eh --params=`$a, x_5, x_6, x_7, x_8, x_9, x_{10}, x_{11}, x_{12}, x_{13}, z_{13}, x_{14}, z_{14}, x_{15}, z_{15}, x_{16}, z_{16}, x_{17}, z_{17}, x_{18}, z_{18}$

Other compounds with this structure

- $Dy_{11}Cd_{45}$, $Er_{11}Cd_{45}$, $Gd_{11}Cd_{45}$, $Ho_{11}Cd_{45}$, $Lu_{11}Cd_{45}$, $Nd_{11}Cd_{45}$, $Tb_{11}Cd_{45}$, $Tm_{11}Cd_{45}$, $Pr_{11}Cd_{45}$, $Y_{11}Cd_{45}$, $Ce_{11}Hg_{45}$, $Nd_{11}Hg_{45}$, $Pr_{11}Hg_{45}$, and $Sm_{11}Hg_{45}$

Face-centered Cubic primitive vectors:

$$\mathbf{a}_1 = \tfrac{1}{2}\,a\,\hat{\mathbf{y}} + \tfrac{1}{2}\,a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{y}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(4*a*)	Sm I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(4*b*)	Cd I
$\mathbf{B}_3$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(4*c*)	Sm II
$\mathbf{B}_4$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(4*d*)	Cd II
$\mathbf{B}_5$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}} + x_5a\,\hat{\mathbf{y}} + x_5a\,\hat{\mathbf{z}}$	(16*e*)	Cd III
$\mathbf{B}_6$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 - 3x_5\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}} - x_5a\,\hat{\mathbf{y}} + x_5a\,\hat{\mathbf{z}}$	(16*e*)	Cd III
$\mathbf{B}_7$	=	$x_5\,\mathbf{a}_1 - 3x_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}} + x_5a\,\hat{\mathbf{y}} - x_5a\,\hat{\mathbf{z}}$	(16*e*)	Cd III
$\mathbf{B}_8$	=	$-3x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}} - x_5a\,\hat{\mathbf{y}} - x_5a\,\hat{\mathbf{z}}$	(16*e*)	Cd III
$\mathbf{B}_9$	=	$x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + x_6\,\mathbf{a}_3$	=	$x_6a\,\hat{\mathbf{x}} + x_6a\,\hat{\mathbf{y}} + x_6a\,\hat{\mathbf{z}}$	(16*e*)	Cd IV
$\mathbf{B}_{10}$	=	$x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 - 3x_6\,\mathbf{a}_3$	=	$-x_6a\,\hat{\mathbf{x}} - x_6a\,\hat{\mathbf{y}} + x_6a\,\hat{\mathbf{z}}$	(16*e*)	Cd IV
$\mathbf{B}_{11}$	=	$x_6\,\mathbf{a}_1 - 3x_6\,\mathbf{a}_2 + x_6\,\mathbf{a}_3$	=	$-x_6a\,\hat{\mathbf{x}} + x_6a\,\hat{\mathbf{y}} - x_6a\,\hat{\mathbf{z}}$	(16*e*)	Cd IV
$\mathbf{B}_{12}$	=	$-3x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + x_6\,\mathbf{a}_3$	=	$x_6a\,\hat{\mathbf{x}} - x_6a\,\hat{\mathbf{y}} - x_6a\,\hat{\mathbf{z}}$	(16*e*)	Cd IV
$\mathbf{B}_{13}$	=	$x_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + x_7\,\mathbf{a}_3$	=	$x_7a\,\hat{\mathbf{x}} + x_7a\,\hat{\mathbf{y}} + x_7a\,\hat{\mathbf{z}}$	(16*e*)	Cd V
$\mathbf{B}_{14}$	=	$x_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 - 3x_7\,\mathbf{a}_3$	=	$-x_7a\,\hat{\mathbf{x}} - x_7a\,\hat{\mathbf{y}} + x_7a\,\hat{\mathbf{z}}$	(16*e*)	Cd V
$\mathbf{B}_{15}$	=	$x_7\,\mathbf{a}_1 - 3x_7\,\mathbf{a}_2 + x_7\,\mathbf{a}_3$	=	$-x_7a\,\hat{\mathbf{x}} + x_7a\,\hat{\mathbf{y}} - x_7a\,\hat{\mathbf{z}}$	(16*e*)	Cd V
$\mathbf{B}_{16}$	=	$-3x_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + x_7\,\mathbf{a}_3$	=	$x_7a\,\hat{\mathbf{x}} - x_7a\,\hat{\mathbf{y}} - x_7a\,\hat{\mathbf{z}}$	(16*e*)	Cd V
$\mathbf{B}_{17}$	=	$x_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 + x_8\,\mathbf{a}_3$	=	$x_8a\,\hat{\mathbf{x}} + x_8a\,\hat{\mathbf{y}} + x_8a\,\hat{\mathbf{z}}$	(16*e*)	Cd VI
$\mathbf{B}_{18}$	=	$x_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 - 3x_8\,\mathbf{a}_3$	=	$-x_8a\,\hat{\mathbf{x}} - x_8a\,\hat{\mathbf{y}} + x_8a\,\hat{\mathbf{z}}$	(16*e*)	Cd VI
$\mathbf{B}_{19}$	=	$x_8\,\mathbf{a}_1 - 3x_8\,\mathbf{a}_2 + x_8\,\mathbf{a}_3$	=	$-x_8a\,\hat{\mathbf{x}} + x_8a\,\hat{\mathbf{y}} - x_8a\,\hat{\mathbf{z}}$	(16*e*)	Cd VI
$\mathbf{B}_{20}$	=	$-3x_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 + x_8\,\mathbf{a}_3$	=	$x_8a\,\hat{\mathbf{x}} - x_8a\,\hat{\mathbf{y}} - x_8a\,\hat{\mathbf{z}}$	(16*e*)	Cd VI
$\mathbf{B}_{21}$	=	$x_9\,\mathbf{a}_1 + x_9\,\mathbf{a}_2 + x_9\,\mathbf{a}_3$	=	$x_9a\,\hat{\mathbf{x}} + x_9a\,\hat{\mathbf{y}} + x_9a\,\hat{\mathbf{z}}$	(16*e*)	Sm III
$\mathbf{B}_{22}$	=	$x_9\,\mathbf{a}_1 + x_9\,\mathbf{a}_2 - 3x_9\,\mathbf{a}_3$	=	$-x_9a\,\hat{\mathbf{x}} - x_9a\,\hat{\mathbf{y}} + x_9a\,\hat{\mathbf{z}}$	(16*e*)	Sm III
$\mathbf{B}_{23}$	=	$x_9\,\mathbf{a}_1 - 3x_9\,\mathbf{a}_2 + x_9\,\mathbf{a}_3$	=	$-x_9a\,\hat{\mathbf{x}} + x_9a\,\hat{\mathbf{y}} - x_9a\,\hat{\mathbf{z}}$	(16*e*)	Sm III
$\mathbf{B}_{24}$	=	$-3x_9\,\mathbf{a}_1 + x_9\,\mathbf{a}_2 + x_9\,\mathbf{a}_3$	=	$x_9a\,\hat{\mathbf{x}} - x_9a\,\hat{\mathbf{y}} - x_9a\,\hat{\mathbf{z}}$	(16*e*)	Sm III
$\mathbf{B}_{25}$	=	$x_{10}\,\mathbf{a}_1 + x_{10}\,\mathbf{a}_2 + x_{10}\,\mathbf{a}_3$	=	$x_{10}a\,\hat{\mathbf{x}} + x_{10}a\,\hat{\mathbf{y}} + x_{10}a\,\hat{\mathbf{z}}$	(16*e*)	Sm IV
$\mathbf{B}_{26}$	=	$x_{10}\,\mathbf{a}_1 + x_{10}\,\mathbf{a}_2 - 3x_{10}\,\mathbf{a}_3$	=	$-x_{10}a\,\hat{\mathbf{x}} - x_{10}a\,\hat{\mathbf{y}} + x_{10}a\,\hat{\mathbf{z}}$	(16*e*)	Sm IV

$\mathbf{B}_{27}$	=	$x_{10}\,\mathbf{a}_1 - 3x_{10}\,\mathbf{a}_2 + x_{10}\,\mathbf{a}_3$	=	$-x_{10}a\,\hat{\mathbf{x}} + x_{10}a\,\hat{\mathbf{y}} - x_{10}a\,\hat{\mathbf{z}}$	(16*e*)	Sm IV
$\mathbf{B}_{28}$	=	$-3x_{10}\,\mathbf{a}_1 + x_{10}\,\mathbf{a}_2 + x_{10}\,\mathbf{a}_3$	=	$x_{10}a\,\hat{\mathbf{x}} - x_{10}a\,\hat{\mathbf{y}} - x_{10}a\,\hat{\mathbf{z}}$	(16*e*)	Sm IV
$\mathbf{B}_{29}$	=	$-x_{11}\,\mathbf{a}_1 + x_{11}\,\mathbf{a}_2 + x_{11}\,\mathbf{a}_3$	=	$x_{11}a\,\hat{\mathbf{x}}$	(24*f*)	Cd VII
$\mathbf{B}_{30}$	=	$x_{11}\,\mathbf{a}_1 - x_{11}\,\mathbf{a}_2 - x_{11}\,\mathbf{a}_3$	=	$-x_{11}a\,\hat{\mathbf{x}}$	(24*f*)	Cd VII
$\mathbf{B}_{31}$	=	$x_{11}\,\mathbf{a}_1 - x_{11}\,\mathbf{a}_2 + x_{11}\,\mathbf{a}_3$	=	$x_{11}a\,\hat{\mathbf{y}}$	(24*f*)	Cd VII
$\mathbf{B}_{32}$	=	$-x_{11}\,\mathbf{a}_1 + x_{11}\,\mathbf{a}_2 - x_{11}\,\mathbf{a}_3$	=	$-x_{11}a\,\hat{\mathbf{y}}$	(24*f*)	Cd VII
$\mathbf{B}_{33}$	=	$x_{11}\,\mathbf{a}_1 + x_{11}\,\mathbf{a}_2 - x_{11}\,\mathbf{a}_3$	=	$x_{11}a\,\hat{\mathbf{z}}$	(24*f*)	Cd VII
$\mathbf{B}_{34}$	=	$-x_{11}\,\mathbf{a}_1 - x_{11}\,\mathbf{a}_2 + x_{11}\,\mathbf{a}_3$	=	$-x_{11}a\,\hat{\mathbf{z}}$	(24*f*)	Cd VII
$\mathbf{B}_{35}$	=	$\left(\frac{1}{2} - x_{12}\right)\,\mathbf{a}_1 + x_{12}\,\mathbf{a}_2 + x_{12}\,\mathbf{a}_3$	=	$x_{12}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(24*g*)	Cd VIII
$\mathbf{B}_{36}$	=	$x_{12}\,\mathbf{a}_1 + \left(\frac{1}{2} - x_{12}\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - x_{12}\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_{12}\right)a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(24*g*)	Cd VIII
$\mathbf{B}_{37}$	=	$x_{12}\,\mathbf{a}_1 + \left(\frac{1}{2} - x_{12}\right)\,\mathbf{a}_2 + x_{12}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + x_{12}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(24*g*)	Cd VIII
$\mathbf{B}_{38}$	=	$\left(\frac{1}{2} - x_{12}\right)\,\mathbf{a}_1 + x_{12}\,\mathbf{a}_2 + \left(\frac{1}{2} - x_{12}\right)\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_{12}\right)a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(24*g*)	Cd VIII
$\mathbf{B}_{39}$	=	$x_{12}\,\mathbf{a}_1 + x_{12}\,\mathbf{a}_2 + \left(\frac{1}{2} - x_{12}\right)\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + x_{12}a\,\hat{\mathbf{z}}$	(24*g*)	Cd VIII
$\mathbf{B}_{40}$	=	$\left(\frac{1}{2} - x_{12}\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - x_{12}\right)\,\mathbf{a}_2 + x_{12}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_{12}\right)a\,\hat{\mathbf{z}}$	(24*g*)	Cd VIII
$\mathbf{B}_{41}$	=	$z_{13}\,\mathbf{a}_1 + z_{13}\,\mathbf{a}_2 + (2x_{13} - z_{13})\,\mathbf{a}_3$	=	$x_{13}a\,\hat{\mathbf{x}} + x_{13}a\,\hat{\mathbf{y}} + z_{13}a\,\hat{\mathbf{z}}$	(48*h*)	Cd IX
$\mathbf{B}_{42}$	=	$z_{13}\,\mathbf{a}_1 + z_{13}\,\mathbf{a}_2 + (-2x_{13} - z_{13})\,\mathbf{a}_3$	=	$-x_{13}a\,\hat{\mathbf{x}} - x_{13}a\,\hat{\mathbf{y}} + z_{13}a\,\hat{\mathbf{z}}$	(48*h*)	Cd IX
$\mathbf{B}_{43}$	=	$(2x_{13} - z_{13})\,\mathbf{a}_1 + (-2x_{13} - z_{13})\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3$	=	$-x_{13}a\,\hat{\mathbf{x}} + x_{13}a\,\hat{\mathbf{y}} - z_{13}a\,\hat{\mathbf{z}}$	(48*h*)	Cd IX
$\mathbf{B}_{44}$	=	$(-2x_{13} - z_{13})\,\mathbf{a}_1 + (2x_{13} - z_{13})\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3$	=	$x_{13}a\,\hat{\mathbf{x}} - x_{13}a\,\hat{\mathbf{y}} - z_{13}a\,\hat{\mathbf{z}}$	(48*h*)	Cd IX
$\mathbf{B}_{45}$	=	$(2x_{13} - z_{13})\,\mathbf{a}_1 + z_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3$	=	$z_{13}a\,\hat{\mathbf{x}} + x_{13}a\,\hat{\mathbf{y}} + x_{13}a\,\hat{\mathbf{z}}$	(48*h*)	Cd IX
$\mathbf{B}_{46}$	=	$(-2x_{13} - z_{13})\,\mathbf{a}_1 + z_{13}\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3$	=	$z_{13}a\,\hat{\mathbf{x}} - x_{13}a\,\hat{\mathbf{y}} - x_{13}a\,\hat{\mathbf{z}}$	(48*h*)	Cd IX
$\mathbf{B}_{47}$	=	$z_{13}\,\mathbf{a}_1 + (2x_{13} - z_{13})\,\mathbf{a}_2 + (-2x_{13} - z_{13})\,\mathbf{a}_3$	=	$-z_{13}a\,\hat{\mathbf{x}} - x_{13}a\,\hat{\mathbf{y}} + x_{13}a\,\hat{\mathbf{z}}$	(48*h*)	Cd IX
$\mathbf{B}_{48}$	=	$z_{13}\,\mathbf{a}_1 + (-2x_{13} - z_{13})\,\mathbf{a}_2 + (2x_{13} - z_{13})\,\mathbf{a}_3$	=	$-z_{13}a\,\hat{\mathbf{x}} + x_{13}a\,\hat{\mathbf{y}} - x_{13}a\,\hat{\mathbf{z}}$	(48*h*)	Cd IX
$\mathbf{B}_{49}$	=	$z_{13}\,\mathbf{a}_1 + (2x_{13} - z_{13})\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3$	=	$x_{13}a\,\hat{\mathbf{x}} + z_{13}a\,\hat{\mathbf{y}} + x_{13}a\,\hat{\mathbf{z}}$	(48*h*)	Cd IX
$\mathbf{B}_{50}$	=	$z_{13}\,\mathbf{a}_1 + (-2x_{13} - z_{13})\,\mathbf{a}_2 + z_{13}\,\mathbf{a}_3$	=	$-x_{13}a\,\hat{\mathbf{x}} + z_{13}a\,\hat{\mathbf{y}} - x_{13}a\,\hat{\mathbf{z}}$	(48*h*)	Cd IX
$\mathbf{B}_{51}$	=	$(-2x_{13} - z_{13})\,\mathbf{a}_1 + z_{13}\,\mathbf{a}_2 + (2x_{13} - z_{13})\,\mathbf{a}_3$	=	$x_{13}a\,\hat{\mathbf{x}} - z_{13}a\,\hat{\mathbf{y}} - x_{13}a\,\hat{\mathbf{z}}$	(48*h*)	Cd IX
$\mathbf{B}_{52}$	=	$(2x_{13} - z_{13})\,\mathbf{a}_1 + z_{13}\,\mathbf{a}_2 + (-2x_{13} - z_{13})\,\mathbf{a}_3$	=	$-x_{13}a\,\hat{\mathbf{x}} - z_{13}a\,\hat{\mathbf{y}} + x_{13}a\,\hat{\mathbf{z}}$	(48*h*)	Cd IX
$\mathbf{B}_{53}$	=	$z_{14}\,\mathbf{a}_1 + z_{14}\,\mathbf{a}_2 + (2x_{14} - z_{14})\,\mathbf{a}_3$	=	$x_{14}a\,\hat{\mathbf{x}} + x_{14}a\,\hat{\mathbf{y}} + z_{14}a\,\hat{\mathbf{z}}$	(48*h*)	Cd X
$\mathbf{B}_{54}$	=	$z_{14}\,\mathbf{a}_1 + z_{14}\,\mathbf{a}_2 + (-2x_{14} - z_{14})\,\mathbf{a}_3$	=	$-x_{14}a\,\hat{\mathbf{x}} - x_{14}a\,\hat{\mathbf{y}} + z_{14}a\,\hat{\mathbf{z}}$	(48*h*)	Cd X
$\mathbf{B}_{55}$	=	$(2x_{14} - z_{14})\,\mathbf{a}_1 + (-2x_{14} - z_{14})\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3$	=	$-x_{14}a\,\hat{\mathbf{x}} + x_{14}a\,\hat{\mathbf{y}} - z_{14}a\,\hat{\mathbf{z}}$	(48*h*)	Cd X
$\mathbf{B}_{56}$	=	$(-2x_{14} - z_{14})\,\mathbf{a}_1 + (2x_{14} - z_{14})\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3$	=	$x_{14}a\,\hat{\mathbf{x}} - x_{14}a\,\hat{\mathbf{y}} - z_{14}a\,\hat{\mathbf{z}}$	(48*h*)	Cd X
$\mathbf{B}_{57}$	=	$(2x_{14} - z_{14})\,\mathbf{a}_1 + z_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3$	=	$z_{14}a\,\hat{\mathbf{x}} + x_{14}a\,\hat{\mathbf{y}} + x_{14}a\,\hat{\mathbf{z}}$	(48*h*)	Cd X
$\mathbf{B}_{58}$	=	$(-2x_{14} - z_{14})\,\mathbf{a}_1 + z_{14}\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3$	=	$z_{14}a\,\hat{\mathbf{x}} - x_{14}a\,\hat{\mathbf{y}} - x_{14}a\,\hat{\mathbf{z}}$	(48*h*)	Cd X
$\mathbf{B}_{59}$	=	$z_{14}\,\mathbf{a}_1 + (2x_{14} - z_{14})\,\mathbf{a}_2 + (-2x_{14} - z_{14})\,\mathbf{a}_3$	=	$-z_{14}a\,\hat{\mathbf{x}} - x_{14}a\,\hat{\mathbf{y}} + x_{14}a\,\hat{\mathbf{z}}$	(48*h*)	Cd X
$\mathbf{B}_{60}$	=	$z_{14}\,\mathbf{a}_1 + (-2x_{14} - z_{14})\,\mathbf{a}_2 + (2x_{14} - z_{14})\,\mathbf{a}_3$	=	$-z_{14}a\,\hat{\mathbf{x}} + x_{14}a\,\hat{\mathbf{y}} - x_{14}a\,\hat{\mathbf{z}}$	(48*h*)	Cd X
$\mathbf{B}_{61}$	=	$z_{14}\,\mathbf{a}_1 + (2x_{14} - z_{14})\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3$	=	$x_{14}a\,\hat{\mathbf{x}} + z_{14}a\,\hat{\mathbf{y}} + x_{14}a\,\hat{\mathbf{z}}$	(48*h*)	Cd X
$\mathbf{B}_{62}$	=	$z_{14}\,\mathbf{a}_1 + (-2x_{14} - z_{14})\,\mathbf{a}_2 + z_{14}\,\mathbf{a}_3$	=	$-x_{14}a\,\hat{\mathbf{x}} + z_{14}a\,\hat{\mathbf{y}} - x_{14}a\,\hat{\mathbf{z}}$	(48*h*)	Cd X

$$\mathbf{B}_{63} = (-2x_{14} - z_{14})\,\mathbf{a}_1 + z_{14}\,\mathbf{a}_2 + (2x_{14} - z_{14})\,\mathbf{a}_3 = x_{14}a\,\hat{\mathbf{x}} - z_{14}a\,\hat{\mathbf{y}} - x_{14}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd X}$$

$$\mathbf{B}_{64} = (2x_{14} - z_{14})\,\mathbf{a}_1 + z_{14}\,\mathbf{a}_2 + (-2x_{14} - z_{14})\,\mathbf{a}_3 = -x_{14}a\,\hat{\mathbf{x}} - z_{14}a\,\hat{\mathbf{y}} + x_{14}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd X}$$

$$\mathbf{B}_{65} = z_{15}\,\mathbf{a}_1 + z_{15}\,\mathbf{a}_2 + (2x_{15} - z_{15})\,\mathbf{a}_3 = x_{15}a\,\hat{\mathbf{x}} + x_{15}a\,\hat{\mathbf{y}} + z_{15}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd XI}$$

$$\mathbf{B}_{66} = z_{15}\,\mathbf{a}_1 + z_{15}\,\mathbf{a}_2 + (-2x_{15} - z_{15})\,\mathbf{a}_3 = -x_{15}a\,\hat{\mathbf{x}} - x_{15}a\,\hat{\mathbf{y}} + z_{15}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd XI}$$

$$\mathbf{B}_{67} = (2x_{15} - z_{15})\,\mathbf{a}_1 + (-2x_{15} - z_{15})\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 = -x_{15}a\,\hat{\mathbf{x}} + x_{15}a\,\hat{\mathbf{y}} - z_{15}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd XI}$$

$$\mathbf{B}_{68} = (-2x_{15} - z_{15})\,\mathbf{a}_1 + (2x_{15} - z_{15})\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 = x_{15}a\,\hat{\mathbf{x}} - x_{15}a\,\hat{\mathbf{y}} - z_{15}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd XI}$$

$$\mathbf{B}_{69} = (2x_{15} - z_{15})\,\mathbf{a}_1 + z_{15}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 = z_{15}a\,\hat{\mathbf{x}} + x_{15}a\,\hat{\mathbf{y}} + x_{15}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd XI}$$

$$\mathbf{B}_{70} = (-2x_{15} - z_{15})\,\mathbf{a}_1 + z_{15}\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 = z_{15}a\,\hat{\mathbf{x}} - x_{15}a\,\hat{\mathbf{y}} - x_{15}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd XI}$$

$$\mathbf{B}_{71} = z_{15}\,\mathbf{a}_1 + (2x_{15} - z_{15})\,\mathbf{a}_2 + (-2x_{15} - z_{15})\,\mathbf{a}_3 = -z_{15}a\,\hat{\mathbf{x}} - x_{15}a\,\hat{\mathbf{y}} + x_{15}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd XI}$$

$$\mathbf{B}_{72} = z_{15}\,\mathbf{a}_1 + (-2x_{15} - z_{15})\,\mathbf{a}_2 + (2x_{15} - z_{15})\,\mathbf{a}_3 = -z_{15}a\,\hat{\mathbf{x}} + x_{15}a\,\hat{\mathbf{y}} - x_{15}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd XI}$$

$$\mathbf{B}_{73} = z_{15}\,\mathbf{a}_1 + (2x_{15} - z_{15})\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 = x_{15}a\,\hat{\mathbf{x}} + z_{15}a\,\hat{\mathbf{y}} + x_{15}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd XI}$$

$$\mathbf{B}_{74} = z_{15}\,\mathbf{a}_1 + (-2x_{15} - z_{15})\,\mathbf{a}_2 + z_{15}\,\mathbf{a}_3 = -x_{15}a\,\hat{\mathbf{x}} + z_{15}a\,\hat{\mathbf{y}} - x_{15}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd XI}$$

$$\mathbf{B}_{75} = (-2x_{15} - z_{15})\,\mathbf{a}_1 + z_{15}\,\mathbf{a}_2 + (2x_{15} - z_{15})\,\mathbf{a}_3 = x_{15}a\,\hat{\mathbf{x}} - z_{15}a\,\hat{\mathbf{y}} - x_{15}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd XI}$$

$$\mathbf{B}_{76} = (2x_{15} - z_{15})\,\mathbf{a}_1 + z_{15}\,\mathbf{a}_2 + (-2x_{15} - z_{15})\,\mathbf{a}_3 = -x_{15}a\,\hat{\mathbf{x}} - z_{15}a\,\hat{\mathbf{y}} + x_{15}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd XI}$$

$$\mathbf{B}_{77} = z_{16}\,\mathbf{a}_1 + z_{16}\,\mathbf{a}_2 + (2x_{16} - z_{16})\,\mathbf{a}_3 = x_{16}a\,\hat{\mathbf{x}} + x_{16}a\,\hat{\mathbf{y}} + z_{16}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd XII}$$

$$\mathbf{B}_{78} = z_{16}\,\mathbf{a}_1 + z_{16}\,\mathbf{a}_2 + (-2x_{16} - z_{16})\,\mathbf{a}_3 = -x_{16}a\,\hat{\mathbf{x}} - x_{16}a\,\hat{\mathbf{y}} + z_{16}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd XII}$$

$$\mathbf{B}_{79} = (2x_{16} - z_{16})\,\mathbf{a}_1 + (-2x_{16} - z_{16})\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 = -x_{16}a\,\hat{\mathbf{x}} + x_{16}a\,\hat{\mathbf{y}} - z_{16}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd XII}$$

$$\mathbf{B}_{80} = (-2x_{16} - z_{16})\,\mathbf{a}_1 + (2x_{16} - z_{16})\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 = x_{16}a\,\hat{\mathbf{x}} - x_{16}a\,\hat{\mathbf{y}} - z_{16}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd XII}$$

$$\mathbf{B}_{81} = (2x_{16} - z_{16})\,\mathbf{a}_1 + z_{16}\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 = z_{16}a\,\hat{\mathbf{x}} + x_{16}a\,\hat{\mathbf{y}} + x_{16}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd XII}$$

$$\mathbf{B}_{82} = (-2x_{16} - z_{16})\,\mathbf{a}_1 + z_{16}\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 = z_{16}a\,\hat{\mathbf{x}} - x_{16}a\,\hat{\mathbf{y}} - x_{16}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd XII}$$

$$\mathbf{B}_{83} = z_{16}\,\mathbf{a}_1 + (2x_{16} - z_{16})\,\mathbf{a}_2 + (-2x_{16} - z_{16})\,\mathbf{a}_3 = -z_{16}a\,\hat{\mathbf{x}} - x_{16}a\,\hat{\mathbf{y}} + x_{16}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd XII}$$

$$\mathbf{B}_{84} = z_{16}\,\mathbf{a}_1 + (-2x_{16} - z_{16})\,\mathbf{a}_2 + (2x_{16} - z_{16})\,\mathbf{a}_3 = -z_{16}a\,\hat{\mathbf{x}} + x_{16}a\,\hat{\mathbf{y}} - x_{16}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd XII}$$

$$\mathbf{B}_{85} = z_{16}\,\mathbf{a}_1 + (2x_{16} - z_{16})\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 = x_{16}a\,\hat{\mathbf{x}} + z_{16}a\,\hat{\mathbf{y}} + x_{16}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd XII}$$

$$\mathbf{B}_{86} = z_{16}\,\mathbf{a}_1 + (-2x_{16} - z_{16})\,\mathbf{a}_2 + z_{16}\,\mathbf{a}_3 = -x_{16}a\,\hat{\mathbf{x}} + z_{16}a\,\hat{\mathbf{y}} - x_{16}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd XII}$$

$$\mathbf{B}_{87} = (-2x_{16} - z_{16})\,\mathbf{a}_1 + z_{16}\,\mathbf{a}_2 + (2x_{16} - z_{16})\,\mathbf{a}_3 = x_{16}a\,\hat{\mathbf{x}} - z_{16}a\,\hat{\mathbf{y}} - x_{16}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd XII}$$

$$\mathbf{B}_{88} = (2x_{16} - z_{16})\,\mathbf{a}_1 + z_{16}\,\mathbf{a}_2 + (-2x_{16} - z_{16})\,\mathbf{a}_3 = -x_{16}a\,\hat{\mathbf{x}} - z_{16}a\,\hat{\mathbf{y}} + x_{16}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd XII}$$

$$\mathbf{B}_{89} = z_{17}\,\mathbf{a}_1 + z_{17}\,\mathbf{a}_2 + (2x_{17} - z_{17})\,\mathbf{a}_3 = x_{17}a\,\hat{\mathbf{x}} + x_{17}a\,\hat{\mathbf{y}} + z_{17}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd XIII}$$

$$\mathbf{B}_{90} = z_{17}\,\mathbf{a}_1 + z_{17}\,\mathbf{a}_2 + (-2x_{17} - z_{17})\,\mathbf{a}_3 = -x_{17}a\,\hat{\mathbf{x}} - x_{17}a\,\hat{\mathbf{y}} + z_{17}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd XIII}$$

$$\mathbf{B}_{91} = (2x_{17} - z_{17})\,\mathbf{a}_1 + (-2x_{17} - z_{17})\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 = -x_{17}a\,\hat{\mathbf{x}} + x_{17}a\,\hat{\mathbf{y}} - z_{17}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd XIII}$$

$$\mathbf{B}_{92} = (-2x_{17} - z_{17})\,\mathbf{a}_1 + (2x_{17} - z_{17})\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 = x_{17}a\,\hat{\mathbf{x}} - x_{17}a\,\hat{\mathbf{y}} - z_{17}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd XIII}$$

$$\mathbf{B}_{93} = (2x_{17} - z_{17})\,\mathbf{a}_1 + z_{17}\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 = z_{17}a\,\hat{\mathbf{x}} + x_{17}a\,\hat{\mathbf{y}} + x_{17}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd XIII}$$

$$\mathbf{B}_{94} = (-2x_{17} - z_{17})\,\mathbf{a}_1 + z_{17}\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 = z_{17}a\,\hat{\mathbf{x}} - x_{17}a\,\hat{\mathbf{y}} - x_{17}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd XIII}$$

$$\mathbf{B}_{95} = z_{17}\,\mathbf{a}_1 + (2x_{17} - z_{17})\,\mathbf{a}_2 + (-2x_{17} - z_{17})\,\mathbf{a}_3 = -z_{17}a\,\hat{\mathbf{x}} - x_{17}a\,\hat{\mathbf{y}} + x_{17}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd XIII}$$

$$\mathbf{B}_{96} = z_{17}\,\mathbf{a}_1 + (-2x_{17} - z_{17})\,\mathbf{a}_2 + (2x_{17} - z_{17})\,\mathbf{a}_3 = -z_{17}a\,\hat{\mathbf{x}} + x_{17}a\,\hat{\mathbf{y}} - x_{17}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd XIII}$$

$$\mathbf{B}_{97} = z_{17}\,\mathbf{a}_1 + (2x_{17} - z_{17})\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 = x_{17}a\,\hat{\mathbf{x}} + z_{17}a\,\hat{\mathbf{y}} + x_{17}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd XIII}$$

$$\mathbf{B}_{98} = z_{17}\,\mathbf{a}_1 + (-2x_{17} - z_{17})\,\mathbf{a}_2 + z_{17}\,\mathbf{a}_3 = -x_{17}a\,\hat{\mathbf{x}} + z_{17}a\,\hat{\mathbf{y}} - x_{17}a\,\hat{\mathbf{z}} \quad (48h) \quad \text{Cd XIII}$$

$\mathbf{B}_{99}$	=	$(-2x_{17}-z_{17})\,\mathbf{a}_1+z_{17}\,\mathbf{a}_2+(2x_{17}-z_{17})\,\mathbf{a}_3$	=	$x_{17}a\,\hat{\mathbf{x}}-z_{17}a\,\hat{\mathbf{y}}-x_{17}a\,\hat{\mathbf{z}}$	(48*h*)	Cd XIII
$\mathbf{B}_{100}$	=	$(2x_{17}-z_{17})\,\mathbf{a}_1+z_{17}\,\mathbf{a}_2+(-2x_{17}-z_{17})\,\mathbf{a}_3$	=	$-x_{17}a\,\hat{\mathbf{x}}-z_{17}a\,\hat{\mathbf{y}}+x_{17}a\,\hat{\mathbf{z}}$	(48*h*)	Cd XIII
$\mathbf{B}_{101}$	=	$z_{18}\,\mathbf{a}_1+z_{18}\,\mathbf{a}_2+(2x_{18}-z_{18})\,\mathbf{a}_3$	=	$x_{18}a\,\hat{\mathbf{x}}+x_{18}a\,\hat{\mathbf{y}}+z_{18}a\,\hat{\mathbf{z}}$	(48*h*)	Sm V
$\mathbf{B}_{102}$	=	$z_{18}\,\mathbf{a}_1+z_{18}\,\mathbf{a}_2+(-2x_{18}-z_{18})\,\mathbf{a}_3$	=	$-x_{18}a\,\hat{\mathbf{x}}-x_{18}a\,\hat{\mathbf{y}}+z_{18}a\,\hat{\mathbf{z}}$	(48*h*)	Sm V
$\mathbf{B}_{103}$	=	$(2x_{18}-z_{18})\,\mathbf{a}_1+(-2x_{18}-z_{18})\,\mathbf{a}_2+z_{18}\,\mathbf{a}_3$	=	$-x_{18}a\,\hat{\mathbf{x}}+x_{18}a\,\hat{\mathbf{y}}-z_{18}a\,\hat{\mathbf{z}}$	(48*h*)	Sm V
$\mathbf{B}_{104}$	=	$(-2x_{18}-z_{18})\,\mathbf{a}_1+(2x_{18}-z_{18})\,\mathbf{a}_2+z_{18}\,\mathbf{a}_3$	=	$x_{18}a\,\hat{\mathbf{x}}-x_{18}a\,\hat{\mathbf{y}}-z_{18}a\,\hat{\mathbf{z}}$	(48*h*)	Sm V
$\mathbf{B}_{105}$	=	$(2x_{18}-z_{18})\,\mathbf{a}_1+z_{18}\,\mathbf{a}_2+z_{18}\,\mathbf{a}_3$	=	$z_{18}a\,\hat{\mathbf{x}}+x_{18}a\,\hat{\mathbf{y}}+x_{18}a\,\hat{\mathbf{z}}$	(48*h*)	Sm V
$\mathbf{B}_{106}$	=	$(-2x_{18}-z_{18})\,\mathbf{a}_1+z_{18}\,\mathbf{a}_2+z_{18}\,\mathbf{a}_3$	=	$z_{18}a\,\hat{\mathbf{x}}-x_{18}a\,\hat{\mathbf{y}}-x_{18}a\,\hat{\mathbf{z}}$	(48*h*)	Sm V
$\mathbf{B}_{107}$	=	$z_{18}\,\mathbf{a}_1+(2x_{18}-z_{18})\,\mathbf{a}_2+(-2x_{18}-z_{18})\,\mathbf{a}_3$	=	$-z_{18}a\,\hat{\mathbf{x}}-x_{18}a\,\hat{\mathbf{y}}+x_{18}a\,\hat{\mathbf{z}}$	(48*h*)	Sm V
$\mathbf{B}_{108}$	=	$z_{18}\,\mathbf{a}_1+(-2x_{18}-z_{18})\,\mathbf{a}_2+(2x_{18}-z_{18})\,\mathbf{a}_3$	=	$-z_{18}a\,\hat{\mathbf{x}}+x_{18}a\,\hat{\mathbf{y}}-x_{18}a\,\hat{\mathbf{z}}$	(48*h*)	Sm V
$\mathbf{B}_{109}$	=	$z_{18}\,\mathbf{a}_1+(2x_{18}-z_{18})\,\mathbf{a}_2+z_{18}\,\mathbf{a}_3$	=	$x_{18}a\,\hat{\mathbf{x}}+z_{18}a\,\hat{\mathbf{y}}+x_{18}a\,\hat{\mathbf{z}}$	(48*h*)	Sm V
$\mathbf{B}_{110}$	=	$z_{18}\,\mathbf{a}_1+(-2x_{18}-z_{18})\,\mathbf{a}_2+z_{18}\,\mathbf{a}_3$	=	$-x_{18}a\,\hat{\mathbf{x}}+z_{18}a\,\hat{\mathbf{y}}-x_{18}a\,\hat{\mathbf{z}}$	(48*h*)	Sm V
$\mathbf{B}_{111}$	=	$(-2x_{18}-z_{18})\,\mathbf{a}_1+z_{18}\,\mathbf{a}_2+(2x_{18}-z_{18})\,\mathbf{a}_3$	=	$x_{18}a\,\hat{\mathbf{x}}-z_{18}a\,\hat{\mathbf{y}}-x_{18}a\,\hat{\mathbf{z}}$	(48*h*)	Sm V
$\mathbf{B}_{112}$	=	$(2x_{18}-z_{18})\,\mathbf{a}_1+z_{18}\,\mathbf{a}_2+(-2x_{18}-z_{18})\,\mathbf{a}_3$	=	$-x_{18}a\,\hat{\mathbf{x}}-z_{18}a\,\hat{\mathbf{y}}+x_{18}a\,\hat{\mathbf{z}}$	(48*h*)	Sm V

References:

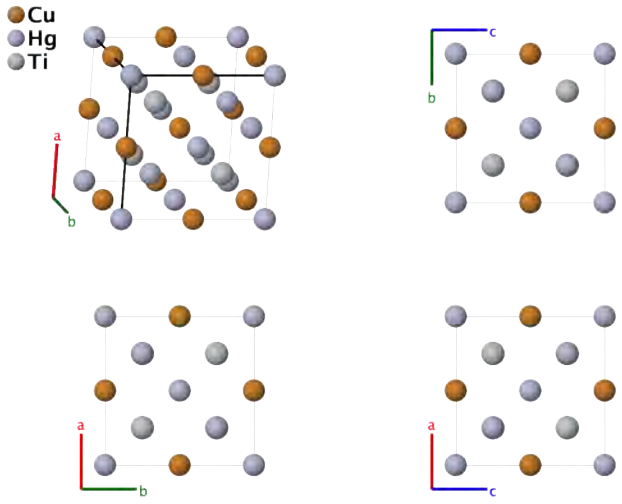
- M. L. Fornasini, B. Chabot, and E. Parthé, *The crystal structure of $Sm_{11}Cd_{45}$ with γ-brass and α-Mn clusters*, Acta Crystallogr. Sect. B Struct. Sci. **34**, 2093–2099 (1978), doi:10.1107/S0567740878007505.

Geometry files:

- CIF: pp. 1793
- POSCAR: pp. 1794

Hg_2TiCu Inverse Heusler Structure: AB2C_cF16_216_b_ad_c

http://aflow.org/prototype-encyclopedia/AB2C_cF16_216_b_ad_c



Prototype	:	$CuHg_2Ti$
AFLOW prototype label	:	AB2C_cF16_216_b_ad_c
***Strukturbericht* designation**	:	None
Pearson symbol	:	cF16
Space group number	:	216
Space group symbol	:	$F\bar{4}3m$
AFLOW prototype command	:	`aflow --proto=AB2C_cF16_216_b_ad_c --params=`a

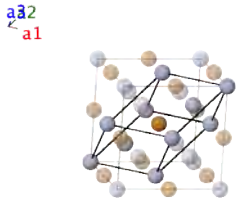
Other compounds with this structure

- Mn_2CoAl, Mn_2CoGa, Mn_2CoIn, Li_2AgSb, Li_2CuSn, Li_2CuSb, Li_2AgAl, Li_2AgIn, Li_2AgSb, Li_2AgSn, Li_2AgPb, Li_2AgBi, Li_2AuGa, Li_2AuIn, Li_2AuSb, Li_2AuSn, Li_2AuPb, and Li_2AuTl

- Most of the literature on Inverse Heusler compounds identifies Hg_2TiCu as the prototype structure, however (Villars, 2016) and (Villars, 2016a) use the Li_2AgSb structure found in (Pauly, 1968) as the prototype. Although it is tempting to use the older paper, we follow the majority and use Hg_2TiCu as the prototype.
- The inverse Heusler structure is a variation of the Quaternary Heusler structure, where the mercury atoms are on adjacent face-centered cubic sublattices, forming a diamond structure. Contrast this with the standard Heusler ($L1_2$) structure, where the like atoms are at second-neighbor positions in the fcc lattice.
- This structure is also referred to as the XA or X_a structure.

Face-centered Cubic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(4*a*)	Hg I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(4*b*)	Cu
$\mathbf{B}_3$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(4*c*)	Ti
$\mathbf{B}_4$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(4*d*)	Hg II

References:

- M. Pušelj and Z. Ban, *The Crystal Structure of $TiCuHg_2$*, Croat. Chem. Acta **41**, 79–83 (1969). http://hrcak.srce.hr/207957.
- H. Pauly, A. Weiss, and H. Witte, *The Crystal Structure of the Ternary Intermetallic Phases Li_2EX (E=Cu, Ag, Au; X=Al, Ga, In, Tl, Si, Ge, Sn, Pb, Sb, Bi)*, Z. Metallkd. **59**, 47–58 (1968).
- P. Villars (Chief Editor), *Li_2AgSb Crystal Structure*, http://materials.springer.com/isp/crystallographic/docs/sd_0461484 (2016). PAULING FILE in: Inorganic Solid Phases, SpringerMaterials (online database), Springer, Heidelberg (ed.).

Found in:

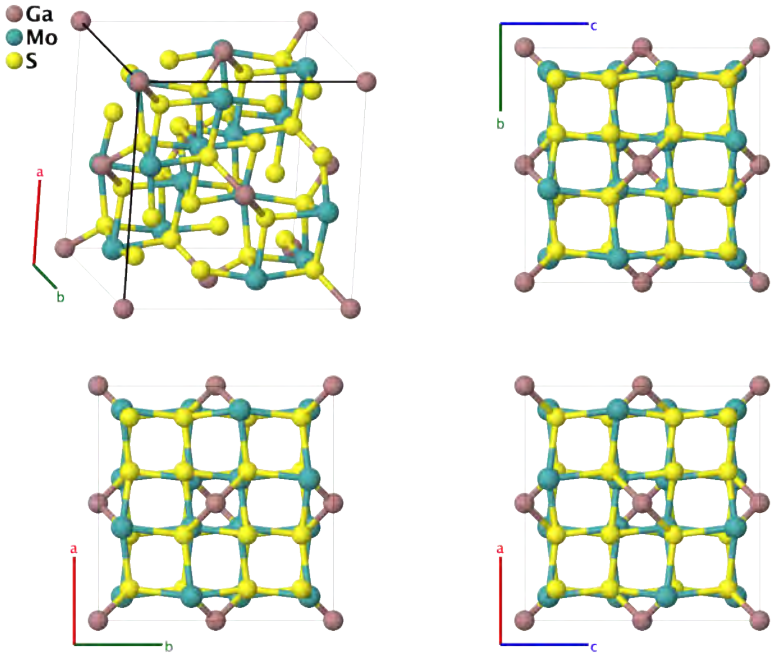
- P. Villars (Chief Editor), *$TiCuHg_2$ ($CuHg_2Ti$) Crystal Structure*, http://materials.springer.com/isp/crystallographic/docs/sd_1214419 (2016). PAULING FILE in: Inorganic Solid Phases, SpringerMaterials (online database), Springer, Heidelberg (ed.).

Geometry files:

- CIF: pp. 1794
- POSCAR: pp. 1795

$GaMo_4S_8$ Structure: AB4C8_cF52_216_a_e_2e

http://aflow.org/prototype-encyclopedia/AB4C8_cF52_216_a_e_2e



Prototype	:	$GaMo_4S_8$
AFLOW prototype label	:	AB4C8_cF52_216_a_e_2e
***Strukturbericht* designation**	:	None
Pearson symbol	:	cF52
Space group number	:	216
Space group symbol	:	$F\bar{4}3m$
AFLOW prototype command	:	`aflow --proto=AB4C8_cF52_216_a_e_2e --params=`a, x_2, x_3, x_4

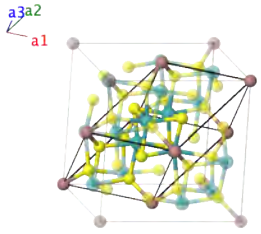
Other compounds with this structure

- $Co(Mo_2Re_2)S_8$, $Fe(Mo_2Re_2)S_8$, $GaMo_4S_4Te_4$, $GaMo_4S_8$, $GaMo_4Se_4Te_4$, $GaMo_4Se_8$, $GaNb_4S_8$, $GaNb_4Se_8$, $GaTa_4S_8$, $GaTa_4Se_8$, GaV_4S_8, GaV_4Se_8, $LaMo_4S_8$, $Ni(Mo_2Re_2)S_8$, and $Zn(Mo_2Re_2)S_8$

- (Ben Yaich, 1984) do not give the lattice constant for $GaMo_4S_8$. We infer it from their interatomic distances.

Face-centered Cubic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(4*a*)	Ga
$\mathbf{B}_2$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$x_2a\,\hat{\mathbf{x}} + x_2a\,\hat{\mathbf{y}} + x_2a\,\hat{\mathbf{z}}$	(16*e*)	Mo
$\mathbf{B}_3$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 - 3x_2\,\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}} - x_2a\,\hat{\mathbf{y}} + x_2a\,\hat{\mathbf{z}}$	(16*e*)	Mo
$\mathbf{B}_4$	=	$x_2\,\mathbf{a}_1 - 3x_2\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}} + x_2a\,\hat{\mathbf{y}} - x_2a\,\hat{\mathbf{z}}$	(16*e*)	Mo
$\mathbf{B}_5$	=	$-3x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$x_2a\,\hat{\mathbf{x}} - x_2a\,\hat{\mathbf{y}} - x_2a\,\hat{\mathbf{z}}$	(16*e*)	Mo
$\mathbf{B}_6$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(16*e*)	S I
$\mathbf{B}_7$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 - 3x_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(16*e*)	S I
$\mathbf{B}_8$	=	$x_3\,\mathbf{a}_1 - 3x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(16*e*)	S I
$\mathbf{B}_9$	=	$-3x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(16*e*)	S I
$\mathbf{B}_{10}$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	(16*e*)	S II
$\mathbf{B}_{11}$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 - 3x_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	(16*e*)	S II
$\mathbf{B}_{12}$	=	$x_4\,\mathbf{a}_1 - 3x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} - x_4a\,\hat{\mathbf{z}}$	(16*e*)	S II
$\mathbf{B}_{13}$	=	$-3x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} - x_4a\,\hat{\mathbf{z}}$	(16*e*)	S II

References:

- H. Ben Yaich, J. C. Jegaden, M. Potel, R. Chevrel, M. Sergent, A. Berton, J. Chaussy, A. K. Rastogi, and R. Tournier, *Nouveaux chalcogenures mixtes $GaMo_4(XX')_8$ (X = S, Se, Te) à clusters tetraedriques Mo_4*, J. Solid State Chem. **51**, 212–217 (1984), doi:10.1016/0022-4596(84)90336-0.

Found in:

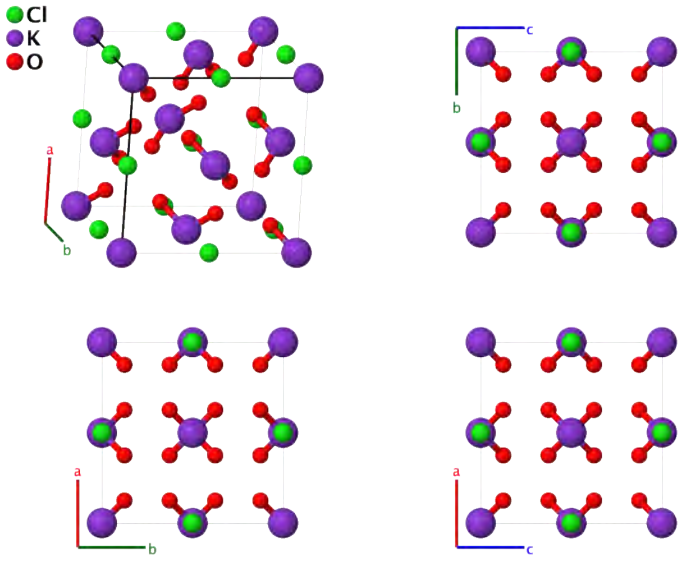
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1795
- POSCAR: pp. 1796

High-Temperature Cubic $KClO_4$ ($H0_5$) Structure: ABC4_cF24_216_b_a_e

http://aflow.org/prototype-encyclopedia/ABC4_cF24_216_b_a_e



Prototype	:	$ClKO_4$
AFLOW prototype label	:	ABC4_cF24_216_b_a_e
***Strukturbericht* designation**	:	$H0_5$
Pearson symbol	:	cF24
Space group number	:	216
Space group symbol	:	$F\bar{4}3m$
AFLOW prototype command	:	`aflow --proto=ABC4_cF24_216_b_a_e --params=`a, x_3

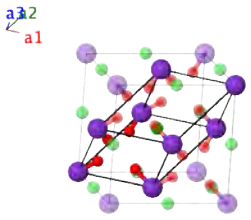
Other compounds with this structure

- $NaClO_4$, $RbClO_4$, $CsClO_4$, NH_4ClO_4, $AgClO_4$, and $TlClO_4$

- This is the high-temperature phase of the listed perchlorate structures. $KClO_4$ transforms from its ground-state orthorhombic structure, $H0_2$ into this structure at 299.5 °C. The transition temperature for the other compounds range from 155 °C ($AgClO_4$) to 308 °C ($NaClO_4$).
- The lattice constant for $KClO_4$ was measured at 310 °C.

Face-centered Cubic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(4*a*)	K
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(4*b*)	Cl
$\mathbf{B}_3$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(16*e*)	O
$\mathbf{B}_4$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 - 3x_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(16*e*)	O
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 - 3x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(16*e*)	O
$\mathbf{B}_6$	=	$-3x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(16*e*)	O

References:

- K. Hermann and W. Ilge, *Röntgenographische Strukturerforschung der kubischen Modifikation der Perchlorate*, Zeitschrift für Kristallographie - Crystalline Materials **75**, 41–66 (1930), doi:10.1515/zkri-1930-0105.

Found in:

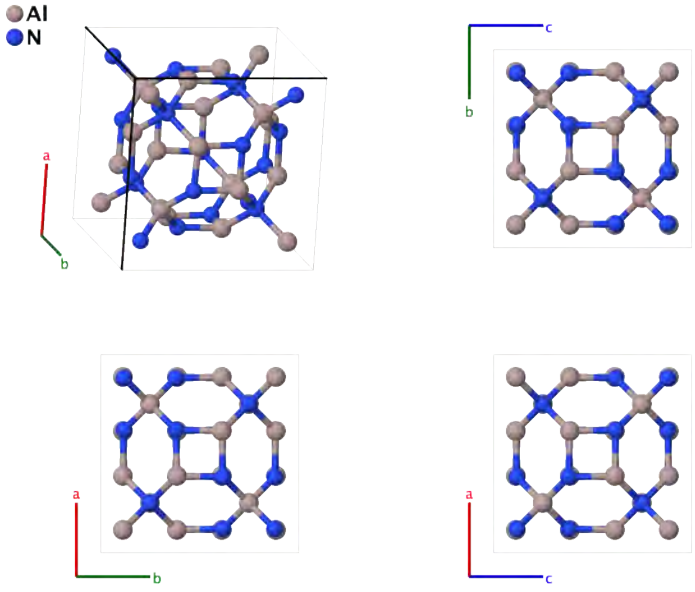
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1796
- POSCAR: pp. 1797

AlN (cF40) Structure: AB_cF40_216_de_ce

http://aflow.org/prototype-encyclopedia/AB_cF40_216_de_ce

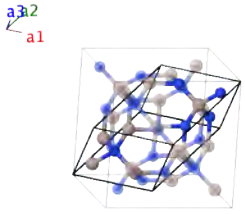


Prototype	:	AlN
AFLOW prototype label	:	AB_cF40_216_de_ce
***Strukturbericht* designation**	:	None
Pearson symbol	:	cF40
Space group number	:	216
Space group symbol	:	$F\bar{4}3m$
AFLOW prototype command	:	`aflow --proto=AB_cF40_216_de_ce --params=`a, x_3, x_4

- AlN naturally occurs in two forms (Liu, 2019): the stable wz-AlN wurtzite ($B4$) structure, and the high-pressure rs-AlN rock salt ($B1$) structure. A metastable zb-AlN zincblende (zb-AlN) structure can be synthesized via a solid-state reaction.
- (Liu, 2019) used a first-principles evolutionary technique to find four possible metastable phases: one in the SC16 structure, and three novel cubic structures, cF40, cI16, and cI24.

Face-centered Cubic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(4*c*)	N I
$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(4*d*)	Al I
$\mathbf{B}_3$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(16*e*)	Al II
$\mathbf{B}_4$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 - 3x_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(16*e*)	Al II
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 - 3x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(16*e*)	Al II
$\mathbf{B}_6$	=	$-3x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(16*e*)	Al II
$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	(16*e*)	N II
$\mathbf{B}_8$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 - 3x_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	(16*e*)	N II
$\mathbf{B}_9$	=	$x_4\,\mathbf{a}_1 - 3x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} - x_4a\,\hat{\mathbf{z}}$	(16*e*)	N II
$\mathbf{B}_{10}$	=	$-3x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} - x_4a\,\hat{\mathbf{z}}$	(16*e*)	N II

References:

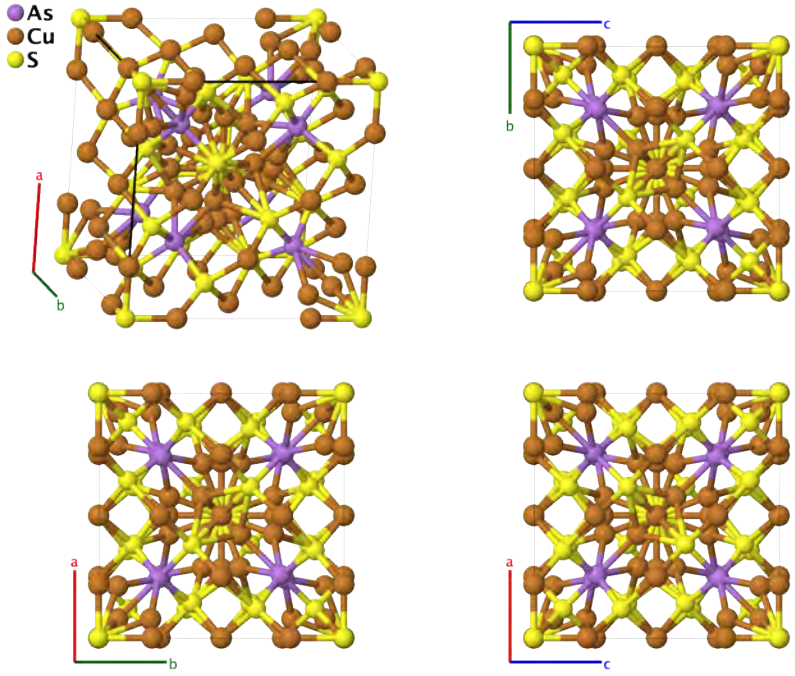
- C. Liu, M. Chen, J. Li, L. Liu, P. Li, M. Ma, C. Shao, J. He, and T. Liang, *A first-principles study of novel cubic AlN phases*, J. Phys. Chem. Solids **130**, 58–66 (2019), doi:10.1016/j.jpcs.2019.02.009.

Geometry files:

- CIF: pp. 1797
- POSCAR: pp. 1798

Tennantite ($Cu_{12}As_4S_{13}$) Structure: A4B24C13_cI82_217_c_deg_ag

http://aflow.org/prototype-encyclopedia/A4B24C13_cI82_217_c_deg_ag



Prototype	:	$As_4Cu_{12}S_{13}$
AFLOW prototype label	:	A4B24C13_cI82_217_c_deg_ag
***Strukturbericht* designation**	:	None
Pearson symbol	:	cI82
Space group number	:	217
Space group symbol	:	$I\bar{4}3m$
AFLOW prototype command	:	`aflow --proto=A4B24C13_cI82_217_c_deg_ag --params=`$a, x_2, x_4, x_5, z_5, x_6, z_6$

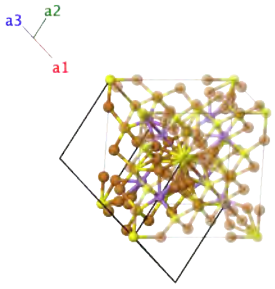
Other compounds with this structure

- $Cu_{14}Sb_4S_{13}$ (tetrahedrite)

- The Cu-II (12*e*) site is only occupied 75.8% of the time, and the Cu-III site is occupied 12.1% of the time, so that these sites only contain twelve atoms.
- Searching (Downs, 2003) shows that natural samples often have antimony substituting for arsenic. The antimony structures (tetrahedrites) contain higher concentrations of copper.
- The AFLOW label models the structure as if the sites were fully occupied.

Body-centered Cubic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= -\tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{y}} + \tfrac{1}{2}\,a\,\hat{\mathbf{z}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2}\,a\,\hat{\mathbf{y}} + \tfrac{1}{2}\,a\,\hat{\mathbf{z}}\\ \mathbf{a}_3 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{y}} - \tfrac{1}{2}\,a\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(2*a*)	S I
$\mathbf{B}_2$	=	$2x_2\,\mathbf{a}_1 + 2x_2\,\mathbf{a}_2 + 2x_2\,\mathbf{a}_3$	=	$x_2a\,\hat{\mathbf{x}} + x_2a\,\hat{\mathbf{y}} + x_2a\,\hat{\mathbf{z}}$	(8*c*)	As
$\mathbf{B}_3$	=	$-2x_2\,\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}} - x_2a\,\hat{\mathbf{y}} + x_2a\,\hat{\mathbf{z}}$	(8*c*)	As
$\mathbf{B}_4$	=	$-2x_2\,\mathbf{a}_2$	=	$-x_2a\,\hat{\mathbf{x}} + x_2a\,\hat{\mathbf{y}} - x_2a\,\hat{\mathbf{z}}$	(8*c*)	As
$\mathbf{B}_5$	=	$-2x_2\,\mathbf{a}_1$	=	$x_2a\,\hat{\mathbf{x}} - x_2a\,\hat{\mathbf{y}} - x_2a\,\hat{\mathbf{z}}$	(8*c*)	As
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$	(12*d*)	Cu I
$\mathbf{B}_7$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(12*d*)	Cu I
$\mathbf{B}_8$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(12*d*)	Cu I
$\mathbf{B}_9$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}}$	(12*d*)	Cu I
$\mathbf{B}_{10}$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(12*d*)	Cu I
$\mathbf{B}_{11}$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(12*d*)	Cu I
$\mathbf{B}_{12}$	=	$x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}}$	(12*e*)	Cu II
$\mathbf{B}_{13}$	=	$-x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}}$	(12*e*)	Cu II
$\mathbf{B}_{14}$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{y}}$	(12*e*)	Cu II
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{y}}$	(12*e*)	Cu II
$\mathbf{B}_{16}$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2$	=	$x_4a\,\hat{\mathbf{z}}$	(12*e*)	Cu II
$\mathbf{B}_{17}$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2$	=	$-x_4a\,\hat{\mathbf{z}}$	(12*e*)	Cu II
$\mathbf{B}_{18}$	=	$(x_5 + z_5)\,\mathbf{a}_1 + (x_5 + z_5)\,\mathbf{a}_2 + 2x_5\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}} + x_5a\,\hat{\mathbf{y}} + z_5a\,\hat{\mathbf{z}}$	(24*g*)	Cu III
$\mathbf{B}_{19}$	=	$(-x_5 + z_5)\,\mathbf{a}_1 + (-x_5 + z_5)\,\mathbf{a}_2 - 2x_5\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}} - x_5a\,\hat{\mathbf{y}} + z_5a\,\hat{\mathbf{z}}$	(24*g*)	Cu III
$\mathbf{B}_{20}$	=	$(x_5 - z_5)\,\mathbf{a}_1 + (-x_5 - z_5)\,\mathbf{a}_2$	=	$-x_5a\,\hat{\mathbf{x}} + x_5a\,\hat{\mathbf{y}} - z_5a\,\hat{\mathbf{z}}$	(24*g*)	Cu III
$\mathbf{B}_{21}$	=	$(-x_5 - z_5)\,\mathbf{a}_1 + (x_5 - z_5)\,\mathbf{a}_2$	=	$x_5a\,\hat{\mathbf{x}} - x_5a\,\hat{\mathbf{y}} - z_5a\,\hat{\mathbf{z}}$	(24*g*)	Cu III
$\mathbf{B}_{22}$	=	$2x_5\,\mathbf{a}_1 + (x_5 + z_5)\,\mathbf{a}_2 + (x_5 + z_5)\,\mathbf{a}_3$	=	$z_5a\,\hat{\mathbf{x}} + x_5a\,\hat{\mathbf{y}} + x_5a\,\hat{\mathbf{z}}$	(24*g*)	Cu III
$\mathbf{B}_{23}$	=	$-2x_5\,\mathbf{a}_1 + (-x_5 + z_5)\,\mathbf{a}_2 + (-x_5 + z_5)\,\mathbf{a}_3$	=	$z_5a\,\hat{\mathbf{x}} - x_5a\,\hat{\mathbf{y}} - x_5a\,\hat{\mathbf{z}}$	(24*g*)	Cu III
$\mathbf{B}_{24}$	=	$(x_5 - z_5)\,\mathbf{a}_2 + (-x_5 - z_5)\,\mathbf{a}_3$	=	$-z_5a\,\hat{\mathbf{x}} - x_5a\,\hat{\mathbf{y}} + x_5a\,\hat{\mathbf{z}}$	(24*g*)	Cu III
$\mathbf{B}_{25}$	=	$(-x_5 - z_5)\,\mathbf{a}_2 + (x_5 - z_5)\,\mathbf{a}_3$	=	$-z_5a\,\hat{\mathbf{x}} + x_5a\,\hat{\mathbf{y}} - x_5a\,\hat{\mathbf{z}}$	(24*g*)	Cu III
$\mathbf{B}_{26}$	=	$(x_5 + z_5)\,\mathbf{a}_1 + 2x_5\,\mathbf{a}_2 + (x_5 + z_5)\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}} + z_5a\,\hat{\mathbf{y}} + x_5a\,\hat{\mathbf{z}}$	(24*g*)	Cu III
$\mathbf{B}_{27}$	=	$(-x_5 + z_5)\,\mathbf{a}_1 - 2x_5\,\mathbf{a}_2 + (-x_5 + z_5)\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}} + z_5a\,\hat{\mathbf{y}} - x_5a\,\hat{\mathbf{z}}$	(24*g*)	Cu III

$\mathbf{B}_{28}$	=	$(-x_5 - z_5)\,\mathbf{a}_1 + (x_5 - z_5)\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} - z_5 a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}}$	(24g)	Cu III
$\mathbf{B}_{29}$	=	$(x_5 - z_5)\,\mathbf{a}_1 + (-x_5 - z_5)\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - z_5 a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}}$	(24g)	Cu III
$\mathbf{B}_{30}$	=	$(x_6 + z_6)\,\mathbf{a}_1 + (x_6 + z_6)\,\mathbf{a}_2 + 2x_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} + z_6 a\,\hat{\mathbf{z}}$	(24g)	S II
$\mathbf{B}_{31}$	=	$(-x_6 + z_6)\,\mathbf{a}_1 + (-x_6 + z_6)\,\mathbf{a}_2 - 2x_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} + z_6 a\,\hat{\mathbf{z}}$	(24g)	S II
$\mathbf{B}_{32}$	=	$(x_6 - z_6)\,\mathbf{a}_1 + (-x_6 - z_6)\,\mathbf{a}_2$	=	$-x_6 a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} - z_6 a\,\hat{\mathbf{z}}$	(24g)	S II
$\mathbf{B}_{33}$	=	$(-x_6 - z_6)\,\mathbf{a}_1 + (x_6 - z_6)\,\mathbf{a}_2$	=	$x_6 a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} - z_6 a\,\hat{\mathbf{z}}$	(24g)	S II
$\mathbf{B}_{34}$	=	$2x_6\,\mathbf{a}_1 + (x_6 + z_6)\,\mathbf{a}_2 + (x_6 + z_6)\,\mathbf{a}_3$	=	$z_6 a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} + x_6 a\,\hat{\mathbf{z}}$	(24g)	S II
$\mathbf{B}_{35}$	=	$-2x_6\,\mathbf{a}_1 + (-x_6 + z_6)\,\mathbf{a}_2 + (-x_6 + z_6)\,\mathbf{a}_3$	=	$z_6 a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} - x_6 a\,\hat{\mathbf{z}}$	(24g)	S II
$\mathbf{B}_{36}$	=	$(x_6 - z_6)\,\mathbf{a}_2 + (-x_6 - z_6)\,\mathbf{a}_3$	=	$-z_6 a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} + x_6 a\,\hat{\mathbf{z}}$	(24g)	S II
$\mathbf{B}_{37}$	=	$(-x_6 - z_6)\,\mathbf{a}_2 + (x_6 - z_6)\,\mathbf{a}_3$	=	$-z_6 a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} - x_6 a\,\hat{\mathbf{z}}$	(24g)	S II
$\mathbf{B}_{38}$	=	$(x_6 + z_6)\,\mathbf{a}_1 + 2x_6\,\mathbf{a}_2 + (x_6 + z_6)\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + z_6 a\,\hat{\mathbf{y}} + x_6 a\,\hat{\mathbf{z}}$	(24g)	S II
$\mathbf{B}_{39}$	=	$(-x_6 + z_6)\,\mathbf{a}_1 - 2x_6\,\mathbf{a}_2 + (-x_6 + z_6)\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + z_6 a\,\hat{\mathbf{y}} - x_6 a\,\hat{\mathbf{z}}$	(24g)	S II
$\mathbf{B}_{40}$	=	$(-x_6 - z_6)\,\mathbf{a}_1 + (x_6 - z_6)\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} - z_6 a\,\hat{\mathbf{y}} - x_6 a\,\hat{\mathbf{z}}$	(24g)	S II
$\mathbf{B}_{41}$	=	$(x_6 - z_6)\,\mathbf{a}_1 + (-x_6 - z_6)\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - z_6 a\,\hat{\mathbf{y}} + x_6 a\,\hat{\mathbf{z}}$	(24g)	S II

References:

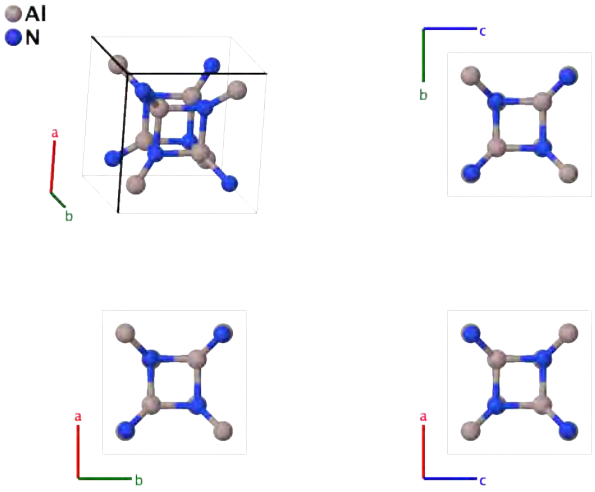
- A. A. Yaroslavzev, A. V. Mironov, A. N. Kuznetsov, A. P. Dudka, and O. N. Khrykina, *Tennantite: multi-temperature crystal structure, phase transition and electronic structure of synthetic* $Cu_{12}As_4S_{13}$, Acta Crystallogr. Sect. B Struct. Sci. **75**, 634–642 (2019), doi:10.1107/S2052520619007595.
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1798
- POSCAR: pp. 1798

AlN (cI16) Structure: AB_cI16_217_c_c

http://aflow.org/prototype-encyclopedia/AB_cI16_217_c_c

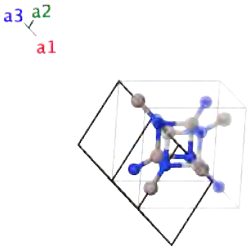


Prototype	:	AlN
AFLOW prototype label	:	AB_cI16_217_c_c
***Strukturbericht* designation**	:	None
Pearson symbol	:	cI16
Space group number	:	217
Space group symbol	:	$I\bar{4}3m$
AFLOW prototype command	:	`aflow --proto=AB_cI16_217_c_c --params=`a, x_1, x_2

- AlN naturally occurs in two forms (Liu, 2019): the stable wz-AlN wurtzite ($B4$) structure, and the high-pressure rs-AlN rock salt ($B1$) structure. A metastable zb-AlN zincblende (zb-AlN) structure can be synthesized via a solid-state reaction.
- (Liu, 2019) used a first-principles evolutionary technique to find four possible metastable phases: one in the SC16 structure, and three novel cubic structures, cF40, cI16, and cI24.

Body-centered Cubic primitive vectors:

$$\mathbf{a}_1 = -\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}a\,\hat{\mathbf{x}} - \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} - \tfrac{1}{2}a\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$ =	$2x_1\,\mathbf{a}_1 + 2x_1\,\mathbf{a}_2 + 2x_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + x_1 a\,\hat{\mathbf{y}} + x_1 a\,\hat{\mathbf{z}}$	($8c$)	Al

$\mathbf{B}_2$	=	$-2x_1\,\mathbf{a}_3$	=	$-x_1a\,\hat{\mathbf{x}} - x_1a\,\hat{\mathbf{y}} + x_1a\,\hat{\mathbf{z}}$	(8*c*)	Al
$\mathbf{B}_3$	=	$-2x_1\,\mathbf{a}_2$	=	$-x_1a\,\hat{\mathbf{x}} + x_1a\,\hat{\mathbf{y}} - x_1a\,\hat{\mathbf{z}}$	(8*c*)	Al
$\mathbf{B}_4$	=	$-2x_1\,\mathbf{a}_1$	=	$x_1a\,\hat{\mathbf{x}} - x_1a\,\hat{\mathbf{y}} - x_1a\,\hat{\mathbf{z}}$	(8*c*)	Al
$\mathbf{B}_5$	=	$2x_2\,\mathbf{a}_1 + 2x_2\,\mathbf{a}_2 + 2x_2\,\mathbf{a}_3$	=	$x_2a\,\hat{\mathbf{x}} + x_2a\,\hat{\mathbf{y}} + x_2a\,\hat{\mathbf{z}}$	(8*c*)	N
$\mathbf{B}_6$	=	$-2x_2\,\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}} - x_2a\,\hat{\mathbf{y}} + x_2a\,\hat{\mathbf{z}}$	(8*c*)	N
$\mathbf{B}_7$	=	$-2x_2\,\mathbf{a}_2$	=	$-x_2a\,\hat{\mathbf{x}} + x_2a\,\hat{\mathbf{y}} - x_2a\,\hat{\mathbf{z}}$	(8*c*)	N
$\mathbf{B}_8$	=	$-2x_2\,\mathbf{a}_1$	=	$x_2a\,\hat{\mathbf{x}} - x_2a\,\hat{\mathbf{y}} - x_2a\,\hat{\mathbf{z}}$	(8*c*)	N

References:

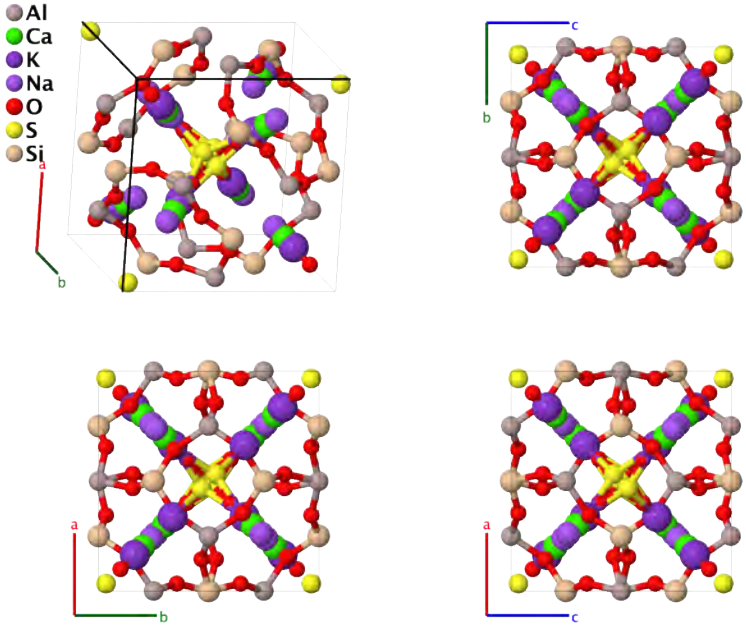
- C. Liu, M. Chen, J. Li, L. Liu, P. Li, M. Ma, C. Shao, J. He, and T. Liang, *A first-principles study of novel cubic AlN phases*, J. Phys. Chem. Solids **130**, 58–66 (2019), doi:10.1016/j.jpcs.2019.02.009.

Geometry files:

- CIF: pp. 1798
- POSCAR: pp. 1799

Hauyne [$(Na_{0.5}Ca_{0.3}K_{0.2})_8(Al_6Si_6O_{24})(SO_4)_{1.5}$, $S6_9$] Structure: A3B4C4D4E16F4G3_cP76_218_c_e_e_e_ei_e_d

http://aflow.org/prototype-encyclopedia/A3B4C4D4E16F4G3_cP76_218_c_e_e_e_ei_e_d



Prototype	:	$Al_6Ca_{2.4}K_{1.6}Na_4O_{30}S_{1.5}Si_6$
AFLOW prototype label	:	A3B4C4D4E16F4G3_cP76_218_c_e_e_e_ei_e_d
***Strukturbericht* designation**	:	$S6_9$
Pearson symbol	:	cP76
Space group number	:	218
Space group symbol	:	$P\bar{4}3n$
AFLOW prototype command	:	`aflow --proto=A3B4C4D4E16F4G3_cP76_218_c_e_e_e_ei_e_d --params=`$a, x_3, x_4, x_5, x_6, x_7, x_8, y_8, z_8$

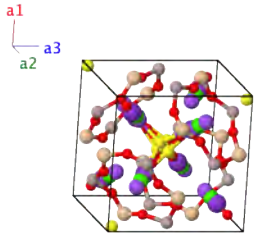
- (Gottfried, 1937) used the work of (Machatschki, 1934) for *Strukturbericht* label $S6_9$. We have used the updated 153 K data of (Hassan, 1991).
- (Machatschki, 1934) puts the sodium, calcium, and potassium atoms at the same (8*e*) Wyckoff position, with partial occupancies of approximately $Na_{0.75}Ca_{0.25}$ and a trace of potassium. The sample studied by (Hassan, 1991) was found to have the three atoms at slightly different (8*e*) positions, with 54% sodium occupation, 30% calcium occupation, and 20% potassium occupation.
- (Machatschki, 1934) put the sulfur atom at the origin, Wyckoff position (2*a*), and assumed that there were two SO_4 molecules per formula unit. (Hassan, 1991) found that there were statistically only 1.5 molecules per formula unit, and that the sulfur atom was slightly displaced from the origin to a (8*e*) site, where each atomic position was occupied only 19% of the time. The corresponding (8*e*) oxygen site (O-I in our notation, O2 in Hassan) is only occupied 75% of the time to maintain the SO_4 stoichiometry.
- We shifted the origin of (Hassan, 1991) so that the resulting atomic positions are close to those reported by (Machatschki, 1934). The latter structure can be recovered from the former by moving all of the sodium, calcium,

and potassium atoms to the same Wyckoff position, moving the sulfur atom to the origin, Wyckoff position ($2a$), and adjusting the occupation of each site appropriately.

- The 153K data presented in (Downs, 2003) has errors in the z coordinates of the CaC2 (our Ca) and O2 (our O-I) atoms. In both cases, the z coordinate should be the same as the x and y coordinates.
- The AFLOW label models the structure as if the sites were fully occupied.

Simple Cubic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = a\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$	($6c$)	Al
$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$	($6c$)	Al
$\mathbf{B}_3$	=	$\frac{1}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	($6c$)	Al
$\mathbf{B}_4$	=	$\frac{3}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	($6c$)	Al
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	($6c$)	Al
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	($6c$)	Al
$\mathbf{B}_7$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	($6d$)	Si
$\mathbf{B}_8$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	($6d$)	Si
$\mathbf{B}_9$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}}$	($6d$)	Si
$\mathbf{B}_{10}$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}}$	($6d$)	Si
$\mathbf{B}_{11}$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	($6d$)	Si
$\mathbf{B}_{12}$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	($6d$)	Si
$\mathbf{B}_{13}$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	($8e$)	Ca
$\mathbf{B}_{14}$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	($8e$)	Ca
$\mathbf{B}_{15}$	=	$-x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	($8e$)	Ca
$\mathbf{B}_{16}$	=	$x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	($8e$)	Ca
$\mathbf{B}_{17}$	=	$\left(\frac{1}{2}+x_3\right)\mathbf{a}_1 + \left(\frac{1}{2}+x_3\right)\mathbf{a}_2 + \left(\frac{1}{2}+x_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{z}}$	($8e$)	Ca
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2}-x_3\right)\mathbf{a}_1 + \left(\frac{1}{2}-x_3\right)\mathbf{a}_2 + \left(\frac{1}{2}+x_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{z}}$	($8e$)	Ca
$\mathbf{B}_{19}$	=	$\left(\frac{1}{2}+x_3\right)\mathbf{a}_1 + \left(\frac{1}{2}-x_3\right)\mathbf{a}_2 + \left(\frac{1}{2}-x_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{z}}$	($8e$)	Ca

$$
\begin{array}{lcccccc}
\mathbf{B}_{20} & = & \left(\frac{1}{2}-x_3\right)\mathbf{a}_1+\left(\frac{1}{2}+x_3\right)\mathbf{a}_2+\left(\frac{1}{2}-x_3\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_3\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-x_3\right)a\,\hat{\mathbf{z}} & (8e) & \text{Ca}\\
\mathbf{B}_{21} & = & x_4\,\mathbf{a}_1+x_4\,\mathbf{a}_2+x_4\,\mathbf{a}_3 & = & x_4a\,\hat{\mathbf{x}}+x_4a\,\hat{\mathbf{y}}+x_4a\,\hat{\mathbf{z}} & (8e) & \text{K}\\
\mathbf{B}_{22} & = & -x_4\,\mathbf{a}_1-x_4\,\mathbf{a}_2+x_4\,\mathbf{a}_3 & = & -x_4a\,\hat{\mathbf{x}}-x_4a\,\hat{\mathbf{y}}+x_4a\,\hat{\mathbf{z}} & (8e) & \text{K}\\
\mathbf{B}_{23} & = & -x_4\,\mathbf{a}_1+x_4\,\mathbf{a}_2-x_4\,\mathbf{a}_3 & = & -x_4a\,\hat{\mathbf{x}}+x_4a\,\hat{\mathbf{y}}-x_4a\,\hat{\mathbf{z}} & (8e) & \text{K}\\
\mathbf{B}_{24} & = & x_4\,\mathbf{a}_1-x_4\,\mathbf{a}_2-x_4\,\mathbf{a}_3 & = & x_4a\,\hat{\mathbf{x}}-x_4a\,\hat{\mathbf{y}}-x_4a\,\hat{\mathbf{z}} & (8e) & \text{K}\\
\mathbf{B}_{25} & = & \left(\frac{1}{2}+x_4\right)\mathbf{a}_1+\left(\frac{1}{2}+x_4\right)\mathbf{a}_2+\left(\frac{1}{2}+x_4\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{z}} & (8e) & \text{K}\\
\mathbf{B}_{26} & = & \left(\frac{1}{2}-x_4\right)\mathbf{a}_1+\left(\frac{1}{2}-x_4\right)\mathbf{a}_2+\left(\frac{1}{2}+x_4\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{z}} & (8e) & \text{K}\\
\mathbf{B}_{27} & = & \left(\frac{1}{2}+x_4\right)\mathbf{a}_1+\left(\frac{1}{2}-x_4\right)\mathbf{a}_2+\left(\frac{1}{2}-x_4\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{z}} & (8e) & \text{K}\\
\mathbf{B}_{28} & = & \left(\frac{1}{2}-x_4\right)\mathbf{a}_1+\left(\frac{1}{2}+x_4\right)\mathbf{a}_2+\left(\frac{1}{2}-x_4\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{z}} & (8e) & \text{K}\\
\mathbf{B}_{29} & = & x_5\,\mathbf{a}_1+x_5\,\mathbf{a}_2+x_5\,\mathbf{a}_3 & = & x_5a\,\hat{\mathbf{x}}+x_5a\,\hat{\mathbf{y}}+x_5a\,\hat{\mathbf{z}} & (8e) & \text{Na}\\
\mathbf{B}_{30} & = & -x_5\,\mathbf{a}_1-x_5\,\mathbf{a}_2+x_5\,\mathbf{a}_3 & = & -x_5a\,\hat{\mathbf{x}}-x_5a\,\hat{\mathbf{y}}+x_5a\,\hat{\mathbf{z}} & (8e) & \text{Na}\\
\mathbf{B}_{31} & = & -x_5\,\mathbf{a}_1+x_5\,\mathbf{a}_2-x_5\,\mathbf{a}_3 & = & -x_5a\,\hat{\mathbf{x}}+x_5a\,\hat{\mathbf{y}}-x_5a\,\hat{\mathbf{z}} & (8e) & \text{Na}\\
\mathbf{B}_{32} & = & x_5\,\mathbf{a}_1-x_5\,\mathbf{a}_2-x_5\,\mathbf{a}_3 & = & x_5a\,\hat{\mathbf{x}}-x_5a\,\hat{\mathbf{y}}-x_5a\,\hat{\mathbf{z}} & (8e) & \text{Na}\\
\mathbf{B}_{33} & = & \left(\frac{1}{2}+x_5\right)\mathbf{a}_1+\left(\frac{1}{2}+x_5\right)\mathbf{a}_2+\left(\frac{1}{2}+x_5\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{z}} & (8e) & \text{Na}\\
\mathbf{B}_{34} & = & \left(\frac{1}{2}-x_5\right)\mathbf{a}_1+\left(\frac{1}{2}-x_5\right)\mathbf{a}_2+\left(\frac{1}{2}+x_5\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{z}} & (8e) & \text{Na}\\
\mathbf{B}_{35} & = & \left(\frac{1}{2}+x_5\right)\mathbf{a}_1+\left(\frac{1}{2}-x_5\right)\mathbf{a}_2+\left(\frac{1}{2}-x_5\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{z}} & (8e) & \text{Na}\\
\mathbf{B}_{36} & = & \left(\frac{1}{2}-x_5\right)\mathbf{a}_1+\left(\frac{1}{2}+x_5\right)\mathbf{a}_2+\left(\frac{1}{2}-x_5\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{z}} & (8e) & \text{Na}\\
\mathbf{B}_{37} & = & x_6\,\mathbf{a}_1+x_6\,\mathbf{a}_2+x_6\,\mathbf{a}_3 & = & x_6a\,\hat{\mathbf{x}}+x_6a\,\hat{\mathbf{y}}+x_6a\,\hat{\mathbf{z}} & (8e) & \text{O I}\\
\mathbf{B}_{38} & = & -x_6\,\mathbf{a}_1-x_6\,\mathbf{a}_2+x_6\,\mathbf{a}_3 & = & -x_6a\,\hat{\mathbf{x}}-x_6a\,\hat{\mathbf{y}}+x_6a\,\hat{\mathbf{z}} & (8e) & \text{O I}\\
\mathbf{B}_{39} & = & -x_6\,\mathbf{a}_1+x_6\,\mathbf{a}_2-x_6\,\mathbf{a}_3 & = & -x_6a\,\hat{\mathbf{x}}+x_6a\,\hat{\mathbf{y}}-x_6a\,\hat{\mathbf{z}} & (8e) & \text{O I}\\
\mathbf{B}_{40} & = & x_6\,\mathbf{a}_1-x_6\,\mathbf{a}_2-x_6\,\mathbf{a}_3 & = & x_6a\,\hat{\mathbf{x}}-x_6a\,\hat{\mathbf{y}}-x_6a\,\hat{\mathbf{z}} & (8e) & \text{O I}\\
\mathbf{B}_{41} & = & \left(\frac{1}{2}+x_6\right)\mathbf{a}_1+\left(\frac{1}{2}+x_6\right)\mathbf{a}_2+\left(\frac{1}{2}+x_6\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{z}} & (8e) & \text{O I}\\
\mathbf{B}_{42} & = & \left(\frac{1}{2}-x_6\right)\mathbf{a}_1+\left(\frac{1}{2}-x_6\right)\mathbf{a}_2+\left(\frac{1}{2}+x_6\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{z}} & (8e) & \text{O I}\\
\mathbf{B}_{43} & = & \left(\frac{1}{2}+x_6\right)\mathbf{a}_1+\left(\frac{1}{2}-x_6\right)\mathbf{a}_2+\left(\frac{1}{2}-x_6\right)\mathbf{a}_3 & = & \left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{z}} & (8e) & \text{O I}\\
\mathbf{B}_{44} & = & \left(\frac{1}{2}-x_6\right)\mathbf{a}_1+\left(\frac{1}{2}+x_6\right)\mathbf{a}_2+\left(\frac{1}{2}-x_6\right)\mathbf{a}_3 & = & \left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{2}+x_6\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-x_6\right)a\,\hat{\mathbf{z}} & (8e) & \text{O I}\\
\mathbf{B}_{45} & = & x_7\,\mathbf{a}_1+x_7\,\mathbf{a}_2+x_7\,\mathbf{a}_3 & = & x_7a\,\hat{\mathbf{x}}+x_7a\,\hat{\mathbf{y}}+x_7a\,\hat{\mathbf{z}} & (8e) & \text{S}
\end{array}
$$

$\mathbf{B}_{46}$	=	$-x_7\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 + x_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} - x_7 a\,\hat{\mathbf{y}} + x_7 a\,\hat{\mathbf{z}}$	(8*e*)	S
$\mathbf{B}_{47}$	=	$-x_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 - x_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} + x_7 a\,\hat{\mathbf{y}} - x_7 a\,\hat{\mathbf{z}}$	(8*e*)	S
$\mathbf{B}_{48}$	=	$x_7\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 - x_7\,\mathbf{a}_3$	=	$x_7 a\,\hat{\mathbf{x}} - x_7 a\,\hat{\mathbf{y}} - x_7 a\,\hat{\mathbf{z}}$	(8*e*)	S
$\mathbf{B}_{49}$	=	$\left(\frac{1}{2}+x_7\right)\mathbf{a}_1 + \left(\frac{1}{2}+x_7\right)\mathbf{a}_2 + \left(\frac{1}{2}+x_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{z}}$	(8*e*)	S
$\mathbf{B}_{50}$	=	$\left(\frac{1}{2}-x_7\right)\mathbf{a}_1 + \left(\frac{1}{2}-x_7\right)\mathbf{a}_2 + \left(\frac{1}{2}+x_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_7\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{z}}$	(8*e*)	S
$\mathbf{B}_{51}$	=	$\left(\frac{1}{2}+x_7\right)\mathbf{a}_1 + \left(\frac{1}{2}-x_7\right)\mathbf{a}_2 + \left(\frac{1}{2}-x_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_7\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-x_7\right)a\,\hat{\mathbf{z}}$	(8*e*)	S
$\mathbf{B}_{52}$	=	$\left(\frac{1}{2}-x_7\right)\mathbf{a}_1 + \left(\frac{1}{2}+x_7\right)\mathbf{a}_2 + \left(\frac{1}{2}-x_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-x_7\right)a\,\hat{\mathbf{z}}$	(8*e*)	S
$\mathbf{B}_{53}$	=	$x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + y_8 a\,\hat{\mathbf{y}} + z_8 a\,\hat{\mathbf{z}}$	(24*i*)	O II
$\mathbf{B}_{54}$	=	$-x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} - y_8 a\,\hat{\mathbf{y}} + z_8 a\,\hat{\mathbf{z}}$	(24*i*)	O II
$\mathbf{B}_{55}$	=	$-x_8\,\mathbf{a}_1 + y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} + y_8 a\,\hat{\mathbf{y}} - z_8 a\,\hat{\mathbf{z}}$	(24*i*)	O II
$\mathbf{B}_{56}$	=	$x_8\,\mathbf{a}_1 - y_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} - y_8 a\,\hat{\mathbf{y}} - z_8 a\,\hat{\mathbf{z}}$	(24*i*)	O II
$\mathbf{B}_{57}$	=	$z_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 + y_8\,\mathbf{a}_3$	=	$z_8 a\,\hat{\mathbf{x}} + x_8 a\,\hat{\mathbf{y}} + y_8 a\,\hat{\mathbf{z}}$	(24*i*)	O II
$\mathbf{B}_{58}$	=	$z_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 - y_8\,\mathbf{a}_3$	=	$z_8 a\,\hat{\mathbf{x}} - x_8 a\,\hat{\mathbf{y}} - y_8 a\,\hat{\mathbf{z}}$	(24*i*)	O II
$\mathbf{B}_{59}$	=	$-z_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 + y_8\,\mathbf{a}_3$	=	$-z_8 a\,\hat{\mathbf{x}} - x_8 a\,\hat{\mathbf{y}} + y_8 a\,\hat{\mathbf{z}}$	(24*i*)	O II
$\mathbf{B}_{60}$	=	$-z_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 - y_8\,\mathbf{a}_3$	=	$-z_8 a\,\hat{\mathbf{x}} + x_8 a\,\hat{\mathbf{y}} - y_8 a\,\hat{\mathbf{z}}$	(24*i*)	O II
$\mathbf{B}_{61}$	=	$y_8\,\mathbf{a}_1 + z_8\,\mathbf{a}_2 + x_8\,\mathbf{a}_3$	=	$y_8 a\,\hat{\mathbf{x}} + z_8 a\,\hat{\mathbf{y}} + x_8 a\,\hat{\mathbf{z}}$	(24*i*)	O II
$\mathbf{B}_{62}$	=	$-y_8\,\mathbf{a}_1 + z_8\,\mathbf{a}_2 - x_8\,\mathbf{a}_3$	=	$-y_8 a\,\hat{\mathbf{x}} + z_8 a\,\hat{\mathbf{y}} - x_8 a\,\hat{\mathbf{z}}$	(24*i*)	O II
$\mathbf{B}_{63}$	=	$y_8\,\mathbf{a}_1 - z_8\,\mathbf{a}_2 - x_8\,\mathbf{a}_3$	=	$y_8 a\,\hat{\mathbf{x}} - z_8 a\,\hat{\mathbf{y}} - x_8 a\,\hat{\mathbf{z}}$	(24*i*)	O II
$\mathbf{B}_{64}$	=	$-y_8\,\mathbf{a}_1 - z_8\,\mathbf{a}_2 + x_8\,\mathbf{a}_3$	=	$-y_8 a\,\hat{\mathbf{x}} - z_8 a\,\hat{\mathbf{y}} + x_8 a\,\hat{\mathbf{z}}$	(24*i*)	O II
$\mathbf{B}_{65}$	=	$\left(\frac{1}{2}+y_8\right)\mathbf{a}_1 + \left(\frac{1}{2}+x_8\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_8\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_8\right)a\,\hat{\mathbf{z}}$	(24*i*)	O II
$\mathbf{B}_{66}$	=	$\left(\frac{1}{2}-y_8\right)\mathbf{a}_1 + \left(\frac{1}{2}-x_8\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_8\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_8\right)a\,\hat{\mathbf{z}}$	(24*i*)	O II
$\mathbf{B}_{67}$	=	$\left(\frac{1}{2}+y_8\right)\mathbf{a}_1 + \left(\frac{1}{2}-x_8\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_8\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_8\right)a\,\hat{\mathbf{z}}$	(24*i*)	O II
$\mathbf{B}_{68}$	=	$\left(\frac{1}{2}-y_8\right)\mathbf{a}_1 + \left(\frac{1}{2}+x_8\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_8\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_8\right)a\,\hat{\mathbf{z}}$	(24*i*)	O II
$\mathbf{B}_{69}$	=	$\left(\frac{1}{2}+x_8\right)\mathbf{a}_1 + \left(\frac{1}{2}+z_8\right)\mathbf{a}_2 + \left(\frac{1}{2}+y_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+z_8\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+y_8\right)a\,\hat{\mathbf{z}}$	(24*i*)	O II
$\mathbf{B}_{70}$	=	$\left(\frac{1}{2}-x_8\right)\mathbf{a}_1 + \left(\frac{1}{2}+z_8\right)\mathbf{a}_2 + \left(\frac{1}{2}-y_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+z_8\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-y_8\right)a\,\hat{\mathbf{z}}$	(24*i*)	O II
$\mathbf{B}_{71}$	=	$\left(\frac{1}{2}-x_8\right)\mathbf{a}_1 + \left(\frac{1}{2}-z_8\right)\mathbf{a}_2 + \left(\frac{1}{2}+y_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-z_8\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+y_8\right)a\,\hat{\mathbf{z}}$	(24*i*)	O II
$\mathbf{B}_{72}$	=	$\left(\frac{1}{2}+x_8\right)\mathbf{a}_1 + \left(\frac{1}{2}-z_8\right)\mathbf{a}_2 + \left(\frac{1}{2}-y_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-z_8\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-y_8\right)a\,\hat{\mathbf{z}}$	(24*i*)	O II

$$\mathbf{B}_{73} = \left(\tfrac{1}{2}+z_8\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+z_8\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_8\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{z}} \quad (24i) \quad \text{O II}$$

$$\mathbf{B}_{74} = \left(\tfrac{1}{2}+z_8\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}-x_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+z_8\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_8\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{z}} \quad (24i) \quad \text{O II}$$

$$\mathbf{B}_{75} = \left(\tfrac{1}{2}-z_8\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}-x_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-z_8\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_8\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{z}} \quad (24i) \quad \text{O II}$$

$$\mathbf{B}_{76} = \left(\tfrac{1}{2}-z_8\right)\mathbf{a}_1+\left(\tfrac{1}{2}-y_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_8\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-z_8\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-y_8\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{z}} \quad (24i) \quad \text{O II}$$

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- F. Machatschki, *Kristallstruktur von Nauyn und Nosean*, Zbl. Mineral. Geol. und Paläont. A pp. 136–144 (1934).
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Found in:

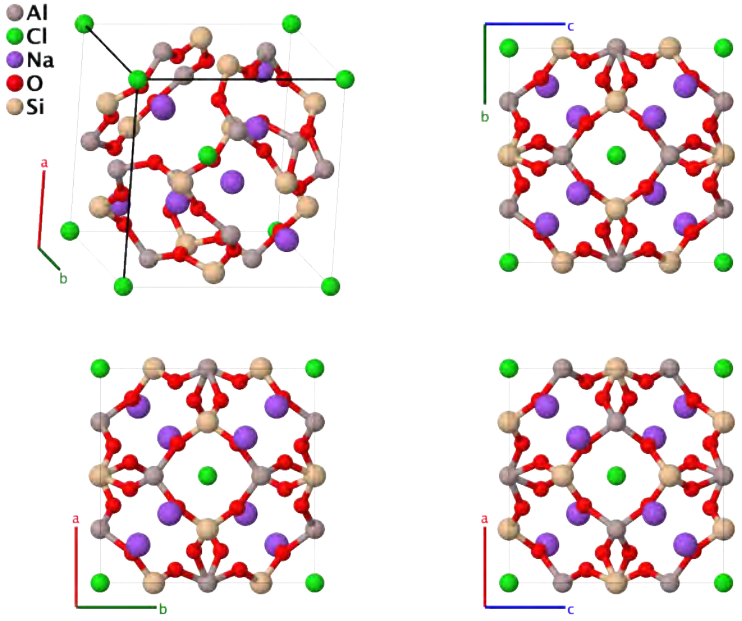
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1799
- POSCAR: pp. 1800

Sodalite [$Na_4(AlSiO_4)_3Cl$, $S6_2$] Structure: A3BC4D12E3_cP46_218_d_a_e_i_c

http://aflow.org/prototype-encyclopedia/A3BC4D12E3_cP46_218_d_a_e_i_c



Prototype	:	$Al_3ClNa_4O_{12}Si_3$
AFLOW prototype label	:	A3BC4D12E3_cP46_218_d_a_e_i_c
***Strukturbericht* designation**	:	$S6_2$
Pearson symbol	:	cP46
Space group number	:	218
Space group symbol	:	$P\bar{4}3n$
AFLOW prototype command	:	`aflow --proto=A3BC4D12E3_cP46_218_d_a_e_i_c --params=`a, x_4, x_5, y_5, z_5

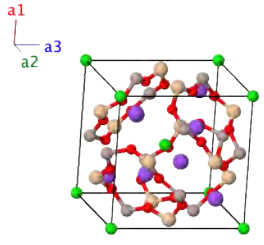
Other compounds with this structure

- $(Na_xK_{1-x})_4(AlSiO_4)_3Cl$, $Li_4(AlSiO_4)_3Cl$, $(Li_xK_{1-x})_4(AlSiO_4)_3Cl$, $Na_4(AlSiO_4)_3Br$, and $Na_4(AlSiO_4)_3I$

- Sodalites of the form $A_4(AlSiO_4)_3B$ with (A,B) = (Li,F), (Li,Br), (Li,I), (Na,F), (K,F), (K,Br) and (Rb,F) have been predicted to form, but not seen experimentally.

Simple Cubic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = a\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($2a$)	Cl
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	($2a$)	Cl
$\mathbf{B}_3$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$	($6c$)	Si
$\mathbf{B}_4$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$	($6c$)	Si
$\mathbf{B}_5$	=	$\frac{1}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	($6c$)	Si
$\mathbf{B}_6$	=	$\frac{3}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	($6c$)	Si
$\mathbf{B}_7$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	($6c$)	Si
$\mathbf{B}_8$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	($6c$)	Si
$\mathbf{B}_9$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	($6d$)	Al
$\mathbf{B}_{10}$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	($6d$)	Al
$\mathbf{B}_{11}$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}}$	($6d$)	Al
$\mathbf{B}_{12}$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}}$	($6d$)	Al
$\mathbf{B}_{13}$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	($6d$)	Al
$\mathbf{B}_{14}$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	($6d$)	Al
$\mathbf{B}_{15}$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	($8e$)	Na
$\mathbf{B}_{16}$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	($8e$)	Na
$\mathbf{B}_{17}$	=	$-x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} - x_4a\,\hat{\mathbf{z}}$	($8e$)	Na
$\mathbf{B}_{18}$	=	$x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} - x_4a\,\hat{\mathbf{z}}$	($8e$)	Na
$\mathbf{B}_{19}$	=	$\left(\frac{1}{2}+x_4\right)\mathbf{a}_1 + \left(\frac{1}{2}+x_4\right)\mathbf{a}_2 + \left(\frac{1}{2}+x_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{z}}$	($8e$)	Na
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2}-x_4\right)\mathbf{a}_1 + \left(\frac{1}{2}-x_4\right)\mathbf{a}_2 + \left(\frac{1}{2}+x_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{z}}$	($8e$)	Na
$\mathbf{B}_{21}$	=	$\left(\frac{1}{2}+x_4\right)\mathbf{a}_1 + \left(\frac{1}{2}-x_4\right)\mathbf{a}_2 + \left(\frac{1}{2}-x_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{z}}$	($8e$)	Na
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2}-x_4\right)\mathbf{a}_1 + \left(\frac{1}{2}+x_4\right)\mathbf{a}_2 + \left(\frac{1}{2}-x_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-x_4\right)a\,\hat{\mathbf{z}}$	($8e$)	Na
$\mathbf{B}_{23}$	=	$x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}} + y_5a\,\hat{\mathbf{y}} + z_5a\,\hat{\mathbf{z}}$	($24i$)	O
$\mathbf{B}_{24}$	=	$-x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}} - y_5a\,\hat{\mathbf{y}} + z_5a\,\hat{\mathbf{z}}$	($24i$)	O
$\mathbf{B}_{25}$	=	$-x_5\,\mathbf{a}_1 + y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}} + y_5a\,\hat{\mathbf{y}} - z_5a\,\hat{\mathbf{z}}$	($24i$)	O
$\mathbf{B}_{26}$	=	$x_5\,\mathbf{a}_1 - y_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}} - y_5a\,\hat{\mathbf{y}} - z_5a\,\hat{\mathbf{z}}$	($24i$)	O
$\mathbf{B}_{27}$	=	$z_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + y_5\,\mathbf{a}_3$	=	$z_5a\,\hat{\mathbf{x}} + x_5a\,\hat{\mathbf{y}} + y_5a\,\hat{\mathbf{z}}$	($24i$)	O
$\mathbf{B}_{28}$	=	$z_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - y_5\,\mathbf{a}_3$	=	$z_5a\,\hat{\mathbf{x}} - x_5a\,\hat{\mathbf{y}} - y_5a\,\hat{\mathbf{z}}$	($24i$)	O
$\mathbf{B}_{29}$	=	$-z_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + y_5\,\mathbf{a}_3$	=	$-z_5a\,\hat{\mathbf{x}} - x_5a\,\hat{\mathbf{y}} + y_5a\,\hat{\mathbf{z}}$	($24i$)	O
$\mathbf{B}_{30}$	=	$-z_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 - y_5\,\mathbf{a}_3$	=	$-z_5a\,\hat{\mathbf{x}} + x_5a\,\hat{\mathbf{y}} - y_5a\,\hat{\mathbf{z}}$	($24i$)	O
$\mathbf{B}_{31}$	=	$y_5\,\mathbf{a}_1 + z_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$y_5a\,\hat{\mathbf{x}} + z_5a\,\hat{\mathbf{y}} + x_5a\,\hat{\mathbf{z}}$	($24i$)	O

$\mathbf{B}_{32}$	=	$-y_5\,\mathbf{a}_1 + z_5\,\mathbf{a}_2 - x_5\,\mathbf{a}_3$	=	$-y_5 a\,\hat{\mathbf{x}} + z_5 a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}}$	(24*i*)	O
$\mathbf{B}_{33}$	=	$y_5\,\mathbf{a}_1 - z_5\,\mathbf{a}_2 - x_5\,\mathbf{a}_3$	=	$y_5 a\,\hat{\mathbf{x}} - z_5 a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}}$	(24*i*)	O
$\mathbf{B}_{34}$	=	$-y_5\,\mathbf{a}_1 - z_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$-y_5 a\,\hat{\mathbf{x}} - z_5 a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}}$	(24*i*)	O
$\mathbf{B}_{35}$	=	$\left(\frac{1}{2}+y_5\right)\mathbf{a}_1 + \left(\frac{1}{2}+x_5\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_5\right)a\,\hat{\mathbf{z}}$	(24*i*)	O
$\mathbf{B}_{36}$	=	$\left(\frac{1}{2}-y_5\right)\mathbf{a}_1 + \left(\frac{1}{2}-x_5\right)\mathbf{a}_2 + \left(\frac{1}{2}+z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+z_5\right)a\,\hat{\mathbf{z}}$	(24*i*)	O
$\mathbf{B}_{37}$	=	$\left(\frac{1}{2}+y_5\right)\mathbf{a}_1 + \left(\frac{1}{2}-x_5\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+y_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_5\right)a\,\hat{\mathbf{z}}$	(24*i*)	O
$\mathbf{B}_{38}$	=	$\left(\frac{1}{2}-y_5\right)\mathbf{a}_1 + \left(\frac{1}{2}+x_5\right)\mathbf{a}_2 + \left(\frac{1}{2}-z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-z_5\right)a\,\hat{\mathbf{z}}$	(24*i*)	O
$\mathbf{B}_{39}$	=	$\left(\frac{1}{2}+x_5\right)\mathbf{a}_1 + \left(\frac{1}{2}+z_5\right)\mathbf{a}_2 + \left(\frac{1}{2}+y_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+z_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+y_5\right)a\,\hat{\mathbf{z}}$	(24*i*)	O
$\mathbf{B}_{40}$	=	$\left(\frac{1}{2}-x_5\right)\mathbf{a}_1 + \left(\frac{1}{2}+z_5\right)\mathbf{a}_2 + \left(\frac{1}{2}-y_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+z_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-y_5\right)a\,\hat{\mathbf{z}}$	(24*i*)	O
$\mathbf{B}_{41}$	=	$\left(\frac{1}{2}-x_5\right)\mathbf{a}_1 + \left(\frac{1}{2}-z_5\right)\mathbf{a}_2 + \left(\frac{1}{2}+y_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-z_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+y_5\right)a\,\hat{\mathbf{z}}$	(24*i*)	O
$\mathbf{B}_{42}$	=	$\left(\frac{1}{2}+x_5\right)\mathbf{a}_1 + \left(\frac{1}{2}-z_5\right)\mathbf{a}_2 + \left(\frac{1}{2}-y_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-z_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-y_5\right)a\,\hat{\mathbf{z}}$	(24*i*)	O
$\mathbf{B}_{43}$	=	$\left(\frac{1}{2}+z_5\right)\mathbf{a}_1 + \left(\frac{1}{2}+y_5\right)\mathbf{a}_2 + \left(\frac{1}{2}+x_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+z_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{z}}$	(24*i*)	O
$\mathbf{B}_{44}$	=	$\left(\frac{1}{2}+z_5\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_5\right)\mathbf{a}_2 + \left(\frac{1}{2}-x_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+z_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{z}}$	(24*i*)	O
$\mathbf{B}_{45}$	=	$\left(\frac{1}{2}-z_5\right)\mathbf{a}_1 + \left(\frac{1}{2}+y_5\right)\mathbf{a}_2 + \left(\frac{1}{2}-x_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-z_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+y_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}-x_5\right)a\,\hat{\mathbf{z}}$	(24*i*)	O
$\mathbf{B}_{46}$	=	$\left(\frac{1}{2}-z_5\right)\mathbf{a}_1 + \left(\frac{1}{2}-y_5\right)\mathbf{a}_2 + \left(\frac{1}{2}+x_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-z_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}-y_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+x_5\right)a\,\hat{\mathbf{z}}$	(24*i*)	O

References:

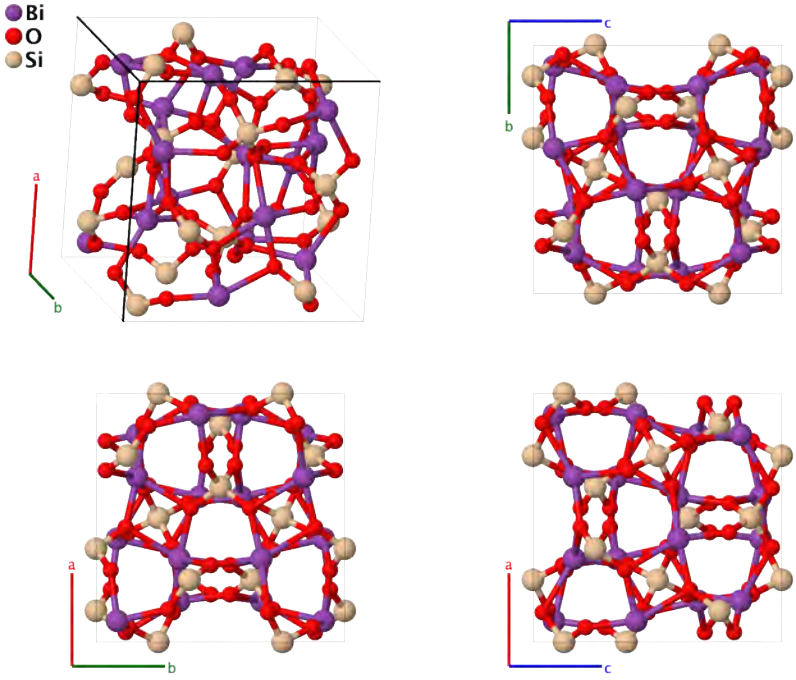
- I. Hassan and H. D. Grundy, *The Crystal Structures of Sodalite-Group Minerals*, Acta Crystallogr. Sect. B Struct. Sci. **40**, 6–13 (1984), doi:10.1107/S0108768184001683.

Geometry files:

- CIF: pp. 1800
- POSCAR: pp. 1800

Eulytine ($Bi_4(SiO_4)_3$, $S\,1_5$) Structure: A4B12C3_cI76_220_c_e_a

http://aflow.org/prototype-encyclopedia/A4B12C3_cI76_220_c_e_a



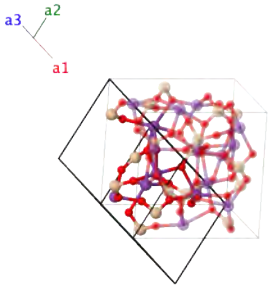
Prototype	:	$Bi_4O_{12}Si_3$
AFLOW prototype label	:	A4B12C3_cI76_220_c_e_a
***Strukturbericht* designation**	:	$S\,1_5$
Pearson symbol	:	cI76
Space group number	:	220
Space group symbol	:	$I\bar{4}3d$
AFLOW prototype command	:	`aflow --proto=A4B12C3_cI76_220_c_e_a --params=`a, x_2, x_3, y_3, z_3

Other compounds with this structure

- $Bi_4(GeO_4)_3$, $Ca_3Bi(PO_4)_3$, and $M_3L(PO_4)_3$ (M = Sr, Ba), L = (La, Nd, Gd, Y, Lu, In, Bi)

Body-centered Cubic primitive vectors:

$$\mathbf{a}_1 = -\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}a\,\hat{\mathbf{x}} - \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} - \tfrac{1}{2}a\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{5}{8}\,\mathbf{a}_2 + \frac{3}{8}\,\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(12*a*)	Si
$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{7}{8}\,\mathbf{a}_2 + \frac{1}{8}\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(12*a*)	Si
$\mathbf{B}_3$	=	$\frac{3}{8}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{5}{8}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}}$	(12*a*)	Si
$\mathbf{B}_4$	=	$\frac{1}{8}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{7}{8}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}}$	(12*a*)	Si
$\mathbf{B}_5$	=	$\frac{5}{8}\,\mathbf{a}_1 + \frac{3}{8}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}}$	(12*a*)	Si
$\mathbf{B}_6$	=	$\frac{7}{8}\,\mathbf{a}_1 + \frac{1}{8}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(12*a*)	Si
$\mathbf{B}_7$	=	$2x_2\,\mathbf{a}_1 + 2x_2\,\mathbf{a}_2 + 2x_2\,\mathbf{a}_3$	=	$x_2a\,\hat{\mathbf{x}} + x_2a\,\hat{\mathbf{y}} + x_2a\,\hat{\mathbf{z}}$	(16*c*)	Bi
$\mathbf{B}_8$	=	$\frac{1}{2}\,\mathbf{a}_1 + \left(\frac{1}{2} - 2x_2\right)\,\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{y}} + x_2a\,\hat{\mathbf{z}}$	(16*c*)	Bi
$\mathbf{B}_9$	=	$\left(\frac{1}{2} - 2x_2\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{x}} + x_2a\,\hat{\mathbf{y}} - x_2a\,\hat{\mathbf{z}}$	(16*c*)	Bi
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - 2x_2\right)\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$x_2a\,\hat{\mathbf{x}} - x_2a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_2\right)a\,\hat{\mathbf{z}}$	(16*c*)	Bi
$\mathbf{B}_{11}$	=	$\left(\frac{1}{2} + 2x_2\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + 2x_2\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + 2x_2\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_2\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + x_2\right)a\,\hat{\mathbf{z}}$	(16*c*)	Bi
$\mathbf{B}_{12}$	=	$\frac{1}{2}\,\mathbf{a}_1 - 2x_2\,\mathbf{a}_3$	=	$-\left(x_2 + \frac{1}{4}\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_2\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + x_2\right)a\,\hat{\mathbf{z}}$	(16*c*)	Bi
$\mathbf{B}_{13}$	=	$-2x_2\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\left(\frac{1}{4} + x_2\right)a\,\hat{\mathbf{x}} - \left(x_2 + \frac{1}{4}\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_2\right)a\,\hat{\mathbf{z}}$	(16*c*)	Bi
$\mathbf{B}_{14}$	=	$-2x_2\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_2\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_2\right)a\,\hat{\mathbf{y}} - \left(x_2 + \frac{1}{4}\right)a\,\hat{\mathbf{z}}$	(16*c*)	Bi
$\mathbf{B}_{15}$	=	$(y_3 + z_3)\,\mathbf{a}_1 + (x_3 + z_3)\,\mathbf{a}_2 + (x_3 + y_3)\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + y_3a\,\hat{\mathbf{y}} + z_3a\,\hat{\mathbf{z}}$	(48*e*)	O
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} - y_3 + z_3\right)\,\mathbf{a}_1 + (-x_3 + z_3)\,\mathbf{a}_2 + \left(\frac{1}{2} - x_3 - y_3\right)\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{y}} + z_3a\,\hat{\mathbf{z}}$	(48*e*)	O
$\mathbf{B}_{17}$	=	$(y_3 - z_3)\,\mathbf{a}_1 + \left(\frac{1}{2} - x_3 - z_3\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - x_3 + y_3\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + y_3a\,\hat{\mathbf{y}} - z_3a\,\hat{\mathbf{z}}$	(48*e*)	O
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} - y_3 - z_3\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + x_3 - z_3\right)\,\mathbf{a}_2 + (x_3 - y_3)\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} - y_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_3\right)a\,\hat{\mathbf{z}}$	(48*e*)	O
$\mathbf{B}_{19}$	=	$(x_3 + y_3)\,\mathbf{a}_1 + (y_3 + z_3)\,\mathbf{a}_2 + (x_3 + z_3)\,\mathbf{a}_3$	=	$z_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + y_3a\,\hat{\mathbf{z}}$	(48*e*)	O
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} - x_3 - y_3\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - y_3 + z_3\right)\,\mathbf{a}_2 + (-x_3 + z_3)\,\mathbf{a}_3$	=	$z_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{z}}$	(48*e*)	O
$\mathbf{B}_{21}$	=	$\left(\frac{1}{2} - x_3 + y_3\right)\,\mathbf{a}_1 + (y_3 - z_3)\,\mathbf{a}_2 + \left(\frac{1}{2} - x_3 - z_3\right)\,\mathbf{a}_3$	=	$-z_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} + y_3a\,\hat{\mathbf{z}}$	(48*e*)	O
$\mathbf{B}_{22}$	=	$(x_3 - y_3)\,\mathbf{a}_1 + \left(\frac{1}{2} - y_3 - z_3\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + x_3 - z_3\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - z_3\right)a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} - y_3a\,\hat{\mathbf{z}}$	(48*e*)	O
$\mathbf{B}_{23}$	=	$(x_3 + z_3)\,\mathbf{a}_1 + (x_3 + y_3)\,\mathbf{a}_2 + (y_3 + z_3)\,\mathbf{a}_3$	=	$y_3a\,\hat{\mathbf{x}} + z_3a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(48*e*)	O
$\mathbf{B}_{24}$	=	$(-x_3 + z_3)\,\mathbf{a}_1 + \left(\frac{1}{2} - x_3 - y_3\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - y_3 + z_3\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - y_3\right)a\,\hat{\mathbf{x}} + z_3a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(48*e*)	O
$\mathbf{B}_{25}$	=	$\left(\frac{1}{2} - x_3 - z_3\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - x_3 + y_3\right)\,\mathbf{a}_2 + (y_3 - z_3)\,\mathbf{a}_3$	=	$y_3a\,\hat{\mathbf{x}} - z_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{z}}$	(48*e*)	O

$\mathbf{B}_{26}$	=	$\left(\frac{1}{2}+x_3-z_3\right)\mathbf{a}_1+(x_3-y_3)\,\mathbf{a}_2+\left(\frac{1}{2}-y_3-z_3\right)\mathbf{a}_3$	=	$-y_3a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-z_3\right)a\,\hat{\mathbf{y}}+x_3a\,\hat{\mathbf{z}}$	(48*e*)	O
$\mathbf{B}_{27}$	=	$\left(\frac{1}{2}+x_3+z_3\right)\mathbf{a}_1+\left(\frac{1}{2}+y_3+z_3\right)\mathbf{a}_2+\left(\frac{1}{2}+x_3+y_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}+y_3\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+x_3\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+z_3\right)a\,\hat{\mathbf{z}}$	(48*e*)	O
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2}-x_3+z_3\right)\mathbf{a}_1+(-y_3+z_3)\,\mathbf{a}_2+(-x_3-y_3)\,\mathbf{a}_3$	=	$-\left(y_3+\frac{1}{4}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-x_3\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+z_3\right)a\,\hat{\mathbf{z}}$	(48*e*)	O
$\mathbf{B}_{29}$	=	$(-x_3-z_3)\,\mathbf{a}_1+\left(\frac{1}{2}+y_3-z_3\right)\mathbf{a}_2+(-x_3+y_3)\,\mathbf{a}_3$	=	$\left(\frac{1}{4}+y_3\right)a\,\hat{\mathbf{x}}-\left(x_3+\frac{1}{4}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-z_3\right)a\,\hat{\mathbf{z}}$	(48*e*)	O
$\mathbf{B}_{30}$	=	$(x_3-z_3)\,\mathbf{a}_1+(-y_3-z_3)\,\mathbf{a}_2+\left(\frac{1}{2}+x_3-y_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}-y_3\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+x_3\right)a\,\hat{\mathbf{y}}-\left(z_3+\frac{1}{4}\right)a\,\hat{\mathbf{z}}$	(48*e*)	O
$\mathbf{B}_{31}$	=	$\left(\frac{1}{2}+y_3+z_3\right)\mathbf{a}_1+\left(\frac{1}{2}+x_3+y_3\right)\mathbf{a}_2+\left(\frac{1}{2}+x_3+z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}+x_3\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+z_3\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+y_3\right)a\,\hat{\mathbf{z}}$	(48*e*)	O
$\mathbf{B}_{32}$	=	$(-y_3+z_3)\,\mathbf{a}_1+(-x_3-y_3)\,\mathbf{a}_2+\left(\frac{1}{2}-x_3+z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}-x_3\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+z_3\right)a\,\hat{\mathbf{y}}-\left(y_3+\frac{1}{4}\right)a\,\hat{\mathbf{z}}$	(48*e*)	O
$\mathbf{B}_{33}$	=	$\left(\frac{1}{2}+y_3-z_3\right)\mathbf{a}_1+(-x_3+y_3)\,\mathbf{a}_2+(-x_3-z_3)\,\mathbf{a}_3$	=	$-\left(x_3+\frac{1}{4}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-z_3\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+y_3\right)a\,\hat{\mathbf{z}}$	(48*e*)	O
$\mathbf{B}_{34}$	=	$(-y_3-z_3)\,\mathbf{a}_1+\left(\frac{1}{2}+x_3-y_3\right)\mathbf{a}_2+(x_3-z_3)\,\mathbf{a}_3$	=	$\left(\frac{1}{4}+x_3\right)a\,\hat{\mathbf{x}}-\left(z_3+\frac{1}{4}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-y_3\right)a\,\hat{\mathbf{z}}$	(48*e*)	O
$\mathbf{B}_{35}$	=	$\left(\frac{1}{2}+x_3+y_3\right)\mathbf{a}_1+\left(\frac{1}{2}+x_3+z_3\right)\mathbf{a}_2+\left(\frac{1}{2}+y_3+z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}+z_3\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+y_3\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+x_3\right)a\,\hat{\mathbf{z}}$	(48*e*)	O
$\mathbf{B}_{36}$	=	$(-x_3-y_3)\,\mathbf{a}_1+\left(\frac{1}{2}-x_3+z_3\right)\mathbf{a}_2+(-y_3+z_3)\,\mathbf{a}_3$	=	$\left(\frac{1}{4}+z_3\right)a\,\hat{\mathbf{x}}-\left(y_3+\frac{1}{4}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-x_3\right)a\,\hat{\mathbf{z}}$	(48*e*)	O
$\mathbf{B}_{37}$	=	$(-x_3+y_3)\,\mathbf{a}_1+(-x_3-z_3)\,\mathbf{a}_2+\left(\frac{1}{2}+y_3-z_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}-z_3\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+y_3\right)a\,\hat{\mathbf{y}}-\left(x_3+\frac{1}{4}\right)a\,\hat{\mathbf{z}}$	(48*e*)	O
$\mathbf{B}_{38}$	=	$\left(\frac{1}{2}+x_3-y_3\right)\mathbf{a}_1+(x_3-z_3)\,\mathbf{a}_2+(-y_3-z_3)\,\mathbf{a}_3$	=	$-\left(z_3+\frac{1}{4}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-y_3\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+x_3\right)a\,\hat{\mathbf{z}}$	(48*e*)	O

References:

- H. Liu and C. Kuo, *Crystal structure of bismuth(III) silicate, $Bi_4(SiO_4)_3$*, Zeitschrift für Kristallographie - Crystalline Materials **212**, 48 (1997), doi:10.1524/zkri.1997.212.1.48.

Found in:

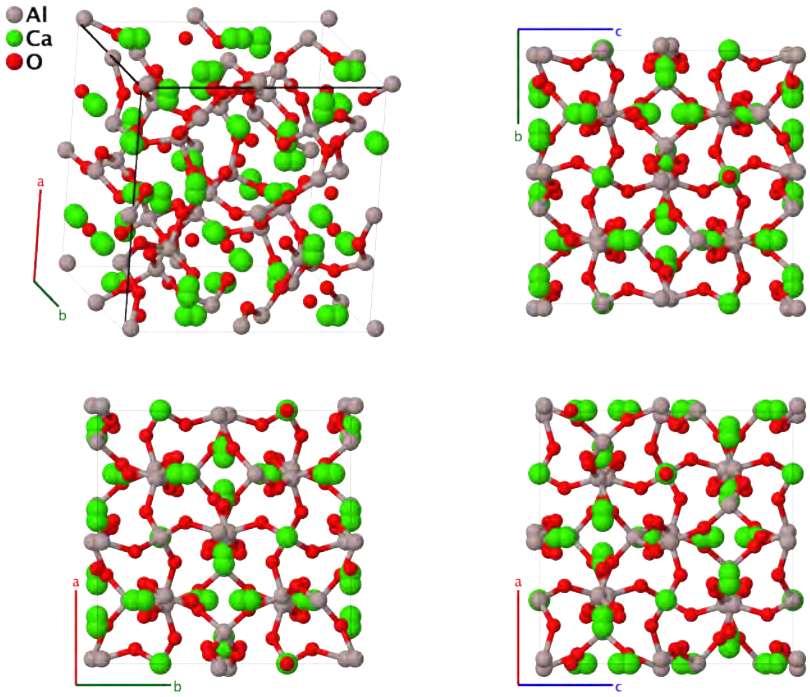
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1801
- POSCAR: pp. 1801

Mayenite ($12CaO{\cdot}7Al_2O_3$, $K7_4$, C12A7) Structure: A7B12C19_cI152_220_bc_2d_ace

http://aflow.org/prototype-encyclopedia/A7B12C19_cI152_220_bc_2d_ace

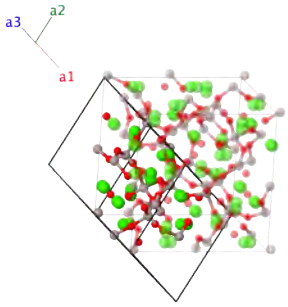


Prototype	:	$Al_{14}Ca_{12}O_{33}$
AFLOW prototype label	:	A7B12C19_cI152_220_bc_2d_ace
***Strukturbericht* designation**	:	$K7_4$
Pearson symbol	:	cI152
Space group number	:	220
Space group symbol	:	$I\bar{4}3d$
AFLOW prototype command	:	`aflow --proto=A7B12C19_cI152_220_bc_2d_ace --params=`$a, x_3, x_4, x_5, x_6, x_7, y_7, z_7$

- We present the structure determined by (Boysen, 2007) with data taken at 293 K. This slightly differs from the original determination of (Büssem, 1936), which was given the $K7_4$ designation by (Gottfried, 1938). In the original work, the calcium atoms were thought to be located at a single (24*d*) site. Newer findings show that calcium is split between two (24*d*) sites, with the site we have labeled Ca-I having 87.5% of the atoms and Ca-II the remainder, although presumably only one of the two sites is occupied in any pair.
- In all works the O-I (12*a*) site is only partially occupied: if this is occupied 1/6 of the time, we get the proper stoichiometry, though (Boysen, 2007) found the occupation was 0.251 at 293 K, dropping as the temperature decreased.
- This structure is often referred to in the literature as C12A7, to distinguish it from other CaO/Al_2O_3 compounds.

Body-centered Cubic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= -\tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{y}} + \tfrac{1}{2}\,a\,\hat{\mathbf{z}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} - \tfrac{1}{2}\,a\,\hat{\mathbf{y}} + \tfrac{1}{2}\,a\,\hat{\mathbf{z}}\\ \mathbf{a}_3 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{y}} - \tfrac{1}{2}\,a\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\tfrac{1}{4}\,\mathbf{a}_1 + \tfrac{5}{8}\,\mathbf{a}_2 + \tfrac{3}{8}\,\mathbf{a}_3$	=	$\tfrac{3}{8}a\,\hat{\mathbf{x}} + \tfrac{1}{4}a\,\hat{\mathbf{z}}$	(12*a*)	O I
$\mathbf{B}_2$	=	$\tfrac{3}{4}\,\mathbf{a}_1 + \tfrac{7}{8}\,\mathbf{a}_2 + \tfrac{1}{8}\,\mathbf{a}_3$	=	$\tfrac{1}{8}a\,\hat{\mathbf{x}} + \tfrac{3}{4}a\,\hat{\mathbf{z}}$	(12*a*)	O I
$\mathbf{B}_3$	=	$\tfrac{3}{8}\,\mathbf{a}_1 + \tfrac{1}{4}\,\mathbf{a}_2 + \tfrac{5}{8}\,\mathbf{a}_3$	=	$\tfrac{1}{4}a\,\hat{\mathbf{x}} + \tfrac{3}{8}a\,\hat{\mathbf{y}}$	(12*a*)	O I
$\mathbf{B}_4$	=	$\tfrac{1}{8}\,\mathbf{a}_1 + \tfrac{3}{4}\,\mathbf{a}_2 + \tfrac{7}{8}\,\mathbf{a}_3$	=	$\tfrac{3}{4}a\,\hat{\mathbf{x}} + \tfrac{1}{8}a\,\hat{\mathbf{y}}$	(12*a*)	O I
$\mathbf{B}_5$	=	$\tfrac{5}{8}\,\mathbf{a}_1 + \tfrac{3}{8}\,\mathbf{a}_2 + \tfrac{1}{4}\,\mathbf{a}_3$	=	$\tfrac{1}{4}a\,\hat{\mathbf{y}} + \tfrac{3}{8}a\,\hat{\mathbf{z}}$	(12*a*)	O I
$\mathbf{B}_6$	=	$\tfrac{7}{8}\,\mathbf{a}_1 + \tfrac{1}{8}\,\mathbf{a}_2 + \tfrac{3}{4}\,\mathbf{a}_3$	=	$\tfrac{3}{4}a\,\hat{\mathbf{y}} + \tfrac{1}{8}a\,\hat{\mathbf{z}}$	(12*a*)	O I
$\mathbf{B}_7$	=	$\tfrac{1}{4}\,\mathbf{a}_1 + \tfrac{1}{8}\,\mathbf{a}_2 + \tfrac{7}{8}\,\mathbf{a}_3$	=	$\tfrac{3}{8}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} - \tfrac{1}{4}a\,\hat{\mathbf{z}}$	(12*b*)	Al I
$\mathbf{B}_8$	=	$\tfrac{3}{4}\,\mathbf{a}_1 + \tfrac{3}{8}\,\mathbf{a}_2 + \tfrac{5}{8}\,\mathbf{a}_3$	=	$\tfrac{1}{8}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{4}a\,\hat{\mathbf{z}}$	(12*b*)	Al I
$\mathbf{B}_9$	=	$\tfrac{7}{8}\,\mathbf{a}_1 + \tfrac{1}{4}\,\mathbf{a}_2 + \tfrac{1}{8}\,\mathbf{a}_3$	=	$-\tfrac{1}{4}a\,\hat{\mathbf{x}} + \tfrac{3}{8}a\,\hat{\mathbf{y}} + \tfrac{1}{2}a\,\hat{\mathbf{z}}$	(12*b*)	Al I
$\mathbf{B}_{10}$	=	$\tfrac{5}{8}\,\mathbf{a}_1 + \tfrac{3}{4}\,\mathbf{a}_2 + \tfrac{3}{8}\,\mathbf{a}_3$	=	$\tfrac{1}{4}a\,\hat{\mathbf{x}} + \tfrac{1}{8}a\,\hat{\mathbf{y}} + \tfrac{1}{2}a\,\hat{\mathbf{z}}$	(12*b*)	Al I
$\mathbf{B}_{11}$	=	$\tfrac{1}{8}\,\mathbf{a}_1 + \tfrac{7}{8}\,\mathbf{a}_2 + \tfrac{1}{4}\,\mathbf{a}_3$	=	$\tfrac{1}{2}a\,\hat{\mathbf{x}} - \tfrac{1}{4}a\,\hat{\mathbf{y}} + \tfrac{3}{8}a\,\hat{\mathbf{z}}$	(12*b*)	Al I
$\mathbf{B}_{12}$	=	$\tfrac{3}{8}\,\mathbf{a}_1 + \tfrac{5}{8}\,\mathbf{a}_2 + \tfrac{3}{4}\,\mathbf{a}_3$	=	$\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{4}a\,\hat{\mathbf{y}} + \tfrac{1}{8}a\,\hat{\mathbf{z}}$	(12*b*)	Al I
$\mathbf{B}_{13}$	=	$2x_3\,\mathbf{a}_1 + 2x_3\,\mathbf{a}_2 + 2x_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(16*c*)	Al II
$\mathbf{B}_{14}$	=	$\tfrac{1}{2}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - 2x_3\right)\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(16*c*)	Al II
$\mathbf{B}_{15}$	=	$\left(\tfrac{1}{2} - 2x_3\right)\mathbf{a}_2 + \tfrac{1}{2}\,\mathbf{a}_3$	=	$\left(\tfrac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(16*c*)	Al II
$\mathbf{B}_{16}$	=	$\left(\tfrac{1}{2} - 2x_3\right)\mathbf{a}_1 + \tfrac{1}{2}\,\mathbf{a}_2$	=	$x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_3\right)a\,\hat{\mathbf{z}}$	(16*c*)	Al II
$\mathbf{B}_{17}$	=	$\left(\tfrac{1}{2} + 2x_3\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + 2x_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + 2x_3\right)\mathbf{a}_3$	=	$\left(\tfrac{1}{4} + x_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_3\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + x_3\right)a\,\hat{\mathbf{z}}$	(16*c*)	Al II
$\mathbf{B}_{18}$	=	$\tfrac{1}{2}\,\mathbf{a}_1 - 2x_3\,\mathbf{a}_3$	=	$-\left(x_3 + \tfrac{1}{4}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_3\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + x_3\right)a\,\hat{\mathbf{z}}$	(16*c*)	Al II
$\mathbf{B}_{19}$	=	$-2x_3\,\mathbf{a}_1 + \tfrac{1}{2}\,\mathbf{a}_2$	=	$\left(\tfrac{1}{4} + x_3\right)a\,\hat{\mathbf{x}} - \left(x_3 + \tfrac{1}{4}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - x_3\right)a\,\hat{\mathbf{z}}$	(16*c*)	Al II
$\mathbf{B}_{20}$	=	$-2x_3\,\mathbf{a}_2 + \tfrac{1}{2}\,\mathbf{a}_3$	=	$\left(\tfrac{1}{4} - x_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_3\right)a\,\hat{\mathbf{y}} - \left(x_3 + \tfrac{1}{4}\right)a\,\hat{\mathbf{z}}$	(16*c*)	Al II
$\mathbf{B}_{21}$	=	$2x_4\,\mathbf{a}_1 + 2x_4\,\mathbf{a}_2 + 2x_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	(16*c*)	O II
$\mathbf{B}_{22}$	=	$\tfrac{1}{2}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - 2x_4\right)\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_4\right)a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	(16*c*)	O II
$\mathbf{B}_{23}$	=	$\left(\tfrac{1}{2} - 2x_4\right)\mathbf{a}_2 + \tfrac{1}{2}\,\mathbf{a}_3$	=	$\left(\tfrac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} - x_4a\,\hat{\mathbf{z}}$	(16*c*)	O II
$\mathbf{B}_{24}$	=	$\left(\tfrac{1}{2} - 2x_4\right)\mathbf{a}_1 + \tfrac{1}{2}\,\mathbf{a}_2$	=	$x_4a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_4\right)a\,\hat{\mathbf{z}}$	(16*c*)	O II

$\mathbf{B}_{25}$	=	$\left(\frac{1}{2}+2x_4\right)\mathbf{a}_1+\left(\frac{1}{2}+2x_4\right)\mathbf{a}_2+\left(\frac{1}{2}+2x_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}+x_4\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+x_4\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+x_4\right)a\,\hat{\mathbf{z}}$	(16*c*)	O II
$\mathbf{B}_{26}$	=	$\frac{1}{2}\mathbf{a}_1-2x_4\,\mathbf{a}_3$	=	$-\left(x_4+\frac{1}{4}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-x_4\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+x_4\right)a\,\hat{\mathbf{z}}$	(16*c*)	O II
$\mathbf{B}_{27}$	=	$-2x_4\,\mathbf{a}_1+\frac{1}{2}\mathbf{a}_2$	=	$\left(\frac{1}{4}+x_4\right)a\,\hat{\mathbf{x}}-\left(x_4+\frac{1}{4}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-x_4\right)a\,\hat{\mathbf{z}}$	(16*c*)	O II
$\mathbf{B}_{28}$	=	$-2x_4\,\mathbf{a}_2+\frac{1}{2}\mathbf{a}_3$	=	$\left(\frac{1}{4}-x_4\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+x_4\right)a\,\hat{\mathbf{y}}-\left(x_4+\frac{1}{4}\right)a\,\hat{\mathbf{z}}$	(16*c*)	O II
$\mathbf{B}_{29}$	=	$\frac{1}{4}\mathbf{a}_1+\left(\frac{1}{4}+x_5\right)\mathbf{a}_2+x_5\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}}+\frac{1}{4}a\,\hat{\mathbf{z}}$	(24*d*)	Ca I
$\mathbf{B}_{30}$	=	$\frac{3}{4}\mathbf{a}_1+\left(\frac{1}{4}-x_5\right)\mathbf{a}_2+\left(\frac{1}{2}-x_5\right)\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}}+\frac{1}{2}a\,\hat{\mathbf{y}}+\frac{1}{4}a\,\hat{\mathbf{z}}$	(24*d*)	Ca I
$\mathbf{B}_{31}$	=	$x_5\,\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+\left(\frac{1}{4}+x_5\right)\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}}+x_5a\,\hat{\mathbf{y}}$	(24*d*)	Ca I
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2}-x_5\right)\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2+\left(\frac{1}{4}-x_5\right)\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}}-x_5a\,\hat{\mathbf{y}}+\frac{1}{2}a\,\hat{\mathbf{z}}$	(24*d*)	Ca I
$\mathbf{B}_{33}$	=	$\left(\frac{1}{4}+x_5\right)\mathbf{a}_1+x_5\,\mathbf{a}_2+\frac{1}{4}\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{y}}+x_5a\,\hat{\mathbf{z}}$	(24*d*)	Ca I
$\mathbf{B}_{34}$	=	$\left(\frac{1}{4}-x_5\right)\mathbf{a}_1+\left(\frac{1}{2}-x_5\right)\mathbf{a}_2+\frac{3}{4}\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}+\frac{1}{4}a\,\hat{\mathbf{y}}-x_5a\,\hat{\mathbf{z}}$	(24*d*)	Ca I
$\mathbf{B}_{35}$	=	$\left(\frac{3}{4}+x_5\right)\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2+\left(\frac{1}{2}+x_5\right)\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+x_5\right)a\,\hat{\mathbf{y}}+\frac{1}{2}a\,\hat{\mathbf{z}}$	(24*d*)	Ca I
$\mathbf{B}_{36}$	=	$\left(\frac{3}{4}-x_5\right)\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2-x_5\,\mathbf{a}_3$	=	$-\frac{1}{4}a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-x_5\right)a\,\hat{\mathbf{y}}+\frac{1}{2}a\,\hat{\mathbf{z}}$	(24*d*)	Ca I
$\mathbf{B}_{37}$	=	$\frac{3}{4}\mathbf{a}_1+\left(\frac{1}{2}+x_5\right)\mathbf{a}_2+\left(\frac{3}{4}+x_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}+x_5\right)a\,\hat{\mathbf{x}}+\frac{1}{2}a\,\hat{\mathbf{y}}+\frac{1}{4}a\,\hat{\mathbf{z}}$	(24*d*)	Ca I
$\mathbf{B}_{38}$	=	$\frac{1}{4}\mathbf{a}_1-x_5\,\mathbf{a}_2+\left(\frac{3}{4}-x_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}-x_5\right)a\,\hat{\mathbf{x}}+\frac{1}{2}a\,\hat{\mathbf{y}}-\frac{1}{4}a\,\hat{\mathbf{z}}$	(24*d*)	Ca I
$\mathbf{B}_{39}$	=	$\left(\frac{1}{2}+x_5\right)\mathbf{a}_1+\left(\frac{3}{4}+x_5\right)\mathbf{a}_2+\frac{3}{4}\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}+\frac{1}{4}a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+x_5\right)a\,\hat{\mathbf{z}}$	(24*d*)	Ca I
$\mathbf{B}_{40}$	=	$-x_5\,\mathbf{a}_1+\left(\frac{3}{4}-x_5\right)\mathbf{a}_2+\frac{1}{4}\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}-\frac{1}{4}a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-x_5\right)a\,\hat{\mathbf{z}}$	(24*d*)	Ca I
$\mathbf{B}_{41}$	=	$\frac{1}{4}\mathbf{a}_1+\left(\frac{1}{4}+x_6\right)\mathbf{a}_2+x_6\,\mathbf{a}_3$	=	$x_6a\,\hat{\mathbf{x}}+\frac{1}{4}a\,\hat{\mathbf{z}}$	(24*d*)	Ca II
$\mathbf{B}_{42}$	=	$\frac{3}{4}\mathbf{a}_1+\left(\frac{1}{4}-x_6\right)\mathbf{a}_2+\left(\frac{1}{2}-x_6\right)\mathbf{a}_3$	=	$-x_6a\,\hat{\mathbf{x}}+\frac{1}{2}a\,\hat{\mathbf{y}}+\frac{1}{4}a\,\hat{\mathbf{z}}$	(24*d*)	Ca II
$\mathbf{B}_{43}$	=	$x_6\,\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2+\left(\frac{1}{4}+x_6\right)\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}}+x_6a\,\hat{\mathbf{y}}$	(24*d*)	Ca II
$\mathbf{B}_{44}$	=	$\left(\frac{1}{2}-x_6\right)\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2+\left(\frac{1}{4}-x_6\right)\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}}-x_6a\,\hat{\mathbf{y}}+\frac{1}{2}a\,\hat{\mathbf{z}}$	(24*d*)	Ca II
$\mathbf{B}_{45}$	=	$\left(\frac{1}{4}+x_6\right)\mathbf{a}_1+x_6\,\mathbf{a}_2+\frac{1}{4}\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{y}}+x_6a\,\hat{\mathbf{z}}$	(24*d*)	Ca II
$\mathbf{B}_{46}$	=	$\left(\frac{1}{4}-x_6\right)\mathbf{a}_1+\left(\frac{1}{2}-x_6\right)\mathbf{a}_2+\frac{3}{4}\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}+\frac{1}{4}a\,\hat{\mathbf{y}}-x_6a\,\hat{\mathbf{z}}$	(24*d*)	Ca II
$\mathbf{B}_{47}$	=	$\left(\frac{3}{4}+x_6\right)\mathbf{a}_1+\frac{3}{4}\mathbf{a}_2+\left(\frac{1}{2}+x_6\right)\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+x_6\right)a\,\hat{\mathbf{y}}+\frac{1}{2}a\,\hat{\mathbf{z}}$	(24*d*)	Ca II
$\mathbf{B}_{48}$	=	$\left(\frac{3}{4}-x_6\right)\mathbf{a}_1+\frac{1}{4}\mathbf{a}_2-x_6\,\mathbf{a}_3$	=	$-\frac{1}{4}a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-x_6\right)a\,\hat{\mathbf{y}}+\frac{1}{2}a\,\hat{\mathbf{z}}$	(24*d*)	Ca II
$\mathbf{B}_{49}$	=	$\frac{3}{4}\mathbf{a}_1+\left(\frac{1}{2}+x_6\right)\mathbf{a}_2+\left(\frac{3}{4}+x_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}+x_6\right)a\,\hat{\mathbf{x}}+\frac{1}{2}a\,\hat{\mathbf{y}}+\frac{1}{4}a\,\hat{\mathbf{z}}$	(24*d*)	Ca II
$\mathbf{B}_{50}$	=	$\frac{1}{4}\mathbf{a}_1-x_6\,\mathbf{a}_2+\left(\frac{3}{4}-x_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}-x_6\right)a\,\hat{\mathbf{x}}+\frac{1}{2}a\,\hat{\mathbf{y}}-\frac{1}{4}a\,\hat{\mathbf{z}}$	(24*d*)	Ca II
$\mathbf{B}_{51}$	=	$\left(\frac{1}{2}+x_6\right)\mathbf{a}_1+\left(\frac{3}{4}+x_6\right)\mathbf{a}_2+\frac{3}{4}\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}+\frac{1}{4}a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+x_6\right)a\,\hat{\mathbf{z}}$	(24*d*)	Ca II
$\mathbf{B}_{52}$	=	$-x_6\,\mathbf{a}_1+\left(\frac{3}{4}-x_6\right)\mathbf{a}_2+\frac{1}{4}\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}-\frac{1}{4}a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-x_6\right)a\,\hat{\mathbf{z}}$	(24*d*)	Ca II
$\mathbf{B}_{53}$	=	$(y_7+z_7)\,\mathbf{a}_1+(x_7+z_7)\,\mathbf{a}_2+(x_7+y_7)\,\mathbf{a}_3$	=	$x_7a\,\hat{\mathbf{x}}+y_7a\,\hat{\mathbf{y}}+z_7a\,\hat{\mathbf{z}}$	(48*e*)	O III
$\mathbf{B}_{54}$	=	$\left(\frac{1}{2}-y_7+z_7\right)\mathbf{a}_1+(-x_7+z_7)\,\mathbf{a}_2+\left(\frac{1}{2}-x_7-y_7\right)\mathbf{a}_3$	=	$-x_7a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_7\right)a\,\hat{\mathbf{y}}+z_7a\,\hat{\mathbf{z}}$	(48*e*)	O III
$\mathbf{B}_{55}$	=	$(y_7-z_7)\,\mathbf{a}_1+\left(\frac{1}{2}-x_7-z_7\right)\mathbf{a}_2+\left(\frac{1}{2}-x_7+y_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_7\right)a\,\hat{\mathbf{x}}+y_7a\,\hat{\mathbf{y}}-z_7a\,\hat{\mathbf{z}}$	(48*e*)	O III

$$\mathbf{B}_{56} = \left(\tfrac{1}{2} - y_7 - z_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_7 - z_7\right)\mathbf{a}_2 + (x_7 - y_7)\,\mathbf{a}_3 = x_7 a\,\hat{\mathbf{x}} - y_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_7\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{57} = (x_7 + y_7)\,\mathbf{a}_1 + (y_7 + z_7)\,\mathbf{a}_2 + (x_7 + z_7)\,\mathbf{a}_3 = z_7 a\,\hat{\mathbf{x}} + x_7 a\,\hat{\mathbf{y}} + y_7 a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{58} = \left(\tfrac{1}{2} - x_7 - y_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_7 + z_7\right)\mathbf{a}_2 + (-x_7 + z_7)\,\mathbf{a}_3 = z_7 a\,\hat{\mathbf{x}} - x_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - y_7\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{59} = \left(\tfrac{1}{2} - x_7 + y_7\right)\mathbf{a}_1 + (y_7 - z_7)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_7 - z_7\right)\mathbf{a}_3 = -z_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_7\right)a\,\hat{\mathbf{y}} + y_7 a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{60} = (x_7 - y_7)\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_7 - z_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_7 - z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - z_7\right)a\,\hat{\mathbf{x}} + x_7 a\,\hat{\mathbf{y}} - y_7 a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{61} = (x_7 + z_7)\,\mathbf{a}_1 + (x_7 + y_7)\,\mathbf{a}_2 + (y_7 + z_7)\,\mathbf{a}_3 = y_7 a\,\hat{\mathbf{x}} + z_7 a\,\hat{\mathbf{y}} + x_7 a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{62} = (-x_7 + z_7)\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_7 - y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - y_7 + z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - y_7\right)a\,\hat{\mathbf{x}} + z_7 a\,\hat{\mathbf{y}} - x_7 a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{63} = \left(\tfrac{1}{2} - x_7 - z_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_7 + y_7\right)\mathbf{a}_2 + (y_7 - z_7)\,\mathbf{a}_3 = y_7 a\,\hat{\mathbf{x}} - z_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_7\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{64} = \left(\tfrac{1}{2} + x_7 - z_7\right)\mathbf{a}_1 + (x_7 - y_7)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - y_7 - z_7\right)\mathbf{a}_3 = -y_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - z_7\right)a\,\hat{\mathbf{y}} + x_7 a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{65} = \left(\tfrac{1}{2} + x_7 + z_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_7 + z_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_7 + y_7\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + y_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_7\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{66} = \left(\tfrac{1}{2} - x_7 + z_7\right)\mathbf{a}_1 + (-y_7 + z_7)\,\mathbf{a}_2 + (-x_7 - y_7)\,\mathbf{a}_3 = -\left(y_7 + \tfrac{1}{4}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_7\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{67} = (-x_7 - z_7)\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_7 - z_7\right)\mathbf{a}_2 + (-x_7 + y_7)\,\mathbf{a}_3 = \left(\tfrac{1}{4} + y_7\right)a\,\hat{\mathbf{x}} - \left(x_7 + \tfrac{1}{4}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - z_7\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{68} = (x_7 - z_7)\,\mathbf{a}_1 + (-y_7 - z_7)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_7 - y_7\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - y_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_7\right)a\,\hat{\mathbf{y}} - \left(z_7 + \tfrac{1}{4}\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{69} = \left(\tfrac{1}{2} + y_7 + z_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_7 + y_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_7 + z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + x_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + z_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + y_7\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{70} = (-y_7 + z_7)\,\mathbf{a}_1 + (-x_7 - y_7)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_7 + z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - x_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + z_7\right)a\,\hat{\mathbf{y}} - \left(y_7 + \tfrac{1}{4}\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{71} = \left(\tfrac{1}{2} + y_7 - z_7\right)\mathbf{a}_1 + (-x_7 + y_7)\,\mathbf{a}_2 + (-x_7 - z_7)\,\mathbf{a}_3 = -\left(x_7 + \tfrac{1}{4}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - z_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + y_7\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{72} = (-y_7 - z_7)\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_7 - y_7\right)\mathbf{a}_2 + (x_7 - z_7)\,\mathbf{a}_3 = \left(\tfrac{1}{4} + x_7\right)a\,\hat{\mathbf{x}} - \left(z_7 + \tfrac{1}{4}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - y_7\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{73} = \left(\tfrac{1}{2} + x_7 + y_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_7 + z_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + y_7 + z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + z_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + y_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + x_7\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{74} = (-x_7 - y_7)\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_7 + z_7\right)\mathbf{a}_2 + (-y_7 + z_7)\,\mathbf{a}_3 = \left(\tfrac{1}{4} + z_7\right)a\,\hat{\mathbf{x}} - \left(y_7 + \tfrac{1}{4}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - x_7\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{75} = (-x_7 + y_7)\,\mathbf{a}_1 + (-x_7 - z_7)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + y_7 - z_7\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - z_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + y_7\right)a\,\hat{\mathbf{y}} - \left(x_7 + \tfrac{1}{4}\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{76} = \left(\tfrac{1}{2} + x_7 - y_7\right)\mathbf{a}_1 + (x_7 - z_7)\,\mathbf{a}_2 + (-y_7 - z_7)\,\mathbf{a}_3 = -\left(z_7 + \tfrac{1}{4}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - y_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + x_7\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

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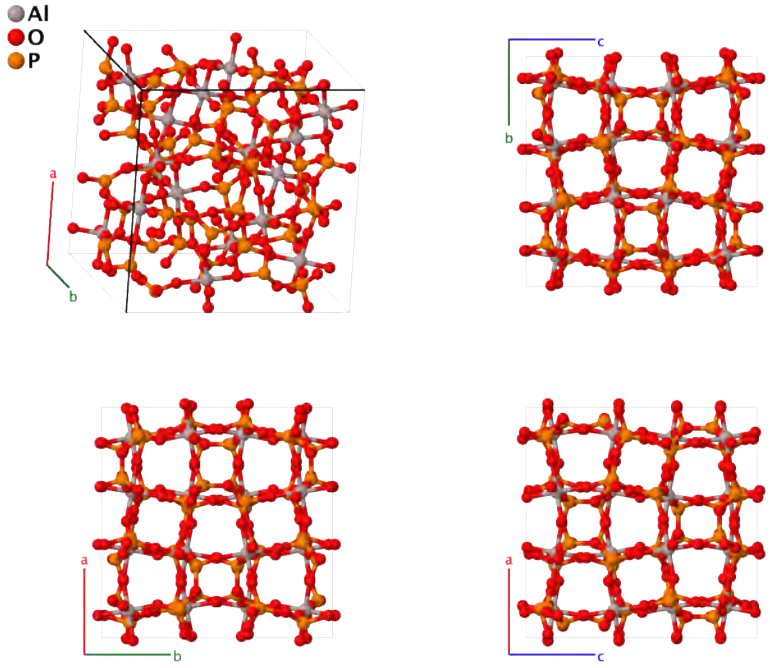
Geometry files:

- CIF: pp. 1801
- POSCAR: pp. 1802

$Al(PO_3)_3$ ($G5_2$) Structure: AB9C3_cI208_220_c_3e_e

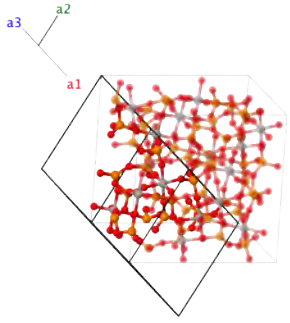
http://aflow.org/prototype-encyclopedia/AB9C3_cI208_220_c_3e_e



Prototype	:	AlO_9P_3
AFLOW prototype label	:	AB9C3_cI208_220_c_3e_e
***Strukturbericht* designation**	:	$G5_2$
Pearson symbol	:	cI208
Space group number	:	220
Space group symbol	:	$I\bar{4}3d$
AFLOW prototype command	:	`aflow --proto=AB9C3_cI208_220_c_3e_e --params=`$a, x_1, x_2, y_2, z_2, x_3, y_3, z_3, x_4, y_4, z_4, x_5, y_5, z_5$

Body-centered Cubic primitive vectors:

$$\mathbf{a}_1 = -\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}a\,\hat{\mathbf{x}} - \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} - \tfrac{1}{2}a\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$2x_1\,\mathbf{a}_1 + 2x_1\,\mathbf{a}_2 + 2x_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + x_1 a\,\hat{\mathbf{y}} + x_1 a\,\hat{\mathbf{z}}$	($16c$)	Al
$\mathbf{B}_2$	=	$\tfrac{1}{2}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - 2x_1\right)\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_1\right)a\,\hat{\mathbf{y}} + x_1 a\,\hat{\mathbf{z}}$	($16c$)	Al

$\mathbf{B}_3$	=	$\left(\frac{1}{2}-2x_1\right)\mathbf{a}_2+\frac{1}{2}\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_1\right)a\,\hat{\mathbf{x}}+x_1a\,\hat{\mathbf{y}}-x_1a\,\hat{\mathbf{z}}$	(16*c*)	Al
$\mathbf{B}_4$	=	$\left(\frac{1}{2}-2x_1\right)\mathbf{a}_1+\frac{1}{2}\mathbf{a}_2$	=	$x_1a\,\hat{\mathbf{x}}-x_1a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-x_1\right)a\,\hat{\mathbf{z}}$	(16*c*)	Al
$\mathbf{B}_5$	=	$\left(\frac{1}{2}+2x_1\right)\mathbf{a}_1+\left(\frac{1}{2}+2x_1\right)\mathbf{a}_2+\left(\frac{1}{2}+2x_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}+x_1\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+x_1\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+x_1\right)a\,\hat{\mathbf{z}}$	(16*c*)	Al
$\mathbf{B}_6$	=	$\frac{1}{2}\mathbf{a}_1-2x_1\,\mathbf{a}_3$	=	$-\left(x_1+\frac{1}{4}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-x_1\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+x_1\right)a\,\hat{\mathbf{z}}$	(16*c*)	Al
$\mathbf{B}_7$	=	$-2x_1\,\mathbf{a}_1+\frac{1}{2}\mathbf{a}_2$	=	$\left(\frac{1}{4}+x_1\right)a\,\hat{\mathbf{x}}-\left(x_1+\frac{1}{4}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-x_1\right)a\,\hat{\mathbf{z}}$	(16*c*)	Al
$\mathbf{B}_8$	=	$-2x_1\,\mathbf{a}_2+\frac{1}{2}\mathbf{a}_3$	=	$\left(\frac{1}{4}-x_1\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+x_1\right)a\,\hat{\mathbf{y}}-\left(x_1+\frac{1}{4}\right)a\,\hat{\mathbf{z}}$	(16*c*)	Al
$\mathbf{B}_9$	=	$(y_2+z_2)\,\mathbf{a}_1+(x_2+z_2)\,\mathbf{a}_2+(x_2+y_2)\,\mathbf{a}_3$	=	$x_2a\,\hat{\mathbf{x}}+y_2a\,\hat{\mathbf{y}}+z_2a\,\hat{\mathbf{z}}$	(48*e*)	O I
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2}-y_2+z_2\right)\mathbf{a}_1+(-x_2+z_2)\,\mathbf{a}_2+\left(\frac{1}{2}-x_2-y_2\right)\mathbf{a}_3$	=	$-x_2a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-y_2\right)a\,\hat{\mathbf{y}}+z_2a\,\hat{\mathbf{z}}$	(48*e*)	O I
$\mathbf{B}_{11}$	=	$(y_2-z_2)\,\mathbf{a}_1+\left(\frac{1}{2}-x_2-z_2\right)\mathbf{a}_2+\left(\frac{1}{2}-x_2+y_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-x_2\right)a\,\hat{\mathbf{x}}+y_2a\,\hat{\mathbf{y}}-z_2a\,\hat{\mathbf{z}}$	(48*e*)	O I
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2}-y_2-z_2\right)\mathbf{a}_1+\left(\frac{1}{2}+x_2-z_2\right)\mathbf{a}_2+(x_2-y_2)\,\mathbf{a}_3$	=	$x_2a\,\hat{\mathbf{x}}-y_2a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-z_2\right)a\,\hat{\mathbf{z}}$	(48*e*)	O I
$\mathbf{B}_{13}$	=	$(x_2+y_2)\,\mathbf{a}_1+(y_2+z_2)\,\mathbf{a}_2+(x_2+z_2)\,\mathbf{a}_3$	=	$z_2a\,\hat{\mathbf{x}}+x_2a\,\hat{\mathbf{y}}+y_2a\,\hat{\mathbf{z}}$	(48*e*)	O I
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2}-x_2-y_2\right)\mathbf{a}_1+\left(\frac{1}{2}-y_2+z_2\right)\mathbf{a}_2+(-x_2+z_2)\,\mathbf{a}_3$	=	$z_2a\,\hat{\mathbf{x}}-x_2a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-y_2\right)a\,\hat{\mathbf{z}}$	(48*e*)	O I
$\mathbf{B}_{15}$	=	$\left(\frac{1}{2}-x_2+y_2\right)\mathbf{a}_1+(y_2-z_2)\,\mathbf{a}_2+\left(\frac{1}{2}-x_2-z_2\right)\mathbf{a}_3$	=	$-z_2a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-x_2\right)a\,\hat{\mathbf{y}}+y_2a\,\hat{\mathbf{z}}$	(48*e*)	O I
$\mathbf{B}_{16}$	=	$(x_2-y_2)\,\mathbf{a}_1+\left(\frac{1}{2}-y_2-z_2\right)\mathbf{a}_2+\left(\frac{1}{2}+x_2-z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-z_2\right)a\,\hat{\mathbf{x}}+x_2a\,\hat{\mathbf{y}}-y_2a\,\hat{\mathbf{z}}$	(48*e*)	O I
$\mathbf{B}_{17}$	=	$(x_2+z_2)\,\mathbf{a}_1+(x_2+y_2)\,\mathbf{a}_2+(y_2+z_2)\,\mathbf{a}_3$	=	$y_2a\,\hat{\mathbf{x}}+z_2a\,\hat{\mathbf{y}}+x_2a\,\hat{\mathbf{z}}$	(48*e*)	O I
$\mathbf{B}_{18}$	=	$(-x_2+z_2)\,\mathbf{a}_1+\left(\frac{1}{2}-x_2-y_2\right)\mathbf{a}_2+\left(\frac{1}{2}-y_2+z_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}-y_2\right)a\,\hat{\mathbf{x}}+z_2a\,\hat{\mathbf{y}}-x_2a\,\hat{\mathbf{z}}$	(48*e*)	O I
$\mathbf{B}_{19}$	=	$\left(\frac{1}{2}-x_2-z_2\right)\mathbf{a}_1+\left(\frac{1}{2}-x_2+y_2\right)\mathbf{a}_2+(y_2-z_2)\,\mathbf{a}_3$	=	$y_2a\,\hat{\mathbf{x}}-z_2a\,\hat{\mathbf{y}}+\left(\frac{1}{2}-x_2\right)a\,\hat{\mathbf{z}}$	(48*e*)	O I
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2}+x_2-z_2\right)\mathbf{a}_1+(x_2-y_2)\,\mathbf{a}_2+\left(\frac{1}{2}-y_2-z_2\right)\mathbf{a}_3$	=	$-y_2a\,\hat{\mathbf{x}}+\left(\frac{1}{2}-z_2\right)a\,\hat{\mathbf{y}}+x_2a\,\hat{\mathbf{z}}$	(48*e*)	O I
$\mathbf{B}_{21}$	=	$\left(\frac{1}{2}+x_2+z_2\right)\mathbf{a}_1+\left(\frac{1}{2}+y_2+z_2\right)\mathbf{a}_2+\left(\frac{1}{2}+x_2+y_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}+y_2\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+x_2\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+z_2\right)a\,\hat{\mathbf{z}}$	(48*e*)	O I
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2}-x_2+z_2\right)\mathbf{a}_1+(-y_2+z_2)\,\mathbf{a}_2+(-x_2-y_2)\,\mathbf{a}_3$	=	$-\left(y_2+\frac{1}{4}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-x_2\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+z_2\right)a\,\hat{\mathbf{z}}$	(48*e*)	O I
$\mathbf{B}_{23}$	=	$(-x_2-z_2)\,\mathbf{a}_1+\left(\frac{1}{2}+y_2-z_2\right)\mathbf{a}_2+(-x_2+y_2)\,\mathbf{a}_3$	=	$\left(\frac{1}{4}+y_2\right)a\,\hat{\mathbf{x}}-\left(x_2+\frac{1}{4}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-z_2\right)a\,\hat{\mathbf{z}}$	(48*e*)	O I
$\mathbf{B}_{24}$	=	$(x_2-z_2)\,\mathbf{a}_1+(-y_2-z_2)\,\mathbf{a}_2+\left(\frac{1}{2}+x_2-y_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}-y_2\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+x_2\right)a\,\hat{\mathbf{y}}-\left(z_2+\frac{1}{4}\right)a\,\hat{\mathbf{z}}$	(48*e*)	O I

$$\mathbf{B}_{25} = \left(\tfrac{1}{2}+y_2+z_2\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+x_2+y_2\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+x_2+z_2\right)\mathbf{a}_3 = \left(\tfrac{1}{4}+x_2\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4}+z_2\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4}+y_2\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O I}$$

$$\mathbf{B}_{26} = (-y_2+z_2)\,\mathbf{a}_1 + (-x_2-y_2)\,\mathbf{a}_2 + \left(\tfrac{1}{2}-x_2+z_2\right)\mathbf{a}_3 = \left(\tfrac{1}{4}-x_2\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4}+z_2\right)a\,\hat{\mathbf{y}} - \left(y_2+\tfrac{1}{4}\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O I}$$

$$\mathbf{B}_{27} = \left(\tfrac{1}{2}+y_2-z_2\right)\mathbf{a}_1 + (-x_2+y_2)\,\mathbf{a}_2 + (-x_2-z_2)\,\mathbf{a}_3 = -\left(x_2+\tfrac{1}{4}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4}-z_2\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4}+y_2\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O I}$$

$$\mathbf{B}_{28} = (-y_2-z_2)\,\mathbf{a}_1 + \left(\tfrac{1}{2}+x_2-y_2\right)\mathbf{a}_2 + (x_2-z_2)\,\mathbf{a}_3 = \left(\tfrac{1}{4}+x_2\right)a\,\hat{\mathbf{x}} - \left(z_2+\tfrac{1}{4}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4}-y_2\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O I}$$

$$\mathbf{B}_{29} = \left(\tfrac{1}{2}+x_2+y_2\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+x_2+z_2\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+y_2+z_2\right)\mathbf{a}_3 = \left(\tfrac{1}{4}+z_2\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4}+y_2\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4}+x_2\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O I}$$

$$\mathbf{B}_{30} = (-x_2-y_2)\,\mathbf{a}_1 + \left(\tfrac{1}{2}-x_2+z_2\right)\mathbf{a}_2 + (-y_2+z_2)\,\mathbf{a}_3 = \left(\tfrac{1}{4}+z_2\right)a\,\hat{\mathbf{x}} - \left(y_2+\tfrac{1}{4}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4}-x_2\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O I}$$

$$\mathbf{B}_{31} = (-x_2+y_2)\,\mathbf{a}_1 + (-x_2-z_2)\,\mathbf{a}_2 + \left(\tfrac{1}{2}+y_2-z_2\right)\mathbf{a}_3 = \left(\tfrac{1}{4}-z_2\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4}+y_2\right)a\,\hat{\mathbf{y}} - \left(x_2+\tfrac{1}{4}\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O I}$$

$$\mathbf{B}_{32} = \left(\tfrac{1}{2}+x_2-y_2\right)\mathbf{a}_1 + (x_2-z_2)\,\mathbf{a}_2 + (-y_2-z_2)\,\mathbf{a}_3 = -\left(z_2+\tfrac{1}{4}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4}-y_2\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4}+x_2\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O I}$$

$$\mathbf{B}_{33} = (y_3+z_3)\,\mathbf{a}_1 + (x_3+z_3)\,\mathbf{a}_2 + (x_3+y_3)\,\mathbf{a}_3 = x_3a\,\hat{\mathbf{x}} + y_3a\,\hat{\mathbf{y}} + z_3a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O II}$$

$$\mathbf{B}_{34} = \left(\tfrac{1}{2}-y_3+z_3\right)\mathbf{a}_1 + (-x_3+z_3)\,\mathbf{a}_2 + \left(\tfrac{1}{2}-x_3-y_3\right)\mathbf{a}_3 = -x_3a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_3\right)a\,\hat{\mathbf{y}} + z_3a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O II}$$

$$\mathbf{B}_{35} = (y_3-z_3)\,\mathbf{a}_1 + \left(\tfrac{1}{2}-x_3-z_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-x_3+y_3\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_3\right)a\,\hat{\mathbf{x}} + y_3a\,\hat{\mathbf{y}} - z_3a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O II}$$

$$\mathbf{B}_{36} = \left(\tfrac{1}{2}-y_3-z_3\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+x_3-z_3\right)\mathbf{a}_2 + (x_3-y_3)\,\mathbf{a}_3 = x_3a\,\hat{\mathbf{x}} - y_3a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_3\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O II}$$

$$\mathbf{B}_{37} = (x_3+y_3)\,\mathbf{a}_1 + (y_3+z_3)\,\mathbf{a}_2 + (x_3+z_3)\,\mathbf{a}_3 = z_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + y_3a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O II}$$

$$\mathbf{B}_{38} = \left(\tfrac{1}{2}-x_3-y_3\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_3+z_3\right)\mathbf{a}_2 + (-x_3+z_3)\,\mathbf{a}_3 = z_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-y_3\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O II}$$

$$\mathbf{B}_{39} = \left(\tfrac{1}{2}-x_3+y_3\right)\mathbf{a}_1 + (y_3-z_3)\,\mathbf{a}_2 + \left(\tfrac{1}{2}-x_3-z_3\right)\mathbf{a}_3 = -z_3a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_3\right)a\,\hat{\mathbf{y}} + y_3a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O II}$$

$$\mathbf{B}_{40} = (x_3-y_3)\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_3-z_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+x_3-z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-z_3\right)a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} - y_3a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O II}$$

$$\mathbf{B}_{41} = (x_3+z_3)\,\mathbf{a}_1 + (x_3+y_3)\,\mathbf{a}_2 + (y_3+z_3)\,\mathbf{a}_3 = y_3a\,\hat{\mathbf{x}} + z_3a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O II}$$

$$\mathbf{B}_{42} = (-x_3+z_3)\,\mathbf{a}_1 + \left(\tfrac{1}{2}-x_3-y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-y_3+z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-y_3\right)a\,\hat{\mathbf{x}} + z_3a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O II}$$

$$\mathbf{B}_{43} = \left(\tfrac{1}{2}-x_3-z_3\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-x_3+y_3\right)\mathbf{a}_2 + (y_3-z_3)\,\mathbf{a}_3 = y_3a\,\hat{\mathbf{x}} - z_3a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-x_3\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O II}$$

$$\mathbf{B}_{44} = \left(\tfrac{1}{2}+x_3-z_3\right)\mathbf{a}_1 + (x_3-y_3)\,\mathbf{a}_2 + \left(\tfrac{1}{2}-y_3-z_3\right)\mathbf{a}_3 = -y_3a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-z_3\right)a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O II}$$

$$\mathbf{B}_{45} = \left(\tfrac{1}{2}+x_3+z_3\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+y_3+z_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+x_3+y_3\right)\mathbf{a}_3 = \left(\tfrac{1}{4}+y_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4}+x_3\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4}+z_3\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O II}$$

$$\mathbf{B}_{46} = \left(\tfrac{1}{2} - x_3 + z_3\right)\mathbf{a}_1 + (-y_3 + z_3)\,\mathbf{a}_2 + (-x_3 - y_3)\,\mathbf{a}_3 = -\left(y_3 + \tfrac{1}{4}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_3\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_3\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O II}$$

$$\mathbf{B}_{47} = (-x_3 - z_3)\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_3 - z_3\right)\mathbf{a}_2 + (-x_3 + y_3)\,\mathbf{a}_3 = \left(\tfrac{1}{4} + y_3\right)a\,\hat{\mathbf{x}} - \left(x_3 + \tfrac{1}{4}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - z_3\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O II}$$

$$\mathbf{B}_{48} = (x_3 - z_3)\,\mathbf{a}_1 + (-y_3 - z_3)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_3 - y_3\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - y_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_3\right)a\,\hat{\mathbf{y}} - \left(z_3 + \tfrac{1}{4}\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O II}$$

$$\mathbf{B}_{49} = \left(\tfrac{1}{2} + y_3 + z_3\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_3 + y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_3 + z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + x_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + z_3\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + y_3\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O II}$$

$$\mathbf{B}_{50} = (-y_3 + z_3)\,\mathbf{a}_1 + (-x_3 - y_3)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_3 + z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - x_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + z_3\right)a\,\hat{\mathbf{y}} - \left(y_3 + \tfrac{1}{4}\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O II}$$

$$\mathbf{B}_{51} = \left(\tfrac{1}{2} + y_3 - z_3\right)\mathbf{a}_1 + (-x_3 + y_3)\,\mathbf{a}_2 + (-x_3 - z_3)\,\mathbf{a}_3 = -\left(x_3 + \tfrac{1}{4}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - z_3\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + y_3\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O II}$$

$$\mathbf{B}_{52} = (-y_3 - z_3)\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_3 - y_3\right)\mathbf{a}_2 + (x_3 - z_3)\,\mathbf{a}_3 = \left(\tfrac{1}{4} + x_3\right)a\,\hat{\mathbf{x}} - \left(z_3 + \tfrac{1}{4}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - y_3\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O II}$$

$$\mathbf{B}_{53} = \left(\tfrac{1}{2} + x_3 + y_3\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_3 + z_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + y_3 + z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + z_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + y_3\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + x_3\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O II}$$

$$\mathbf{B}_{54} = (-x_3 - y_3)\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_3 + z_3\right)\mathbf{a}_2 + (-y_3 + z_3)\,\mathbf{a}_3 = \left(\tfrac{1}{4} + z_3\right)a\,\hat{\mathbf{x}} - \left(y_3 + \tfrac{1}{4}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - x_3\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O II}$$

$$\mathbf{B}_{55} = (-x_3 + y_3)\,\mathbf{a}_1 + (-x_3 - z_3)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + y_3 - z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - z_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + y_3\right)a\,\hat{\mathbf{y}} - \left(x_3 + \tfrac{1}{4}\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O II}$$

$$\mathbf{B}_{56} = \left(\tfrac{1}{2} + x_3 - y_3\right)\mathbf{a}_1 + (x_3 - z_3)\,\mathbf{a}_2 + (-y_3 - z_3)\,\mathbf{a}_3 = -\left(z_3 + \tfrac{1}{4}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - y_3\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + x_3\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O II}$$

$$\mathbf{B}_{57} = (y_4 + z_4)\,\mathbf{a}_1 + (x_4 + z_4)\,\mathbf{a}_2 + (x_4 + y_4)\,\mathbf{a}_3 = x_4 a\,\hat{\mathbf{x}} + y_4 a\,\hat{\mathbf{y}} + z_4 a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{58} = \left(\tfrac{1}{2} - y_4 + z_4\right)\mathbf{a}_1 + (-x_4 + z_4)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_4 - y_4\right)\mathbf{a}_3 = -x_4 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_4\right)a\,\hat{\mathbf{y}} + z_4 a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{59} = (y_4 - z_4)\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_4 - z_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_4 + y_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + y_4 a\,\hat{\mathbf{y}} - z_4 a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{60} = \left(\tfrac{1}{2} - y_4 - z_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_4 - z_4\right)\mathbf{a}_2 + (x_4 - y_4)\,\mathbf{a}_3 = x_4 a\,\hat{\mathbf{x}} - y_4 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_4\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{61} = (x_4 + y_4)\,\mathbf{a}_1 + (y_4 + z_4)\,\mathbf{a}_2 + (x_4 + z_4)\,\mathbf{a}_3 = z_4 a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}} + y_4 a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{62} = \left(\tfrac{1}{2} - x_4 - y_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_4 + z_4\right)\mathbf{a}_2 + (-x_4 + z_4)\,\mathbf{a}_3 = z_4 a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - y_4\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{63} = \left(\tfrac{1}{2} - x_4 + y_4\right)\mathbf{a}_1 + (y_4 - z_4)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_4 - z_4\right)\mathbf{a}_3 = -z_4 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_4\right)a\,\hat{\mathbf{y}} + y_4 a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{64} = (x_4 - y_4)\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_4 - z_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_4 - z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - z_4\right)a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}} - y_4 a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{65} = (x_4 + z_4)\,\mathbf{a}_1 + (x_4 + y_4)\,\mathbf{a}_2 + (y_4 + z_4)\,\mathbf{a}_3 = y_4 a\,\hat{\mathbf{x}} + z_4 a\,\hat{\mathbf{y}} + x_4 a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{66} = (-x_4 + z_4)\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_4 - y_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - y_4 + z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - y_4\right)a\,\hat{\mathbf{x}} + z_4 a\,\hat{\mathbf{y}} - x_4 a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{67} = \left(\tfrac{1}{2} - x_4 - z_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_4 + y_4\right)\mathbf{a}_2 + (y_4 - z_4)\,\mathbf{a}_3 = y_4 a\,\hat{\mathbf{x}} - z_4 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_4\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{68} = \left(\tfrac{1}{2} + x_4 - z_4\right)\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - y_4 - z_4\right)\mathbf{a}_3 = -y_4 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - z_4\right)a\,\hat{\mathbf{y}} + x_4 a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{69} = \left(\tfrac{1}{2} + x_4 + z_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_4 + z_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_4 + y_4\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + y_4\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_4\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{70} = \left(\tfrac{1}{2} - x_4 + z_4\right)\mathbf{a}_1 + (-y_4 + z_4)\,\mathbf{a}_2 + (-x_4 - y_4)\,\mathbf{a}_3 = -\left(y_4 + \tfrac{1}{4}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_4\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{71} = (-x_4 - z_4)\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_4 - z_4\right)\mathbf{a}_2 + (-x_4 + y_4)\,\mathbf{a}_3 = \left(\tfrac{1}{4} + y_4\right)a\,\hat{\mathbf{x}} - \left(x_4 + \tfrac{1}{4}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - z_4\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{72} = (x_4 - z_4)\,\mathbf{a}_1 + (-y_4 - z_4)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_4 - y_4\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - y_4\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_4\right)a\,\hat{\mathbf{y}} - \left(z_4 + \tfrac{1}{4}\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{73} = \left(\tfrac{1}{2} + y_4 + z_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_4 + y_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_4 + z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + x_4\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + z_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + y_4\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{74} = (-y_4 + z_4)\,\mathbf{a}_1 + (-x_4 - y_4)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_4 + z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - x_4\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + z_4\right)a\,\hat{\mathbf{y}} - \left(y_4 + \tfrac{1}{4}\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{75} = \left(\tfrac{1}{2} + y_4 - z_4\right)\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 + (-x_4 - z_4)\,\mathbf{a}_3 = -\left(x_4 + \tfrac{1}{4}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - z_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + y_4\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{76} = (-y_4 - z_4)\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_4 - y_4\right)\mathbf{a}_2 + (x_4 - z_4)\,\mathbf{a}_3 = \left(\tfrac{1}{4} + x_4\right)a\,\hat{\mathbf{x}} - \left(z_4 + \tfrac{1}{4}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - y_4\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{77} = \left(\tfrac{1}{2} + x_4 + y_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_4 + z_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + y_4 + z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + z_4\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + y_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + x_4\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{78} = (-x_4 - y_4)\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_4 + z_4\right)\mathbf{a}_2 + (-y_4 + z_4)\,\mathbf{a}_3 = \left(\tfrac{1}{4} + z_4\right)a\,\hat{\mathbf{x}} - \left(y_4 + \tfrac{1}{4}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - x_4\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{79} = (-x_4 + y_4)\,\mathbf{a}_1 + (-x_4 - z_4)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + y_4 - z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - z_4\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + y_4\right)a\,\hat{\mathbf{y}} - \left(x_4 + \tfrac{1}{4}\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{80} = \left(\tfrac{1}{2} + x_4 - y_4\right)\mathbf{a}_1 + (x_4 - z_4)\,\mathbf{a}_2 + (-y_4 - z_4)\,\mathbf{a}_3 = -\left(z_4 + \tfrac{1}{4}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - y_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + x_4\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{O III}$$

$$\mathbf{B}_{81} = (y_5 + z_5)\,\mathbf{a}_1 + (x_5 + z_5)\,\mathbf{a}_2 + (x_5 + y_5)\,\mathbf{a}_3 = x_5 a\,\hat{\mathbf{x}} + y_5 a\,\hat{\mathbf{y}} + z_5 a\,\hat{\mathbf{z}} \quad (48e) \quad \text{P}$$

$$\mathbf{B}_{82} = \left(\tfrac{1}{2} - y_5 + z_5\right)\mathbf{a}_1 + (-x_5 + z_5)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_5 - y_5\right)\mathbf{a}_3 = -x_5 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_5\right)a\,\hat{\mathbf{y}} + z_5 a\,\hat{\mathbf{z}} \quad (48e) \quad \text{P}$$

$$\mathbf{B}_{83} = (y_5 - z_5)\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_5 - z_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_5 + y_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + y_5 a\,\hat{\mathbf{y}} - z_5 a\,\hat{\mathbf{z}} \quad (48e) \quad \text{P}$$

$$\mathbf{B}_{84} = \left(\tfrac{1}{2} - y_5 - z_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_5 - z_5\right)\mathbf{a}_2 + (x_5 - y_5)\,\mathbf{a}_3 = x_5 a\,\hat{\mathbf{x}} - y_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_5\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{P}$$

$$\mathbf{B}_{85} = (x_5 + y_5)\,\mathbf{a}_1 + (y_5 + z_5)\,\mathbf{a}_2 + (x_5 + z_5)\,\mathbf{a}_3 = z_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + y_5 a\,\hat{\mathbf{z}} \quad (48e) \quad \text{P}$$

$$\mathbf{B}_{86} = \left(\tfrac{1}{2} - x_5 - y_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_5 + z_5\right)\mathbf{a}_2 + (-x_5 + z_5)\,\mathbf{a}_3 = z_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - y_5\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{P}$$

$$\mathbf{B}_{87} = \left(\tfrac{1}{2} - x_5 + y_5\right)\mathbf{a}_1 + (y_5 - z_5)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_5 - z_5\right)\mathbf{a}_3 = -z_5 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_5\right)a\,\hat{\mathbf{y}} + y_5 a\,\hat{\mathbf{z}} \quad (48e) \quad \text{P}$$

$$\mathbf{B}_{88} = (x_5 - y_5)\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_5 - z_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_5 - z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - z_5\right)a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} - y_5 a\,\hat{\mathbf{z}} \quad (48e) \quad \text{P}$$

$$\mathbf{B}_{89} = (x_5 + z_5)\,\mathbf{a}_1 + (x_5 + y_5)\,\mathbf{a}_2 + (y_5 + z_5)\,\mathbf{a}_3 = y_5 a\,\hat{\mathbf{x}} + z_5 a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}} \quad (48e) \quad \text{P}$$

$$\mathbf{B}_{90} = (-x_5 + z_5)\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_5 - y_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - y_5 + z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - y_5\right)a\,\hat{\mathbf{x}} + z_5 a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}} \quad (48e) \quad \text{P}$$

$$\mathbf{B}_{91} = \left(\tfrac{1}{2} - x_5 - z_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_5 + y_5\right)\mathbf{a}_2 + (y_5 - z_5)\,\mathbf{a}_3 = y_5 a\,\hat{\mathbf{x}} - z_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_5\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{P}$$

$$\mathbf{B}_{92} = \left(\tfrac{1}{2} + x_5 - z_5\right)\mathbf{a}_1 + (x_5 - y_5)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - y_5 - z_5\right)\mathbf{a}_3 = -y_5 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - z_5\right)a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}} \quad (48e) \quad \text{P}$$

$$\mathbf{B}_{93} = \left(\tfrac{1}{2} + x_5 + z_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_5 + z_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_5 + y_5\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + y_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_5\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{P}$$

$$\mathbf{B}_{94} = \left(\tfrac{1}{2} - x_5 + z_5\right)\mathbf{a}_1 + (-y_5 + z_5)\,\mathbf{a}_2 + (-x_5 - y_5)\,\mathbf{a}_3 = -\left(y_5 + \tfrac{1}{4}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_5\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{P}$$

$$\mathbf{B}_{95} = (-x_5 - z_5)\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_5 - z_5\right)\mathbf{a}_2 + (-x_5 + y_5)\,\mathbf{a}_3 = \left(\tfrac{1}{4} + y_5\right)a\,\hat{\mathbf{x}} - \left(x_5 + \tfrac{1}{4}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - z_5\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{P}$$

$$\mathbf{B}_{96} = (x_5 - z_5)\,\mathbf{a}_1 + (-y_5 - z_5)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_5 - y_5\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - y_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_5\right)a\,\hat{\mathbf{y}} - \left(z_5 + \tfrac{1}{4}\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{P}$$

$$\mathbf{B}_{97} = \left(\tfrac{1}{2} + y_5 + z_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_5 + y_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_5 + z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + x_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + z_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + y_5\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{P}$$

$$\mathbf{B}_{98} = (-y_5 + z_5)\,\mathbf{a}_1 + (-x_5 - y_5)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_5 + z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - x_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + z_5\right)a\,\hat{\mathbf{y}} - \left(y_5 + \tfrac{1}{4}\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{P}$$

$$\mathbf{B}_{99} = \left(\tfrac{1}{2} + y_5 - z_5\right)\mathbf{a}_1 + (-x_5 + y_5)\,\mathbf{a}_2 + (-x_5 - z_5)\,\mathbf{a}_3 = -\left(x_5 + \tfrac{1}{4}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - z_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + y_5\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{P}$$

$$\mathbf{B}_{100} = (-y_5 - z_5)\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_5 - y_5\right)\mathbf{a}_2 + (x_5 - z_5)\,\mathbf{a}_3 = \left(\tfrac{1}{4} + x_5\right)a\,\hat{\mathbf{x}} - \left(z_5 + \tfrac{1}{4}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - y_5\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{P}$$

$$\mathbf{B}_{101} = \left(\tfrac{1}{2} + x_5 + y_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_5 + z_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + y_5 + z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + z_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + y_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + x_5\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{P}$$

$$\mathbf{B}_{102} = (-x_5 - y_5)\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_5 + z_5\right)\mathbf{a}_2 + (-y_5 + z_5)\,\mathbf{a}_3 = \left(\tfrac{1}{4} + z_5\right)a\,\hat{\mathbf{x}} - \left(y_5 + \tfrac{1}{4}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - x_5\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{P}$$

$$\mathbf{B}_{103} = (-x_5 + y_5)\,\mathbf{a}_1 + (-x_5 - z_5)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + y_5 - z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - z_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + y_5\right)a\,\hat{\mathbf{y}} - \left(x_5 + \tfrac{1}{4}\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{P}$$

$$\mathbf{B}_{104} = \left(\tfrac{1}{2} + x_5 - y_5\right)\mathbf{a}_1 + (x_5 - z_5)\,\mathbf{a}_2 + (-y_5 - z_5)\,\mathbf{a}_3 = -\left(z_5 + \tfrac{1}{4}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - y_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + x_5\right)a\,\hat{\mathbf{z}} \quad (48e) \quad \text{P}$$

References:

- L. Pauling and J. Sherman, *The Crystal Structure of Aluminum Metaphosphate, Al(PO_3)$_3$*, Zeitschrift für Kristallographie - Crystalline Materials **96**, 481–487 (1937), doi:10.1524/zkri.1937.96.1.481.

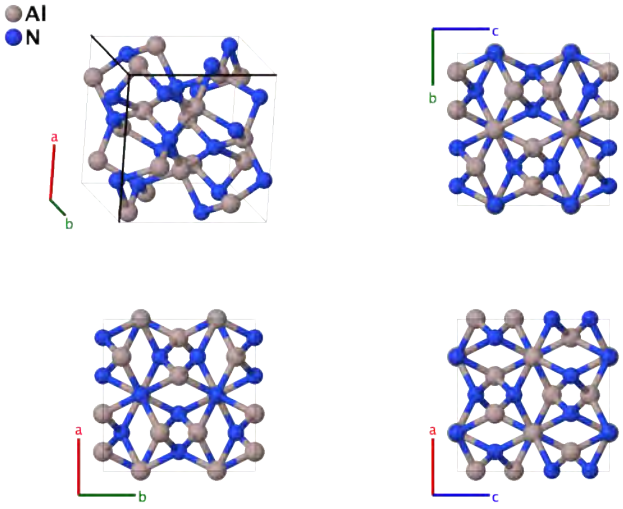
Found in:

- H. van der Meer, *The crystal structure of a monoclinic form of aluminium metaphosphate, Al(PO_3)$_3$*, Acta Crystallogr. Sect. B Struct. Sci. **32**, 2423–2426 (1976), doi:10.1107/S0567740876007899.

Geometry files:
- CIF: pp. 1802
- POSCAR: pp. 1803

AlN (cI24) Structure: AB_cI24_220_a_b

http://aflow.org/prototype-encyclopedia/AB_cI24_220_a_b

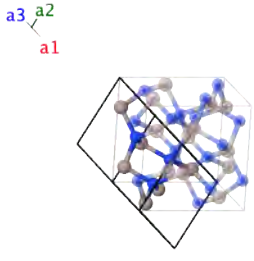


Prototype	:	AlN
AFLOW prototype label	:	AB_cI24_220_a_b
***Strukturbericht* designation**	:	None
Pearson symbol	:	cI24
Space group number	:	220
Space group symbol	:	$I\bar{4}3d$
AFLOW prototype command	:	`aflow --proto=AB_cI24_220_a_b --params=`a

- AlN naturally occurs in two forms (Liu, 2019): the stable wz-AlN wurtzite ($B4$) structure, and the high-pressure rs-AlN rock salt ($B1$) structure. A metastable zb-AlN zincblende (zb-AlN) structure can be synthesized via a solid-state reaction.
- (Liu, 2019) used a first-principles evolutionary technique to find four possible metastable phases: one in the SC16 structure, and three novel cubic structures, cF40, cI16, and cI24.

Body-centered Cubic primitive vectors:

$$\mathbf{a}_1 = -\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} - \frac{1}{2}a\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$ =	$\frac{1}{4}\mathbf{a}_1 + \frac{5}{8}\mathbf{a}_2 + \frac{3}{8}\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	($12a$)	Al

$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{7}{8}\,\mathbf{a}_2 + \frac{1}{8}\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(12*a*)	Al
$\mathbf{B}_3$	=	$\frac{3}{8}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{5}{8}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}}$	(12*a*)	Al
$\mathbf{B}_4$	=	$\frac{1}{8}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{7}{8}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}}$	(12*a*)	Al
$\mathbf{B}_5$	=	$\frac{5}{8}\,\mathbf{a}_1 + \frac{3}{8}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}}$	(12*a*)	Al
$\mathbf{B}_6$	=	$\frac{7}{8}\,\mathbf{a}_1 + \frac{1}{8}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(12*a*)	Al
$\mathbf{B}_7$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{8}\,\mathbf{a}_2 + \frac{7}{8}\,\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} - \frac{1}{4}a\,\hat{\mathbf{z}}$	(12*b*)	N
$\mathbf{B}_8$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{8}\,\mathbf{a}_2 + \frac{5}{8}\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(12*b*)	N
$\mathbf{B}_9$	=	$\frac{7}{8}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{8}\,\mathbf{a}_3$	=	$-\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(12*b*)	N
$\mathbf{B}_{10}$	=	$\frac{5}{8}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{3}{8}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(12*b*)	N
$\mathbf{B}_{11}$	=	$\frac{1}{8}\,\mathbf{a}_1 + \frac{7}{8}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}}$	(12*b*)	N
$\mathbf{B}_{12}$	=	$\frac{3}{8}\,\mathbf{a}_1 + \frac{5}{8}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(12*b*)	N

References:

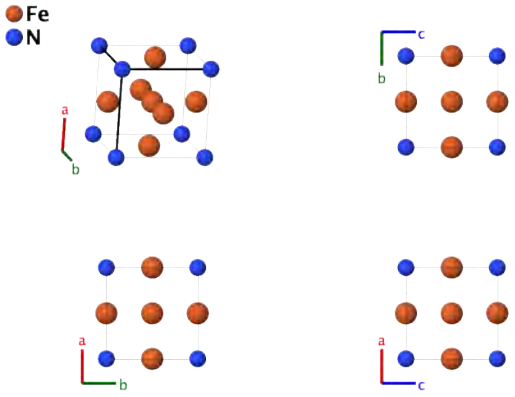
- C. Liu, M. Chen, J. Li, L. Liu, P. Li, M. Ma, C. Shao, J. He, and T. Liang, *A first-principles study of novel cubic AlN phases*, J. Phys. Chem. Solids **130**, 58–66 (2019), doi:10.1016/j.jpcs.2019.02.009.

Geometry files:

- CIF: pp. 1804
- POSCAR: pp. 1804

γ-Fe_4N ($L'1_0$) Structure: A4B_cP5_221_bc_a

http://aflow.org/prototype-encyclopedia/A4B_cP5_221_bc_a

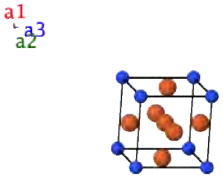


Prototype	:	Fe_4N
AFLOW prototype label	:	A4B_cP5_221_bc_a
***Strukturbericht* designation**	:	$L'1_0$
Pearson symbol	:	cP5
Space group number	:	221
Space group symbol	:	$Pm\bar{3}m$
AFLOW prototype command	:	`aflow --proto=A4B_cP5_221_bc_a --params=`a

- This is a two-component form of the cubic perovskite structure (AB3C_cP5_221_a_c_b), *Strukturbericht* designation $E2_1$.

Simple Cubic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = a\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($1a$)	N
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	($1b$)	Fe I
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	($3c$)	Fe II
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	($3c$)	Fe II
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$	($3c$)	Fe II

References:

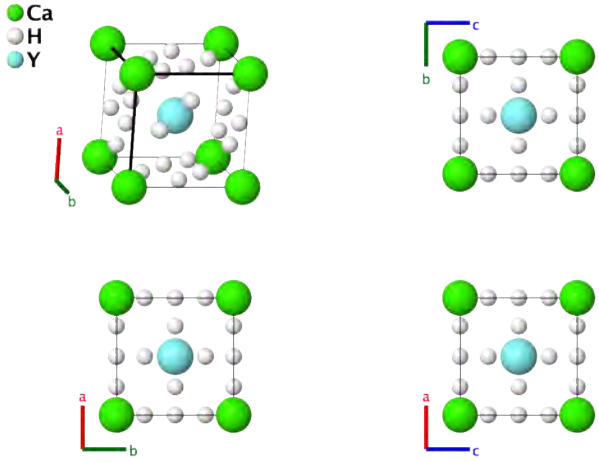
- H. Jacobs, R. Rechenbach, and U. Zachwieja, *Structure determination of γ'-Fe_4N and ϵ-Fe_3N*, J. Alloys Compd. **227**, 10–17 (1995), doi:10.1016/0925-8388(95)01610-4.

Geometry files:

- CIF: pp. 1804
- POSCAR: pp. 1805

Predicted High-Pressure $YCaH_{12}$ Structure: AB12C_cP14_221_a_h_b

http://aflow.org/prototype-encyclopedia/AB12C_cP14_221_a_h_b

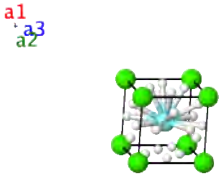


Prototype	:	$CaH_{12}Y$
AFLOW prototype label	:	AB12C_cP14_221_a_h_b
***Strukturbericht* designation**	:	None
Pearson symbol	:	cP14
Space group number	:	221
Space group symbol	:	$Pm\bar{3}m$
AFLOW prototype command	:	`aflow --proto=AB12C_cP14_221_a_h_b --params=`a, x_3

- This structure was determined by *ab initio* methods and is predicted to be stable in the pressure range 180-257 GPa, with T_c = 230 K at 180 GPa. We show the predicted structure at 200 GPa.

Simple Cubic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = a\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($1a$)	Ca
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	($1b$)	Y
$\mathbf{B}_3$	=	$x_3\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$x_3 a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$	($12h$)	H

$\mathbf{B}_4$	=	$-x_3\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$-x_3 a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$	$(12h)$	H
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	$(12h)$	H
$\mathbf{B}_6$	=	$-x_3\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	$(12h)$	H
$\mathbf{B}_7$	=	$\frac{1}{2}\,\mathbf{a}_1 + x_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{z}}$	$(12h)$	H
$\mathbf{B}_8$	=	$\frac{1}{2}\,\mathbf{a}_1 - x_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{z}}$	$(12h)$	H
$\mathbf{B}_9$	=	$\frac{1}{2}\,\mathbf{a}_1 + x_3\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + x_3 a\,\hat{\mathbf{y}}$	$(12h)$	H
$\mathbf{B}_{10}$	=	$\frac{1}{2}\,\mathbf{a}_1 - x_3\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} - x_3 a\,\hat{\mathbf{y}}$	$(12h)$	H
$\mathbf{B}_{11}$	=	$x_3\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$x_3 a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	$(12h)$	H
$\mathbf{B}_{12}$	=	$-x_3\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$-x_3 a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	$(12h)$	H
$\mathbf{B}_{13}$	=	$\frac{1}{2}\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} - x_3 a\,\hat{\mathbf{z}}$	$(12h)$	H
$\mathbf{B}_{14}$	=	$\frac{1}{2}\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + x_3 a\,\hat{\mathbf{z}}$	$(12h)$	H

References:

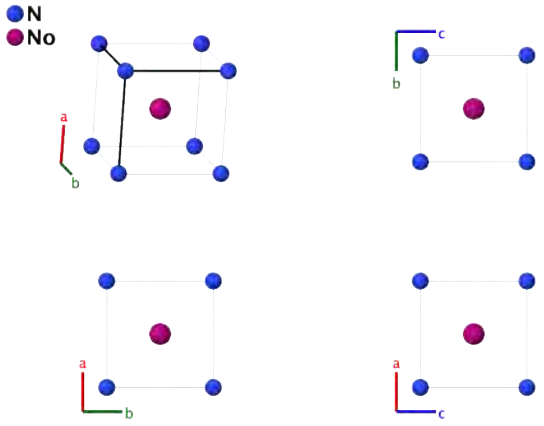
- H. Xie, D. Duan, Z. Shao, H. Song, Y. Wang, X. Xiao, D. Li, F. Tian, B. Liu, and T. Cui, *High-temperature superconductivity in ternary clathrate $YCaH_{12}$ under high pressures*, J. Phys.: Condens. Matter **31**, 245404 (2019), doi:10.1088/1361-648X/ab09b4.

Geometry files:

- CIF: pp. 1805
- POSCAR: pp. 1805

NH_4NO_3 I ($G0_8$) Structure: AB_cP2_221_a_b

http://aflow.org/prototype-encyclopedia/AB_cP2_221_a_b.NH4.NO3



Prototype	:	NH_4NO_3
AFLOW prototype label	:	AB_cP2_221_a_b
***Strukturbericht* designation**	:	$G0_8$
Pearson symbol	:	cP2
Space group number	:	221
Space group symbol	:	$Pm\bar{3}m$
AFLOW prototype command	:	`aflow --proto=AB_cP2_221_a_b --params=`a

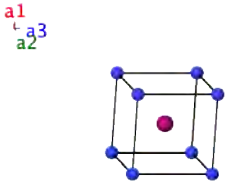
- Ammonium Nitrate exists in a variety of forms, (Hermann, 1937) depending on the temperature:

Phase	Temperature °C	*Strukturbericht*	Page	
I	125 – 170	$G0_8$	AB_cP2_221_a_b.NH4.NO3	(this structure)
II	84 – 125	$G0_9$	ABC3_tP10_100_b_a_bc	
III	32 – 84	$G0_{10}$	ABC3_oP20_62_c_c_cd.N.NH4.O	
IV	-17 – 32	$G0_{11}$	A4B2C3_oP18_59_ef_ab_af	
V	< −17	Gwihabaite	A4B2C3_tP72_77_8d_ab2c2d_6d2	

- In the $G0_8$ structure, both the NH_4^+ and NO_3^- ions rotate freely about their centers of mass (Kracek, 1931). The two ions sit on the same sites as atoms in the CsCl ($B2$) structure.

Simple Cubic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = a\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(1a)$	NH_4
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	$(1b)$	NO_3

References:

- F. C. Kracek, S. B. Hendricks, and E. Posnjak, *Group Rotation in Solid Ammonium and Calcium Nitrates*, Nature **128**, 410–411 (1931), doi:10.1038/128410b0.

Found in:

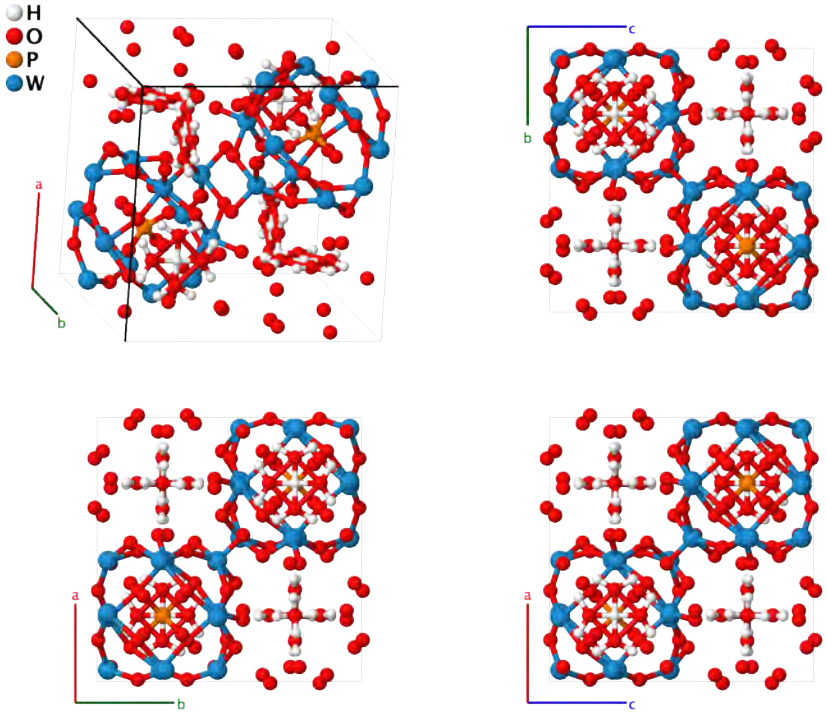
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1806
- POSCAR: pp. 1806

Dodecatungstophosphoric Acid Hexahydrate [$H_3PW_{12}O_{40}\cdot 6H_2O$] Structure: A27B52CD12_cP184_224_dl_eh3k_a_k

http://aflow.org/prototype-encyclopedia/A27B52CD12_cP184_224_dl_eh3k_a_k

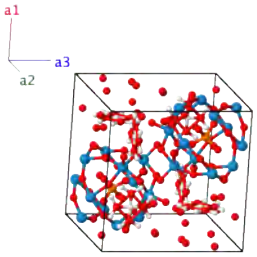


Prototype	:	$H_{15}O_{46}PW_{12}$
AFLOW prototype label	:	A27B52CD12_cP184_224_dl_eh3k_a_k
***Strukturbericht* designation**	:	None
Pearson symbol	:	cP184
Space group number	:	224
Space group symbol	:	$Pn\bar{3}m$
AFLOW prototype command	:	`aflow --proto=A27B52CD12_cP184_224_dl_eh3k_a_k --params=`$a, x_3, x_4, x_5, z_5, x_6, z_6, x_7, z_7, x_8, z_8, x_9, y_9, z_9$

- (Brown, 1977) presents this as an improvement on the $H4_{16}$ structure, $H_3PW_{12}O_{40}\cdot 5H_2O$. The primary difference is the addition of a sixth water molecule and the location of the hydrogen molecules not directly attached to a water molecule.
- The water molecules are formed by the H-II and O-II atoms, and the (24*h*) (O-II) and (48*l*) (H-II) Wyckoff sites are only occupied half of the time. Presumably this means that the nearly flat H-O molecular ions in this structure actually consist of one water molecule, the central hydrogen atom (H-I), and a water molecule on the other side of the central hydrogen, with the other water positions empty. Exactly which water molecules are occupied on around each H-I atom is completely up to chance. (Brown, 1977) state that the molecule has a positive charge, and write it as $H_5O_2^+$.
- We use the neutron data from (Brown, 1977) to locate the non-hydrogen atoms.
- This structure is a partially dehydrated form of $H_3PW_{12}O_{40}\cdot 29H_2O$ ($H4_{21}$). Further dehydration produces the $H_3PW_{12}O_{40}\cdot 3H_2O$ structure.

Simple Cubic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = a\,\hat{\mathbf{z}}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(2*a*)	P
$\mathbf{B}_2$	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(2*a*)	P
$\mathbf{B}_3$	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(6*d*)	H I
$\mathbf{B}_4$	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(6*d*)	H I
$\mathbf{B}_5$	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(6*d*)	H I
$\mathbf{B}_6$	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(6*d*)	H I
$\mathbf{B}_7$	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(6*d*)	H I
$\mathbf{B}_8$	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(6*d*)	H I
$\mathbf{B}_9$	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	$x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{10}$	$\left(\frac{1}{2} - x_3\right)\,\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{11}$	$\left(\frac{1}{2} - x_3\right)\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \left(\frac{1}{2} - x_3\right)\,\mathbf{a}_3$	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{12}$	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\,\mathbf{a}_2 + \left(\frac{1}{2} - x_3\right)\,\mathbf{a}_3$	$x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{13}$	$\left(\frac{1}{2} + x_3\right)\,\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{14}$	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	$-x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{15}$	$\left(\frac{1}{2} + x_3\right)\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\frac{1}{2} + x_3\right)\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{16}$	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\,\mathbf{a}_2 + \left(\frac{1}{2} + x_3\right)\,\mathbf{a}_3$	$-x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{17}$	$x_4\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	$x_4a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(24*h*)	O II
$\mathbf{B}_{18}$	$\left(\frac{1}{2} - x_4\right)\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(24*h*)	O II
$\mathbf{B}_{19}$	$\frac{3}{4}\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	$\frac{3}{4}a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(24*h*)	O II
$\mathbf{B}_{20}$	$\frac{3}{4}\,\mathbf{a}_1 + \left(\frac{1}{2} - x_4\right)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	$\frac{3}{4}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(24*h*)	O II
$\mathbf{B}_{21}$	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	(24*h*)	O II
$\mathbf{B}_{22}$	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \left(\frac{1}{2} - x_4\right)\,\mathbf{a}_3$	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{z}}$	(24*h*)	O II
$\mathbf{B}_{23}$	$\frac{3}{4}\,\mathbf{a}_1 + \left(\frac{1}{2} + x_4\right)\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	$\frac{3}{4}a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(24*h*)	O II
$\mathbf{B}_{24}$	$\frac{3}{4}\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	$\frac{3}{4}a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(24*h*)	O II
$\mathbf{B}_{25}$	$\left(\frac{1}{2} + x_4\right)\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(24*h*)	O II
$\mathbf{B}_{26}$	$-x_4\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	$-x_4a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(24*h*)	O II
$\mathbf{B}_{27}$	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} - x_4a\,\hat{\mathbf{z}}$	(24*h*)	O II

$$\mathbf{B}_{28} = \tfrac{1}{4}\,\mathbf{a}_1 + \tfrac{3}{4}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_4\right)\mathbf{a}_3 = \tfrac{1}{4}a\,\hat{\mathbf{x}} + \tfrac{3}{4}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_4\right)a\,\hat{\mathbf{z}} \quad (24h) \quad \text{O II}$$

$$\mathbf{B}_{29} = -x_4\,\mathbf{a}_1 + \tfrac{3}{4}\,\mathbf{a}_2 + \tfrac{1}{4}\,\mathbf{a}_3 = -x_4 a\,\hat{\mathbf{x}} + \tfrac{3}{4}a\,\hat{\mathbf{y}} + \tfrac{1}{4}a\,\hat{\mathbf{z}} \quad (24h) \quad \text{O II}$$

$$\mathbf{B}_{30} = \left(\tfrac{1}{2} + x_4\right)\mathbf{a}_1 + \tfrac{3}{4}\,\mathbf{a}_2 + \tfrac{1}{4}\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \tfrac{3}{4}a\,\hat{\mathbf{y}} + \tfrac{1}{4}a\,\hat{\mathbf{z}} \quad (24h) \quad \text{O II}$$

$$\mathbf{B}_{31} = \tfrac{1}{4}\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \tfrac{3}{4}\,\mathbf{a}_3 = \tfrac{1}{4}a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}} + \tfrac{3}{4}a\,\hat{\mathbf{z}} \quad (24h) \quad \text{O II}$$

$$\mathbf{B}_{32} = \tfrac{1}{4}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_4\right)\mathbf{a}_2 + \tfrac{3}{4}\,\mathbf{a}_3 = \tfrac{1}{4}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_4\right)a\,\hat{\mathbf{y}} + \tfrac{3}{4}a\,\hat{\mathbf{z}} \quad (24h) \quad \text{O II}$$

$$\mathbf{B}_{33} = \tfrac{3}{4}\,\mathbf{a}_1 + \tfrac{1}{4}\,\mathbf{a}_2 - x_4\,\mathbf{a}_3 = \tfrac{3}{4}a\,\hat{\mathbf{x}} + \tfrac{1}{4}a\,\hat{\mathbf{y}} - x_4 a\,\hat{\mathbf{z}} \quad (24h) \quad \text{O II}$$

$$\mathbf{B}_{34} = \tfrac{3}{4}\,\mathbf{a}_1 + \tfrac{1}{4}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_4\right)\mathbf{a}_3 = \tfrac{3}{4}a\,\hat{\mathbf{x}} + \tfrac{1}{4}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_4\right)a\,\hat{\mathbf{z}} \quad (24h) \quad \text{O II}$$

$$\mathbf{B}_{35} = \tfrac{1}{4}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_4\right)\mathbf{a}_2 + \tfrac{3}{4}\,\mathbf{a}_3 = \tfrac{1}{4}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_4\right)a\,\hat{\mathbf{y}} + \tfrac{3}{4}a\,\hat{\mathbf{z}} \quad (24h) \quad \text{O II}$$

$$\mathbf{B}_{36} = \tfrac{1}{4}\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \tfrac{3}{4}\,\mathbf{a}_3 = \tfrac{1}{4}a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}} + \tfrac{3}{4}a\,\hat{\mathbf{z}} \quad (24h) \quad \text{O II}$$

$$\mathbf{B}_{37} = \left(\tfrac{1}{2} - x_4\right)\mathbf{a}_1 + \tfrac{3}{4}\,\mathbf{a}_2 + \tfrac{1}{4}\,\mathbf{a}_3 = \left(\tfrac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \tfrac{3}{4}a\,\hat{\mathbf{y}} + \tfrac{1}{4}a\,\hat{\mathbf{z}} \quad (24h) \quad \text{O II}$$

$$\mathbf{B}_{38} = x_4\,\mathbf{a}_1 + \tfrac{3}{4}\,\mathbf{a}_2 + \tfrac{1}{4}\,\mathbf{a}_3 = x_4 a\,\hat{\mathbf{x}} + \tfrac{3}{4}a\,\hat{\mathbf{y}} + \tfrac{1}{4}a\,\hat{\mathbf{z}} \quad (24h) \quad \text{O II}$$

$$\mathbf{B}_{39} = \tfrac{3}{4}\,\mathbf{a}_1 + \tfrac{1}{4}\,\mathbf{a}_2 + x_4\,\mathbf{a}_3 = \tfrac{3}{4}a\,\hat{\mathbf{x}} + \tfrac{1}{4}a\,\hat{\mathbf{y}} + x_4 a\,\hat{\mathbf{z}} \quad (24h) \quad \text{O II}$$

$$\mathbf{B}_{40} = \tfrac{3}{4}\,\mathbf{a}_1 + \tfrac{1}{4}\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_4\right)\mathbf{a}_3 = \tfrac{3}{4}a\,\hat{\mathbf{x}} + \tfrac{1}{4}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_4\right)a\,\hat{\mathbf{z}} \quad (24h) \quad \text{O II}$$

$$\mathbf{B}_{41} = x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3 = x_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + z_5 a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{42} = \left(\tfrac{1}{2} - x_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_5\right)\mathbf{a}_2 + z_5\,\mathbf{a}_3 = \left(\tfrac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_5\right)a\,\hat{\mathbf{y}} + z_5 a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{43} = \left(\tfrac{1}{2} - x_5\right)\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_5\right)a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{44} = x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_5\right)\mathbf{a}_3 = x_5 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_5\right)a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{45} = z_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3 = z_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{46} = z_5\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_5\right)\mathbf{a}_3 = z_5 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_5\right)a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{47} = \left(\tfrac{1}{2} - z_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_5\right)\mathbf{a}_2 + x_5\,\mathbf{a}_3 = \left(\tfrac{1}{2} - z_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_5\right)a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{48} = \left(\tfrac{1}{2} - z_5\right)\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - z_5\right)a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_5\right)a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{49} = x_5\,\mathbf{a}_1 + z_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3 = x_5 a\,\hat{\mathbf{x}} + z_5 a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{50} = \left(\tfrac{1}{2} - x_5\right)\mathbf{a}_1 + z_5\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + z_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_5\right)a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{51} = x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2} - z_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_5\right)\mathbf{a}_3 = x_5 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - z_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_5\right)a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{52} = \left(\tfrac{1}{2} - x_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - z_5\right)\mathbf{a}_2 + x_5\,\mathbf{a}_3 = \left(\tfrac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - z_5\right)a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{53} = \left(\tfrac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_5\right)a\,\hat{\mathbf{y}} - z_5 a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{54} = -x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3 = -x_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} - z_5 a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{55} = \left(\tfrac{1}{2} + x_5\right)\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_5\right)a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{56} = -x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_3 = -x_5 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_5\right)a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{57} = \left(\tfrac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_2 - x_5\,\mathbf{a}_3 = \left(\tfrac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + z_5\right)a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{58} = -x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_5\right)\mathbf{a}_3 = -x_5 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + z_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_5\right)a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{59} = -x_5\,\mathbf{a}_1 - z_5\,\mathbf{a}_2 - x_5\,\mathbf{a}_3 = -x_5 a\,\hat{\mathbf{x}} - z_5 a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{60} = \left(\tfrac{1}{2} + x_5\right)\mathbf{a}_1 - z_5\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_5\right)a\,\hat{\mathbf{x}} - z_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_5\right)a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{61} = \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_5\right)\mathbf{a}_2 - x_5\,\mathbf{a}_3 = \left(\tfrac{1}{2} + z_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_5\right)a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{62} = \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + z_5\right)a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_5\right)a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{63} = -z_5\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_5\right)\mathbf{a}_3 = -z_5 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_5\right)a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

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\begin{array}{lcccccc}
\mathbf{B}_{64} & = & -z_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - x_5\,\mathbf{a}_3 & = & -z_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}} & (24k) & \text{O III} \\
\mathbf{B}_{65} & = & x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 & = & x_6 a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} + z_6 a\,\hat{\mathbf{z}} & (24k) & \text{O IV} \\
\mathbf{B}_{66} & = & \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_2 + z_6\,\mathbf{a}_3 & = & \left(\tfrac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_6\right) a\,\hat{\mathbf{y}} + z_6 a\,\hat{\mathbf{z}} & (24k) & \text{O IV} \\
\mathbf{B}_{67} & = & \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_1 + x_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_6\right) a\,\hat{\mathbf{z}} & (24k) & \text{O IV} \\
\mathbf{B}_{68} & = & x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_3 & = & x_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_6\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_6\right) a\,\hat{\mathbf{z}} & (24k) & \text{O IV} \\
\mathbf{B}_{69} & = & z_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + x_6\,\mathbf{a}_3 & = & z_6 a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} + x_6 a\,\hat{\mathbf{z}} & (24k) & \text{O IV} \\
\mathbf{B}_{70} & = & z_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_3 & = & z_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_6\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_6\right) a\,\hat{\mathbf{z}} & (24k) & \text{O IV} \\
\mathbf{B}_{71} & = & \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_2 + x_6\,\mathbf{a}_3 & = & \left(\tfrac{1}{2} - z_6\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_6\right) a\,\hat{\mathbf{y}} + x_6 a\,\hat{\mathbf{z}} & (24k) & \text{O IV} \\
\mathbf{B}_{72} & = & \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_1 + x_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} - z_6\right) a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_6\right) a\,\hat{\mathbf{z}} & (24k) & \text{O IV} \\
\mathbf{B}_{73} & = & x_6\,\mathbf{a}_1 + z_6\,\mathbf{a}_2 + x_6\,\mathbf{a}_3 & = & x_6 a\,\hat{\mathbf{x}} + z_6 a\,\hat{\mathbf{y}} + x_6 a\,\hat{\mathbf{z}} & (24k) & \text{O IV} \\
\mathbf{B}_{74} & = & \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_1 + z_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} + z_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_6\right) a\,\hat{\mathbf{z}} & (24k) & \text{O IV} \\
\mathbf{B}_{75} & = & x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_3 & = & x_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - z_6\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_6\right) a\,\hat{\mathbf{z}} & (24k) & \text{O IV} \\
\mathbf{B}_{76} & = & \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_2 + x_6\,\mathbf{a}_3 & = & \left(\tfrac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - z_6\right) a\,\hat{\mathbf{y}} + x_6 a\,\hat{\mathbf{z}} & (24k) & \text{O IV} \\
\mathbf{B}_{77} & = & \left(\tfrac{1}{2} + x_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3 & = & \left(\tfrac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_6\right) a\,\hat{\mathbf{y}} - z_6 a\,\hat{\mathbf{z}} & (24k) & \text{O IV} \\
\mathbf{B}_{78} & = & -x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3 & = & -x_6 a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} - z_6 a\,\hat{\mathbf{z}} & (24k) & \text{O IV} \\
\mathbf{B}_{79} & = & \left(\tfrac{1}{2} + x_6\right)\mathbf{a}_1 - x_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right) a\,\hat{\mathbf{z}} & (24k) & \text{O IV} \\
\mathbf{B}_{80} & = & -x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 & = & -x_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_6\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right) a\,\hat{\mathbf{z}} & (24k) & \text{O IV} \\
\mathbf{B}_{81} & = & \left(\tfrac{1}{2} + x_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_2 - x_6\,\mathbf{a}_3 & = & \left(\tfrac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + z_6\right) a\,\hat{\mathbf{y}} - x_6 a\,\hat{\mathbf{z}} & (24k) & \text{O IV} \\
\mathbf{B}_{82} & = & -x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_6\right)\mathbf{a}_3 & = & -x_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + z_6\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_6\right) a\,\hat{\mathbf{z}} & (24k) & \text{O IV} \\
\mathbf{B}_{83} & = & -x_6\,\mathbf{a}_1 - z_6\,\mathbf{a}_2 - x_6\,\mathbf{a}_3 & = & -x_6 a\,\hat{\mathbf{x}} - z_6 a\,\hat{\mathbf{y}} - x_6 a\,\hat{\mathbf{z}} & (24k) & \text{O IV} \\
\mathbf{B}_{84} & = & \left(\tfrac{1}{2} + x_6\right)\mathbf{a}_1 - z_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_6\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} - z_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_6\right) a\,\hat{\mathbf{z}} & (24k) & \text{O IV} \\
\mathbf{B}_{85} & = & \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_6\right)\mathbf{a}_2 - x_6\,\mathbf{a}_3 & = & \left(\tfrac{1}{2} + z_6\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_6\right) a\,\hat{\mathbf{y}} - x_6 a\,\hat{\mathbf{z}} & (24k) & \text{O IV} \\
\mathbf{B}_{86} & = & \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_1 - x_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_6\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} + z_6\right) a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_6\right) a\,\hat{\mathbf{z}} & (24k) & \text{O IV} \\
\mathbf{B}_{87} & = & -z_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_6\right)\mathbf{a}_3 & = & -z_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_6\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_6\right) a\,\hat{\mathbf{z}} & (24k) & \text{O IV} \\
\mathbf{B}_{88} & = & -z_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 - x_6\,\mathbf{a}_3 & = & -z_6 a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} - x_6 a\,\hat{\mathbf{z}} & (24k) & \text{O IV} \\
\mathbf{B}_{89} & = & x_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 & = & x_7 a\,\hat{\mathbf{x}} + x_7 a\,\hat{\mathbf{y}} + z_7 a\,\hat{\mathbf{z}} & (24k) & \text{O V} \\
\mathbf{B}_{90} & = & \left(\tfrac{1}{2} - x_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_7\right)\mathbf{a}_2 + z_7\,\mathbf{a}_3 & = & \left(\tfrac{1}{2} - x_7\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_7\right) a\,\hat{\mathbf{y}} + z_7 a\,\hat{\mathbf{z}} & (24k) & \text{O V} \\
\mathbf{B}_{91} & = & \left(\tfrac{1}{2} - x_7\right)\mathbf{a}_1 + x_7\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_7\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} - x_7\right) a\,\hat{\mathbf{x}} + x_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_7\right) a\,\hat{\mathbf{z}} & (24k) & \text{O V} \\
\mathbf{B}_{92} & = & x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_7\right)\mathbf{a}_3 & = & x_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_7\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_7\right) a\,\hat{\mathbf{z}} & (24k) & \text{O V} \\
\mathbf{B}_{93} & = & z_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + x_7\,\mathbf{a}_3 & = & z_7 a\,\hat{\mathbf{x}} + x_7 a\,\hat{\mathbf{y}} + x_7 a\,\hat{\mathbf{z}} & (24k) & \text{O V} \\
\mathbf{B}_{94} & = & z_7\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_7\right)\mathbf{a}_3 & = & z_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_7\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_7\right) a\,\hat{\mathbf{z}} & (24k) & \text{O V} \\
\mathbf{B}_{95} & = & \left(\tfrac{1}{2} - z_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_7\right)\mathbf{a}_2 + x_7\,\mathbf{a}_3 & = & \left(\tfrac{1}{2} - z_7\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_7\right) a\,\hat{\mathbf{y}} + x_7 a\,\hat{\mathbf{z}} & (24k) & \text{O V} \\
\mathbf{B}_{96} & = & \left(\tfrac{1}{2} - z_7\right)\mathbf{a}_1 + x_7\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_7\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} - z_7\right) a\,\hat{\mathbf{x}} + x_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_7\right) a\,\hat{\mathbf{z}} & (24k) & \text{O V} \\
\mathbf{B}_{97} & = & x_7\,\mathbf{a}_1 + z_7\,\mathbf{a}_2 + x_7\,\mathbf{a}_3 & = & x_7 a\,\hat{\mathbf{x}} + z_7 a\,\hat{\mathbf{y}} + x_7 a\,\hat{\mathbf{z}} & (24k) & \text{O V} \\
\mathbf{B}_{98} & = & \left(\tfrac{1}{2} - x_7\right)\mathbf{a}_1 + z_7\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_7\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} - x_7\right) a\,\hat{\mathbf{x}} + z_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_7\right) a\,\hat{\mathbf{z}} & (24k) & \text{O V} \\
\mathbf{B}_{99} & = & x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2} - z_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_7\right)\mathbf{a}_3 & = & x_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - z_7\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_7\right) a\,\hat{\mathbf{z}} & (24k) & \text{O V}
\end{array}
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\begin{aligned}
\mathbf{B}_{100} &= \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}-z_7\right)\mathbf{a}_2+x_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-z_7\right)a\,\hat{\mathbf{y}}+x_7a\,\hat{\mathbf{z}} && (24k) && \text{O V}\\
\mathbf{B}_{101} &= \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_7\right)\mathbf{a}_2-z_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{y}}-z_7a\,\hat{\mathbf{z}} && (24k) && \text{O V}\\
\mathbf{B}_{102} &= -x_7\,\mathbf{a}_1-x_7\,\mathbf{a}_2-z_7\,\mathbf{a}_3 &&= -x_7a\,\hat{\mathbf{x}}-x_7a\,\hat{\mathbf{y}}-z_7a\,\hat{\mathbf{z}} && (24k) && \text{O V}\\
\mathbf{B}_{103} &= \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1-x_7\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{x}}-x_7a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{z}} && (24k) && \text{O V}\\
\mathbf{B}_{104} &= -x_7\,\mathbf{a}_1+\left(\tfrac{1}{2}+x_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 &&= -x_7a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{z}} && (24k) && \text{O V}\\
\mathbf{B}_{105} &= \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}+z_7\right)\mathbf{a}_2-x_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{y}}-x_7a\,\hat{\mathbf{z}} && (24k) && \text{O V}\\
\mathbf{B}_{106} &= -x_7\,\mathbf{a}_1+\left(\tfrac{1}{2}+z_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_7\right)\mathbf{a}_3 &&= -x_7a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{z}} && (24k) && \text{O V}\\
\mathbf{B}_{107} &= -x_7\,\mathbf{a}_1-z_7\,\mathbf{a}_2-x_7\,\mathbf{a}_3 &&= -x_7a\,\hat{\mathbf{x}}-z_7a\,\hat{\mathbf{y}}-x_7a\,\hat{\mathbf{z}} && (24k) && \text{O V}\\
\mathbf{B}_{108} &= \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1-z_7\,\mathbf{a}_2+\left(\tfrac{1}{2}+x_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{x}}-z_7a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{z}} && (24k) && \text{O V}\\
\mathbf{B}_{109} &= \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_7\right)\mathbf{a}_2-x_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{y}}-x_7a\,\hat{\mathbf{z}} && (24k) && \text{O V}\\
\mathbf{B}_{110} &= \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_1-x_7\,\mathbf{a}_2+\left(\tfrac{1}{2}+x_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{x}}-x_7a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{z}} && (24k) && \text{O V}\\
\mathbf{B}_{111} &= -z_7\,\mathbf{a}_1+\left(\tfrac{1}{2}+x_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_7\right)\mathbf{a}_3 &&= -z_7a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{z}} && (24k) && \text{O V}\\
\mathbf{B}_{112} &= -z_7\,\mathbf{a}_1-x_7\,\mathbf{a}_2-x_7\,\mathbf{a}_3 &&= -z_7a\,\hat{\mathbf{x}}-x_7a\,\hat{\mathbf{y}}-x_7a\,\hat{\mathbf{z}} && (24k) && \text{O V}\\
\mathbf{B}_{113} &= x_8\,\mathbf{a}_1+x_8\,\mathbf{a}_2+z_8\,\mathbf{a}_3 &&= x_8a\,\hat{\mathbf{x}}+x_8a\,\hat{\mathbf{y}}+z_8a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{114} &= \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_1+\left(\tfrac{1}{2}-x_8\right)\mathbf{a}_2+z_8\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{y}}+z_8a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{115} &= \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_1+x_8\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_8\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{x}}+x_8a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_8\right)a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{116} &= x_8\,\mathbf{a}_1+\left(\tfrac{1}{2}-x_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_8\right)\mathbf{a}_3 &&= x_8a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_8\right)a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{117} &= z_8\,\mathbf{a}_1+x_8\,\mathbf{a}_2+x_8\,\mathbf{a}_3 &&= z_8a\,\hat{\mathbf{x}}+x_8a\,\hat{\mathbf{y}}+x_8a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{118} &= z_8\,\mathbf{a}_1+\left(\tfrac{1}{2}-x_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}-x_8\right)\mathbf{a}_3 &&= z_8a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{119} &= \left(\tfrac{1}{2}-z_8\right)\mathbf{a}_1+\left(\tfrac{1}{2}-x_8\right)\mathbf{a}_2+x_8\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-z_8\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{y}}+x_8a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{120} &= \left(\tfrac{1}{2}-z_8\right)\mathbf{a}_1+x_8\,\mathbf{a}_2+\left(\tfrac{1}{2}-x_8\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-z_8\right)a\,\hat{\mathbf{x}}+x_8a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{121} &= x_8\,\mathbf{a}_1+z_8\,\mathbf{a}_2+x_8\,\mathbf{a}_3 &&= x_8a\,\hat{\mathbf{x}}+z_8a\,\hat{\mathbf{y}}+x_8a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{122} &= \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_1+z_8\,\mathbf{a}_2+\left(\tfrac{1}{2}-x_8\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{x}}+z_8a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{123} &= x_8\,\mathbf{a}_1+\left(\tfrac{1}{2}-z_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}-x_8\right)\mathbf{a}_3 &&= x_8a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-z_8\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{124} &= \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_1+\left(\tfrac{1}{2}-z_8\right)\mathbf{a}_2+x_8\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-z_8\right)a\,\hat{\mathbf{y}}+x_8a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{125} &= \left(\tfrac{1}{2}+x_8\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_8\right)\mathbf{a}_2-z_8\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{y}}-z_8a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{126} &= -x_8\,\mathbf{a}_1-x_8\,\mathbf{a}_2-z_8\,\mathbf{a}_3 &&= -x_8a\,\hat{\mathbf{x}}-x_8a\,\hat{\mathbf{y}}-z_8a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{127} &= \left(\tfrac{1}{2}+x_8\right)\mathbf{a}_1-x_8\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_8\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{x}}-x_8a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_8\right)a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{128} &= -x_8\,\mathbf{a}_1+\left(\tfrac{1}{2}+x_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_8\right)\mathbf{a}_3 &&= -x_8a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_8\right)a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{129} &= \left(\tfrac{1}{2}+x_8\right)\mathbf{a}_1+\left(\tfrac{1}{2}+z_8\right)\mathbf{a}_2-x_8\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+z_8\right)a\,\hat{\mathbf{y}}-x_8a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{130} &= -x_8\,\mathbf{a}_1+\left(\tfrac{1}{2}+z_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_8\right)\mathbf{a}_3 &&= -x_8a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+z_8\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{131} &= -x_8\,\mathbf{a}_1-z_8\,\mathbf{a}_2-x_8\,\mathbf{a}_3 &&= -x_8a\,\hat{\mathbf{x}}-z_8a\,\hat{\mathbf{y}}-x_8a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{132} &= \left(\tfrac{1}{2}+x_8\right)\mathbf{a}_1-z_8\,\mathbf{a}_2+\left(\tfrac{1}{2}+x_8\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{x}}-z_8a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{133} &= \left(\tfrac{1}{2}+z_8\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_8\right)\mathbf{a}_2-x_8\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+z_8\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{y}}-x_8a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{134} &= \left(\tfrac{1}{2}+z_8\right)\mathbf{a}_1-x_8\,\mathbf{a}_2+\left(\tfrac{1}{2}+x_8\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+z_8\right)a\,\hat{\mathbf{x}}-x_8a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{135} &= -z_8\,\mathbf{a}_1+\left(\tfrac{1}{2}+x_8\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_8\right)\mathbf{a}_3 &&= -z_8a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+x_8\right)a\,\hat{\mathbf{z}} && (24k) && \text{W}
\end{aligned}
$$

$$
\begin{array}{lcccll}
\mathbf{B}_{136} & = & -z_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 - x_8\,\mathbf{a}_3 & = & -z_8 a\,\hat{\mathbf{x}} - x_8 a\,\hat{\mathbf{y}} - x_8 a\,\hat{\mathbf{z}} & (24k) \quad \text{W} \\
\mathbf{B}_{137} & = & x_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 & = & x_9 a\,\hat{\mathbf{x}} + y_9 a\,\hat{\mathbf{y}} + z_9 a\,\hat{\mathbf{z}} & (48l) \quad \text{H II} \\
\mathbf{B}_{138} & = & \left(\tfrac{1}{2} - x_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_9\right)\mathbf{a}_2 + z_9\,\mathbf{a}_3 & = & \left(\tfrac{1}{2} - x_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_9\right)a\,\hat{\mathbf{y}} + z_9 a\,\hat{\mathbf{z}} & (48l) \quad \text{H II} \\
\mathbf{B}_{139} & = & \left(\tfrac{1}{2} - x_9\right)\mathbf{a}_1 + y_9\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_9\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} - x_9\right)a\,\hat{\mathbf{x}} + y_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_9\right)a\,\hat{\mathbf{z}} & (48l) \quad \text{H II} \\
\mathbf{B}_{140} & = & x_9\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_9\right)\mathbf{a}_3 & = & x_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_9\right)a\,\hat{\mathbf{z}} & (48l) \quad \text{H II} \\
\mathbf{B}_{141} & = & z_9\,\mathbf{a}_1 + x_9\,\mathbf{a}_2 + y_9\,\mathbf{a}_3 & = & z_9 a\,\hat{\mathbf{x}} + x_9 a\,\hat{\mathbf{y}} + y_9 a\,\hat{\mathbf{z}} & (48l) \quad \text{H II} \\
\mathbf{B}_{142} & = & z_9\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - y_9\right)\mathbf{a}_3 & = & z_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - y_9\right)a\,\hat{\mathbf{z}} & (48l) \quad \text{H II} \\
\mathbf{B}_{143} & = & \left(\tfrac{1}{2} - z_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_9\right)\mathbf{a}_2 + y_9\,\mathbf{a}_3 & = & \left(\tfrac{1}{2} - z_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_9\right)a\,\hat{\mathbf{y}} + y_9 a\,\hat{\mathbf{z}} & (48l) \quad \text{H II} \\
\mathbf{B}_{144} & = & \left(\tfrac{1}{2} - z_9\right)\mathbf{a}_1 + x_9\,\mathbf{a}_2 + \left(\tfrac{1}{2} - y_9\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} - z_9\right)a\,\hat{\mathbf{x}} + x_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - y_9\right)a\,\hat{\mathbf{z}} & (48l) \quad \text{H II} \\
\mathbf{B}_{145} & = & y_9\,\mathbf{a}_1 + z_9\,\mathbf{a}_2 + x_9\,\mathbf{a}_3 & = & y_9 a\,\hat{\mathbf{x}} + z_9 a\,\hat{\mathbf{y}} + x_9 a\,\hat{\mathbf{z}} & (48l) \quad \text{H II} \\
\mathbf{B}_{146} & = & \left(\tfrac{1}{2} - y_9\right)\mathbf{a}_1 + z_9\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_9\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} - y_9\right)a\,\hat{\mathbf{x}} + z_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_9\right)a\,\hat{\mathbf{z}} & (48l) \quad \text{H II} \\
\mathbf{B}_{147} & = & y_9\,\mathbf{a}_1 + \left(\tfrac{1}{2} - z_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_9\right)\mathbf{a}_3 & = & y_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - z_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_9\right)a\,\hat{\mathbf{z}} & (48l) \quad \text{H II} \\
\mathbf{B}_{148} & = & \left(\tfrac{1}{2} - y_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - z_9\right)\mathbf{a}_2 + x_9\,\mathbf{a}_3 & = & \left(\tfrac{1}{2} - y_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - z_9\right)a\,\hat{\mathbf{y}} + x_9 a\,\hat{\mathbf{z}} & (48l) \quad \text{H II} \\
\mathbf{B}_{149} & = & \left(\tfrac{1}{2} + y_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_9\right)\mathbf{a}_2 - z_9\,\mathbf{a}_3 & = & \left(\tfrac{1}{2} + y_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_9\right)a\,\hat{\mathbf{y}} - z_9 a\,\hat{\mathbf{z}} & (48l) \quad \text{H II} \\
\mathbf{B}_{150} & = & -y_9\,\mathbf{a}_1 - x_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3 & = & -y_9 a\,\hat{\mathbf{x}} - x_9 a\,\hat{\mathbf{y}} - z_9 a\,\hat{\mathbf{z}} & (48l) \quad \text{H II} \\
\mathbf{B}_{151} & = & \left(\tfrac{1}{2} + y_9\right)\mathbf{a}_1 - x_9\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} + y_9\right)a\,\hat{\mathbf{x}} - x_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_9\right)a\,\hat{\mathbf{z}} & (48l) \quad \text{H II} \\
\mathbf{B}_{152} & = & -y_9\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_3 & = & -y_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_9\right)a\,\hat{\mathbf{z}} & (48l) \quad \text{H II} \\
\mathbf{B}_{153} & = & \left(\tfrac{1}{2} + x_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_2 - y_9\,\mathbf{a}_3 & = & \left(\tfrac{1}{2} + x_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + z_9\right)a\,\hat{\mathbf{y}} - y_9 a\,\hat{\mathbf{z}} & (48l) \quad \text{H II} \\
\mathbf{B}_{154} & = & -x_9\,\mathbf{a}_1 + \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + y_9\right)\mathbf{a}_3 & = & -x_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + z_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + y_9\right)a\,\hat{\mathbf{z}} & (48l) \quad \text{H II} \\
\mathbf{B}_{155} & = & -x_9\,\mathbf{a}_1 - z_9\,\mathbf{a}_2 - y_9\,\mathbf{a}_3 & = & -x_9 a\,\hat{\mathbf{x}} - z_9 a\,\hat{\mathbf{y}} - y_9 a\,\hat{\mathbf{z}} & (48l) \quad \text{H II} \\
\mathbf{B}_{156} & = & \left(\tfrac{1}{2} + x_9\right)\mathbf{a}_1 - z_9\,\mathbf{a}_2 + \left(\tfrac{1}{2} + y_9\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} + x_9\right)a\,\hat{\mathbf{x}} - z_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + y_9\right)a\,\hat{\mathbf{z}} & (48l) \quad \text{H II} \\
\mathbf{B}_{157} & = & \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_9\right)\mathbf{a}_2 - x_9\,\mathbf{a}_3 & = & \left(\tfrac{1}{2} + z_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_9\right)a\,\hat{\mathbf{y}} - x_9 a\,\hat{\mathbf{z}} & (48l) \quad \text{H II} \\
\mathbf{B}_{158} & = & \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_1 - y_9\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_9\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} + z_9\right)a\,\hat{\mathbf{x}} - y_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_9\right)a\,\hat{\mathbf{z}} & (48l) \quad \text{H II} \\
\mathbf{B}_{159} & = & -z_9\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_9\right)\mathbf{a}_3 & = & -z_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_9\right)a\,\hat{\mathbf{z}} & (48l) \quad \text{H II} \\
\mathbf{B}_{160} & = & -z_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 - x_9\,\mathbf{a}_3 & = & -z_9 a\,\hat{\mathbf{x}} - y_9 a\,\hat{\mathbf{y}} - x_9 a\,\hat{\mathbf{z}} & (48l) \quad \text{H II} \\
\mathbf{B}_{161} & = & -x_9\,\mathbf{a}_1 - y_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3 & = & -x_9 a\,\hat{\mathbf{x}} - y_9 a\,\hat{\mathbf{y}} - z_9 a\,\hat{\mathbf{z}} & (48l) \quad \text{H II} \\
\mathbf{B}_{162} & = & \left(\tfrac{1}{2} + x_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_9\right)\mathbf{a}_2 - z_9\,\mathbf{a}_3 & = & \left(\tfrac{1}{2} + x_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_9\right)a\,\hat{\mathbf{y}} - z_9 a\,\hat{\mathbf{z}} & (48l) \quad \text{H II} \\
\mathbf{B}_{163} & = & \left(\tfrac{1}{2} + x_9\right)\mathbf{a}_1 - y_9\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} + x_9\right)a\,\hat{\mathbf{x}} - y_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_9\right)a\,\hat{\mathbf{z}} & (48l) \quad \text{H II} \\
\mathbf{B}_{164} & = & -x_9\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_3 & = & -x_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_9\right)a\,\hat{\mathbf{z}} & (48l) \quad \text{H II} \\
\mathbf{B}_{165} & = & -z_9\,\mathbf{a}_1 - x_9\,\mathbf{a}_2 - y_9\,\mathbf{a}_3 & = & -z_9 a\,\hat{\mathbf{x}} - x_9 a\,\hat{\mathbf{y}} - y_9 a\,\hat{\mathbf{z}} & (48l) \quad \text{H II} \\
\mathbf{B}_{166} & = & -z_9\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + y_9\right)\mathbf{a}_3 & = & -z_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + y_9\right)a\,\hat{\mathbf{z}} & (48l) \quad \text{H II} \\
\mathbf{B}_{167} & = & \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_9\right)\mathbf{a}_2 - y_9\,\mathbf{a}_3 & = & \left(\tfrac{1}{2} + z_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_9\right)a\,\hat{\mathbf{y}} - y_9 a\,\hat{\mathbf{z}} & (48l) \quad \text{H II} \\
\mathbf{B}_{168} & = & \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_1 - x_9\,\mathbf{a}_2 + \left(\tfrac{1}{2} + y_9\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} + z_9\right)a\,\hat{\mathbf{x}} - x_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + y_9\right)a\,\hat{\mathbf{z}} & (48l) \quad \text{H II} \\
\mathbf{B}_{169} & = & -y_9\,\mathbf{a}_1 - z_9\,\mathbf{a}_2 - x_9\,\mathbf{a}_3 & = & -y_9 a\,\hat{\mathbf{x}} - z_9 a\,\hat{\mathbf{y}} - x_9 a\,\hat{\mathbf{z}} & (48l) \quad \text{H II} \\
\mathbf{B}_{170} & = & \left(\tfrac{1}{2} + y_9\right)\mathbf{a}_1 - z_9\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_9\right)\mathbf{a}_3 & = & \left(\tfrac{1}{2} + y_9\right)a\,\hat{\mathbf{x}} - z_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_9\right)a\,\hat{\mathbf{z}} & (48l) \quad \text{H II} \\
\mathbf{B}_{171} & = & -y_9\,\mathbf{a}_1 + \left(\tfrac{1}{2} + z_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_9\right)\mathbf{a}_3 & = & -y_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + z_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_9\right)a\,\hat{\mathbf{z}} & (48l) \quad \text{H II}
\end{array}
$$

$$\mathbf{B}_{172} = \left(\tfrac{1}{2}+y_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+z_9\right)\mathbf{a}_2 - x_9\,\mathbf{a}_3 = \left(\tfrac{1}{2}+y_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+z_9\right)a\,\hat{\mathbf{y}} - x_9 a\,\hat{\mathbf{z}} \quad (48l) \quad \text{H II}$$

$$\mathbf{B}_{173} = \left(\tfrac{1}{2}-y_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-x_9\right)\mathbf{a}_2 + z_9\,\mathbf{a}_3 = \left(\tfrac{1}{2}-y_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_9\right)a\,\hat{\mathbf{y}} + z_9 a\,\hat{\mathbf{z}} \quad (48l) \quad \text{H II}$$

$$\mathbf{B}_{174} = y_9\,\mathbf{a}_1 + x_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 = y_9 a\,\hat{\mathbf{x}} + x_9 a\,\hat{\mathbf{y}} + z_9 a\,\hat{\mathbf{z}} \quad (48l) \quad \text{H II}$$

$$\mathbf{B}_{175} = \left(\tfrac{1}{2}-y_9\right)\mathbf{a}_1 + x_9\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-y_9\right)a\,\hat{\mathbf{x}} + x_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_9\right)a\,\hat{\mathbf{z}} \quad (48l) \quad \text{H II}$$

$$\mathbf{B}_{176} = y_9\,\mathbf{a}_1 + \left(\tfrac{1}{2}-x_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_9\right)\mathbf{a}_3 = y_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_9\right)a\,\hat{\mathbf{z}} \quad (48l) \quad \text{H II}$$

$$\mathbf{B}_{177} = \left(\tfrac{1}{2}-x_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-z_9\right)\mathbf{a}_2 + y_9\,\mathbf{a}_3 = \left(\tfrac{1}{2}-x_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-z_9\right)a\,\hat{\mathbf{y}} + y_9 a\,\hat{\mathbf{z}} \quad (48l) \quad \text{H II}$$

$$\mathbf{B}_{178} = x_9\,\mathbf{a}_1 + \left(\tfrac{1}{2}-z_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-y_9\right)\mathbf{a}_3 = x_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-z_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-y_9\right)a\,\hat{\mathbf{z}} \quad (48l) \quad \text{H II}$$

$$\mathbf{B}_{179} = x_9\,\mathbf{a}_1 + z_9\,\mathbf{a}_2 + y_9\,\mathbf{a}_3 = x_9 a\,\hat{\mathbf{x}} + z_9 a\,\hat{\mathbf{y}} + y_9 a\,\hat{\mathbf{z}} \quad (48l) \quad \text{H II}$$

$$\mathbf{B}_{180} = \left(\tfrac{1}{2}-x_9\right)\mathbf{a}_1 + z_9\,\mathbf{a}_2 + \left(\tfrac{1}{2}-y_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_9\right)a\,\hat{\mathbf{x}} + z_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-y_9\right)a\,\hat{\mathbf{z}} \quad (48l) \quad \text{H II}$$

$$\mathbf{B}_{181} = \left(\tfrac{1}{2}-z_9\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_9\right)\mathbf{a}_2 + x_9\,\mathbf{a}_3 = \left(\tfrac{1}{2}-z_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_9\right)a\,\hat{\mathbf{y}} + x_9 a\,\hat{\mathbf{z}} \quad (48l) \quad \text{H II}$$

$$\mathbf{B}_{182} = \left(\tfrac{1}{2}-z_9\right)\mathbf{a}_1 + y_9\,\mathbf{a}_2 + \left(\tfrac{1}{2}-x_9\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-z_9\right)a\,\hat{\mathbf{x}} + y_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-x_9\right)a\,\hat{\mathbf{z}} \quad (48l) \quad \text{H II}$$

$$\mathbf{B}_{183} = z_9\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_9\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-x_9\right)\mathbf{a}_3 = z_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-y_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-x_9\right)a\,\hat{\mathbf{z}} \quad (48l) \quad \text{H II}$$

$$\mathbf{B}_{184} = z_9\,\mathbf{a}_1 + y_9\,\mathbf{a}_2 + x_9\,\mathbf{a}_3 = z_9 a\,\hat{\mathbf{x}} + y_9 a\,\hat{\mathbf{y}} + x_9 a\,\hat{\mathbf{z}} \quad (48l) \quad \text{H II}$$

References:

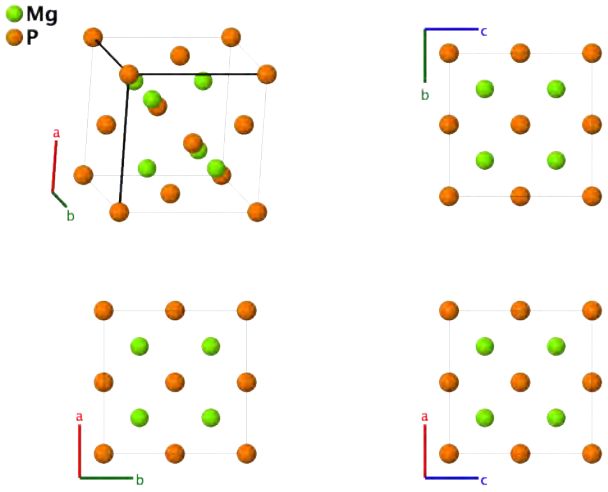
- G. M. Brown, M.-R. Noe-Spirlet, W. R. Busing, and H. A. Levy, *Dodecatungstophosphoric acid hexahydrate, $(H_5O_2^+)_3(PW_{12}O_{40}^{3-})$. The true structure of Keggin's 'pentahydrate' from single-crystal X-ray and neutron diffraction data*, Acta Crystallogr. Sect. B Struct. Sci. **33**, 1038–1046 (1977), doi:10.1107/S0567740877005330.

Geometry files:

- CIF: pp. 1806
- POSCAR: pp. 1807

Mg_3P_2 ($D5_5$) Structure: A3B2_cP10_224_d_b

http://aflow.org/prototype-encyclopedia/A3B2_cP10_224_d_b



Prototype	:	Mg_3P_2
AFLOW prototype label	:	A3B2_cP10_224_d_b
***Strukturbericht* designation**	:	$D5_5$
Pearson symbol	:	cP10
Space group number	:	224
Space group symbol	:	$Pn\bar{3}m$
AFLOW prototype command	:	`aflow --proto=A3B2_cP10_224_d_b --params=`a

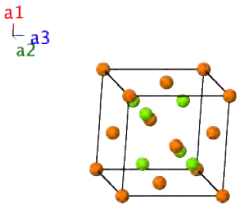
Other compounds with this structure

- Al_3P_2, Cd_3P_2, Zn_3P_2, Mg_3As_2, Zn_3As_2, and Ag_2O_3

- (Parthé, 1993) lists As_2O_3 as the prototype for $D5_5$.
- (Passerini, 1928) gives the atomic positions in setting 1 of space group $Pn\bar{3}m$ #224. We have shifted this to the standard, setting 2.

Simple Cubic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = a\,\hat{\mathbf{z}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(4*b*)	P
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$	(4*b*)	P
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(4*b*)	P
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(4*b*)	P
$\mathbf{B}_5$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(6*d*)	Mg
$\mathbf{B}_6$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(6*d*)	Mg
$\mathbf{B}_7$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(6*d*)	Mg
$\mathbf{B}_8$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(6*d*)	Mg
$\mathbf{B}_9$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(6*d*)	Mg
$\mathbf{B}_{10}$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(6*d*)	Mg

References:

- L. Passerini, *Struttura cristallina di alcuni fosfuri di metalli bivalenti e trivalenti*, Gazz. Chim. Ital. **58**, 655–664 (1928).
- E. Parthé, L. Gelato, B. Chabot, M. Penso, K. Cenzual, and R. Gladyshevskii, in *Standardized Data and Crystal Chemical Characterization of Inorganic Structure Types* (Springer-Verlag, Berlin, Heidelberg, 1993), *Gmelin Handbook of Inorganic and Organometallic Chemistry*, vol. 2, chap. Crystal Chemical Characterization of Inorganic Structure Types, 8 edn., doi:10.1007/978-3-662-02909-1_3.

Found in:

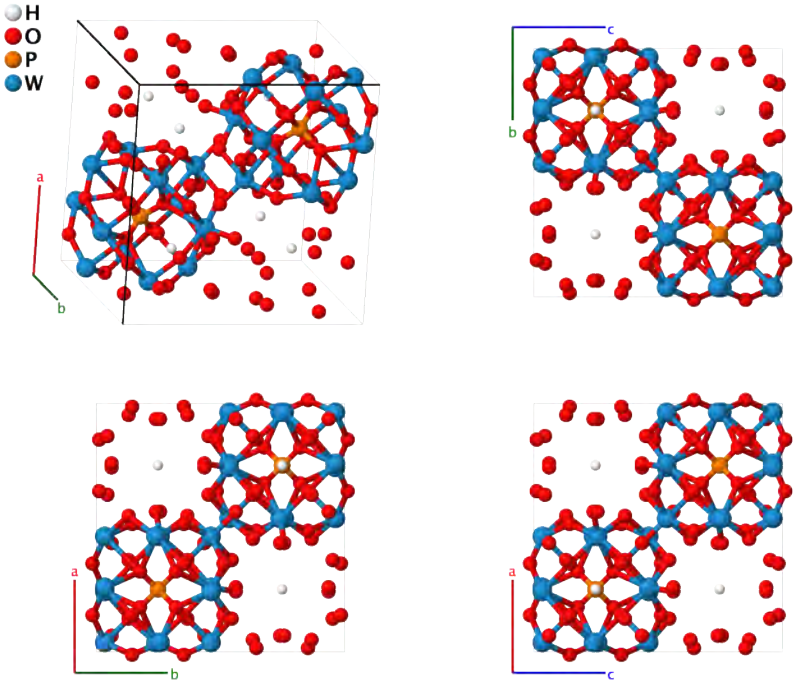
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1808
- POSCAR: pp. 1808

$H_3PW_{12}O_{40}\cdot 3H_2O$ Structure: A3B40CD12_cP112_224_d_e3k_a_k

http://aflow.org/prototype-encyclopedia/A3B40CD12_cP112_224_d_e3k_a_k

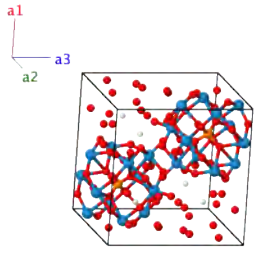


Prototype	:	$(H_3O)_3O_{40}PW_{12}$
AFLOW prototype label	:	A3B40CD12_cP112_224_d_e3k_a_k
***Strukturbericht* designation**	:	None
Pearson symbol	:	cP112
Space group number	:	224
Space group symbol	:	$Pn\bar{3}m$
AFLOW prototype command	:	`aflow --proto=A3B40CD12_cP112_224_d_e3k_a_k --params=`$a, x_3, x_4, z_4, x_5, z_5, x_6, z_6, x_7, z_7$

- This is a partially dehydrated version of $H_3PW_{12}O_{40}\cdot 29H_2O$ ($H4_{21}$) and $H_3PW_{12}O_{40}\cdot 5H_2O$ ($H4_{16}$) or $H_3PW_{12}O_{40}\cdot 6H_2O$. The six $H_5O_2^+$ ions of the latter structure have been replaced by three H_3O^+ ions.

Simple Cubic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = a\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(2*a*)	P
$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(2*a*)	P
$\mathbf{B}_3$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(6*d*)	H_3O
$\mathbf{B}_4$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(6*d*)	H_3O
$\mathbf{B}_5$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(6*d*)	H_3O
$\mathbf{B}_6$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(6*d*)	H_3O
$\mathbf{B}_7$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(6*d*)	H_3O
$\mathbf{B}_8$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(6*d*)	H_3O
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{11}$	=	$\left(\frac{1}{2} - x_3\right)\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{12}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_3\right)\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_3\right)a\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{13}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{14}$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{15}$	=	$\left(\frac{1}{2} + x_3\right)\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{16}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_3\right)\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_3\right)a\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{17}$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + z_4a\,\hat{\mathbf{z}}$	(24*k*)	O II
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{y}} + z_4a\,\hat{\mathbf{z}}$	(24*k*)	O II
$\mathbf{B}_{19}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)a\,\hat{\mathbf{z}}$	(24*k*)	O II
$\mathbf{B}_{20}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_4\right)a\,\hat{\mathbf{z}}$	(24*k*)	O II
$\mathbf{B}_{21}$	=	$z_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$z_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	(24*k*)	O II
$\mathbf{B}_{22}$	=	$z_4\,\mathbf{a}_1 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_3$	=	$z_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{z}}$	(24*k*)	O II
$\mathbf{B}_{23}$	=	$\left(\frac{1}{2} - z_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - z_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	(24*k*)	O II
$\mathbf{B}_{24}$	=	$\left(\frac{1}{2} - z_4\right)\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - z_4\right)a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{z}}$	(24*k*)	O II
$\mathbf{B}_{25}$	=	$x_4\,\mathbf{a}_1 + z_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + z_4a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	(24*k*)	O II
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + z_4\,\mathbf{a}_2 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + z_4a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{z}}$	(24*k*)	O II
$\mathbf{B}_{27}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - z_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{z}}$	(24*k*)	O II
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - z_4\right)\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - z_4\right)a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	(24*k*)	O II
$\mathbf{B}_{29}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_4\right)\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{y}} - z_4a\,\hat{\mathbf{z}}$	(24*k*)	O II
$\mathbf{B}_{30}$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - z_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} - z_4a\,\hat{\mathbf{z}}$	(24*k*)	O II
$\mathbf{B}_{31}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)a\,\hat{\mathbf{z}}$	(24*k*)	O II
$\mathbf{B}_{32}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + x_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_4\right)a\,\hat{\mathbf{z}}$	(24*k*)	O II
$\mathbf{B}_{33}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + z_4\right)a\,\hat{\mathbf{y}} - x_4a\,\hat{\mathbf{z}}$	(24*k*)	O II
$\mathbf{B}_{34}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + z_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_4\right)\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + z_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{z}}$	(24*k*)	O II

$$\mathbf{B}_{35} = -x_4\,\mathbf{a}_1 - z_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3 = -x_4 a\,\hat{\mathbf{x}} - z_4 a\,\hat{\mathbf{y}} - x_4 a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O II}$$

$$\mathbf{B}_{36} = \left(\tfrac{1}{2}+x_4\right)\mathbf{a}_1 - z_4\,\mathbf{a}_2 + \left(\tfrac{1}{2}+x_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_4\right)a\,\hat{\mathbf{x}} - z_4 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_4\right)a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O II}$$

$$\mathbf{B}_{37} = \left(\tfrac{1}{2}+z_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+x_4\right)\mathbf{a}_2 - x_4\,\mathbf{a}_3 = \left(\tfrac{1}{2}+z_4\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_4\right)a\,\hat{\mathbf{y}} - x_4 a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O II}$$

$$\mathbf{B}_{38} = \left(\tfrac{1}{2}+z_4\right)\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \left(\tfrac{1}{2}+x_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+z_4\right)a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_4\right)a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O II}$$

$$\mathbf{B}_{39} = -z_4\,\mathbf{a}_1 + \left(\tfrac{1}{2}+x_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+x_4\right)\mathbf{a}_3 = -z_4 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_4\right)a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O II}$$

$$\mathbf{B}_{40} = -z_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3 = -z_4 a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}} - x_4 a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O II}$$

$$\mathbf{B}_{41} = x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3 = x_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + z_5 a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{42} = \left(\tfrac{1}{2}-x_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-x_5\right)\mathbf{a}_2 + z_5\,\mathbf{a}_3 = \left(\tfrac{1}{2}-x_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_5\right)a\,\hat{\mathbf{y}} + z_5 a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{43} = \left(\tfrac{1}{2}-x_5\right)\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_5\right)a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_5\right)a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{44} = x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2}-x_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_5\right)\mathbf{a}_3 = x_5 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_5\right)a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{45} = z_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3 = z_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{46} = z_5\,\mathbf{a}_1 + \left(\tfrac{1}{2}-x_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-x_5\right)\mathbf{a}_3 = z_5 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-x_5\right)a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{47} = \left(\tfrac{1}{2}-z_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-x_5\right)\mathbf{a}_2 + x_5\,\mathbf{a}_3 = \left(\tfrac{1}{2}-z_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_5\right)a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{48} = \left(\tfrac{1}{2}-z_5\right)\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \left(\tfrac{1}{2}-x_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-z_5\right)a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-x_5\right)a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{49} = x_5\,\mathbf{a}_1 + z_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3 = x_5 a\,\hat{\mathbf{x}} + z_5 a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{50} = \left(\tfrac{1}{2}-x_5\right)\mathbf{a}_1 + z_5\,\mathbf{a}_2 + \left(\tfrac{1}{2}-x_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_5\right)a\,\hat{\mathbf{x}} + z_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-x_5\right)a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{51} = x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2}-z_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-x_5\right)\mathbf{a}_3 = x_5 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-z_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-x_5\right)a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{52} = \left(\tfrac{1}{2}-x_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-z_5\right)\mathbf{a}_2 + x_5\,\mathbf{a}_3 = \left(\tfrac{1}{2}-x_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-z_5\right)a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{53} = \left(\tfrac{1}{2}+x_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+x_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3 = \left(\tfrac{1}{2}+x_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_5\right)a\,\hat{\mathbf{y}} - z_5 a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{54} = -x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3 = -x_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} - z_5 a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{55} = \left(\tfrac{1}{2}+x_5\right)\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_5\right)a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_5\right)a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{56} = -x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2}+x_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_5\right)\mathbf{a}_3 = -x_5 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_5\right)a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{57} = \left(\tfrac{1}{2}+x_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+z_5\right)\mathbf{a}_2 - x_5\,\mathbf{a}_3 = \left(\tfrac{1}{2}+x_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+z_5\right)a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{58} = -x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2}+z_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+x_5\right)\mathbf{a}_3 = -x_5 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+z_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_5\right)a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{59} = -x_5\,\mathbf{a}_1 - z_5\,\mathbf{a}_2 - x_5\,\mathbf{a}_3 = -x_5 a\,\hat{\mathbf{x}} - z_5 a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{60} = \left(\tfrac{1}{2}+x_5\right)\mathbf{a}_1 - z_5\,\mathbf{a}_2 + \left(\tfrac{1}{2}+x_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_5\right)a\,\hat{\mathbf{x}} - z_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_5\right)a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{61} = \left(\tfrac{1}{2}+z_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+x_5\right)\mathbf{a}_2 - x_5\,\mathbf{a}_3 = \left(\tfrac{1}{2}+z_5\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_5\right)a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{62} = \left(\tfrac{1}{2}+z_5\right)\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \left(\tfrac{1}{2}+x_5\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+z_5\right)a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_5\right)a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{63} = -z_5\,\mathbf{a}_1 + \left(\tfrac{1}{2}+x_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+x_5\right)\mathbf{a}_3 = -z_5 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_5\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_5\right)a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{64} = -z_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - x_5\,\mathbf{a}_3 = -z_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O III}$$

$$\mathbf{B}_{65} = x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 = x_6 a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} + z_6 a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O IV}$$

$$\mathbf{B}_{66} = \left(\tfrac{1}{2}-x_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-x_6\right)\mathbf{a}_2 + z_6\,\mathbf{a}_3 = \left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{y}} + z_6 a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O IV}$$

$$\mathbf{B}_{67} = \left(\tfrac{1}{2}-x_6\right)\mathbf{a}_1 + x_6\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_6\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_6\right)a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O IV}$$

$$\mathbf{B}_{68} = x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2}-x_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_6\right)\mathbf{a}_3 = x_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_6\right)a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O IV}$$

$$\mathbf{B}_{69} = z_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + x_6\,\mathbf{a}_3 = z_6 a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} + x_6 a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O IV}$$

$$\mathbf{B}_{70} = z_6\,\mathbf{a}_1 + \left(\tfrac{1}{2}-x_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-x_6\right)\mathbf{a}_3 = z_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{z}} \quad (24k) \quad \text{O IV}$$

$$
\begin{aligned}
\mathbf{B}_{71} &= \left(\tfrac{1}{2}-z_6\right)\mathbf{a}_1+\left(\tfrac{1}{2}-x_6\right)\mathbf{a}_2+x_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-z_6\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{y}}+x_6a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{72} &= \left(\tfrac{1}{2}-z_6\right)\mathbf{a}_1+x_6\,\mathbf{a}_2+\left(\tfrac{1}{2}-x_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-z_6\right)a\,\hat{\mathbf{x}}+x_6a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{73} &= x_6\,\mathbf{a}_1+z_6\,\mathbf{a}_2+x_6\,\mathbf{a}_3 &&= x_6a\,\hat{\mathbf{x}}+z_6a\,\hat{\mathbf{y}}+x_6a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{74} &= \left(\tfrac{1}{2}-x_6\right)\mathbf{a}_1+z_6\,\mathbf{a}_2+\left(\tfrac{1}{2}-x_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{x}}+z_6a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{75} &= x_6\,\mathbf{a}_1+\left(\tfrac{1}{2}-z_6\right)\mathbf{a}_2+\left(\tfrac{1}{2}-x_6\right)\mathbf{a}_3 &&= x_6a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-z_6\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{76} &= \left(\tfrac{1}{2}-x_6\right)\mathbf{a}_1+\left(\tfrac{1}{2}-z_6\right)\mathbf{a}_2+x_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_6\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-z_6\right)a\,\hat{\mathbf{y}}+x_6a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{77} &= \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_6\right)\mathbf{a}_2-z_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{y}}-z_6a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{78} &= -x_6\,\mathbf{a}_1-x_6\,\mathbf{a}_2-z_6\,\mathbf{a}_3 &&= -x_6a\,\hat{\mathbf{x}}-x_6a\,\hat{\mathbf{y}}-z_6a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{79} &= \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_1-x_6\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{x}}-x_6a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_6\right)a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{80} &= -x_6\,\mathbf{a}_1+\left(\tfrac{1}{2}+x_6\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_6\right)\mathbf{a}_3 &&= -x_6a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_6\right)a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{81} &= \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_1+\left(\tfrac{1}{2}+z_6\right)\mathbf{a}_2-x_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+z_6\right)a\,\hat{\mathbf{y}}-x_6a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{82} &= -x_6\,\mathbf{a}_1+\left(\tfrac{1}{2}+z_6\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_6\right)\mathbf{a}_3 &&= -x_6a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+z_6\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{83} &= -x_6\,\mathbf{a}_1-z_6\,\mathbf{a}_2-x_6\,\mathbf{a}_3 &&= -x_6a\,\hat{\mathbf{x}}-z_6a\,\hat{\mathbf{y}}-x_6a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{84} &= \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_1-z_6\,\mathbf{a}_2+\left(\tfrac{1}{2}+x_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{x}}-z_6a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{85} &= \left(\tfrac{1}{2}+z_6\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_6\right)\mathbf{a}_2-x_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+z_6\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{y}}-x_6a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{86} &= \left(\tfrac{1}{2}+z_6\right)\mathbf{a}_1-x_6\,\mathbf{a}_2+\left(\tfrac{1}{2}+x_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+z_6\right)a\,\hat{\mathbf{x}}-x_6a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{87} &= -z_6\,\mathbf{a}_1+\left(\tfrac{1}{2}+x_6\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_6\right)\mathbf{a}_3 &&= -z_6a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{88} &= -z_6\,\mathbf{a}_1-x_6\,\mathbf{a}_2-x_6\,\mathbf{a}_3 &&= -z_6a\,\hat{\mathbf{x}}-x_6a\,\hat{\mathbf{y}}-x_6a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{89} &= x_7\,\mathbf{a}_1+x_7\,\mathbf{a}_2+z_7\,\mathbf{a}_3 &&= x_7a\,\hat{\mathbf{x}}+x_7a\,\hat{\mathbf{y}}+z_7a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{90} &= \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}-x_7\right)\mathbf{a}_2+z_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{y}}+z_7a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{91} &= \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_1+x_7\,\mathbf{a}_2+\left(\tfrac{1}{2}-z_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{x}}+x_7a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_7\right)a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{92} &= x_7\,\mathbf{a}_1+\left(\tfrac{1}{2}-x_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}-z_7\right)\mathbf{a}_3 &&= x_7a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-z_7\right)a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{93} &= z_7\,\mathbf{a}_1+x_7\,\mathbf{a}_2+x_7\,\mathbf{a}_3 &&= z_7a\,\hat{\mathbf{x}}+x_7a\,\hat{\mathbf{y}}+x_7a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{94} &= z_7\,\mathbf{a}_1+\left(\tfrac{1}{2}-x_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}-x_7\right)\mathbf{a}_3 &&= z_7a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{95} &= \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}-x_7\right)\mathbf{a}_2+x_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-z_7\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{y}}+x_7a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{96} &= \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_1+x_7\,\mathbf{a}_2+\left(\tfrac{1}{2}-x_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-z_7\right)a\,\hat{\mathbf{x}}+x_7a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{97} &= x_7\,\mathbf{a}_1+z_7\,\mathbf{a}_2+x_7\,\mathbf{a}_3 &&= x_7a\,\hat{\mathbf{x}}+z_7a\,\hat{\mathbf{y}}+x_7a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{98} &= \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_1+z_7\,\mathbf{a}_2+\left(\tfrac{1}{2}-x_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{x}}+z_7a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{99} &= x_7\,\mathbf{a}_1+\left(\tfrac{1}{2}-z_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}-x_7\right)\mathbf{a}_3 &&= x_7a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-z_7\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{100} &= \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}-z_7\right)\mathbf{a}_2+x_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}-z_7\right)a\,\hat{\mathbf{y}}+x_7a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{101} &= \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_7\right)\mathbf{a}_2-z_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{y}}-z_7a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{102} &= -x_7\,\mathbf{a}_1-x_7\,\mathbf{a}_2-z_7\,\mathbf{a}_3 &&= -x_7a\,\hat{\mathbf{x}}-x_7a\,\hat{\mathbf{y}}-z_7a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{103} &= \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1-x_7\,\mathbf{a}_2+\left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{x}}-x_7a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{104} &= -x_7\,\mathbf{a}_1+\left(\tfrac{1}{2}+x_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 &&= -x_7a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{105} &= \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1+\left(\tfrac{1}{2}+z_7\right)\mathbf{a}_2-x_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{y}}-x_7a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{106} &= -x_7\,\mathbf{a}_1+\left(\tfrac{1}{2}+z_7\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_7\right)\mathbf{a}_3 &&= -x_7a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{z}} && (24k) && \text{W}
\end{aligned}
$$

$\mathbf{B}_{107}$	=	$-x_7\,\mathbf{a}_1 - z_7\,\mathbf{a}_2 - x_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} - z_7 a\,\hat{\mathbf{y}} - x_7 a\,\hat{\mathbf{z}}$	(24*k*)	W
$\mathbf{B}_{108}$	=	$\left(\frac{1}{2}+x_7\right)\mathbf{a}_1 - z_7\,\mathbf{a}_2 + \left(\frac{1}{2}+x_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{x}} - z_7 a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{z}}$	(24*k*)	W
$\mathbf{B}_{109}$	=	$\left(\frac{1}{2}+z_7\right)\mathbf{a}_1 + \left(\frac{1}{2}+x_7\right)\mathbf{a}_2 - x_7\,\mathbf{a}_3$	=	$\left(\frac{1}{2}+z_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{y}} - x_7 a\,\hat{\mathbf{z}}$	(24*k*)	W
$\mathbf{B}_{110}$	=	$\left(\frac{1}{2}+z_7\right)\mathbf{a}_1 - x_7\,\mathbf{a}_2 + \left(\frac{1}{2}+x_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{2}+z_7\right)a\,\hat{\mathbf{x}} - x_7 a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{z}}$	(24*k*)	W
$\mathbf{B}_{111}$	=	$-z_7\,\mathbf{a}_1 + \left(\frac{1}{2}+x_7\right)\mathbf{a}_2 + \left(\frac{1}{2}+x_7\right)\mathbf{a}_3$	=	$-z_7 a\,\hat{\mathbf{x}} + \left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2}+x_7\right)a\,\hat{\mathbf{z}}$	(24*k*)	W
$\mathbf{B}_{112}$	=	$-z_7\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 - x_7\,\mathbf{a}_3$	=	$-z_7 a\,\hat{\mathbf{x}} - x_7 a\,\hat{\mathbf{y}} - x_7 a\,\hat{\mathbf{z}}$	(24*k*)	W

References:

- L. Marosi, E. E. Platero, J. Cifre, and C. O. Areán, *Thermal dehydration of $H_{3+x}PV_xM_{12-x}O_{40}{\cdot}yH_2O$ Keggin type heteropolyacids; formation, thermal stability and structure of the anhydrous acids $H_3PM_{12}O_{40}$, of the corresponding anhydrides $PM_{12}O_{38.5}$ and of a novel trihydrate $H_3PW_{12}O_{40}{\cdot}3H_2O$*, J. Mater. Chem. **10**, 1949–1955 (2000), doi:10.1039/b001476l.

Found in:

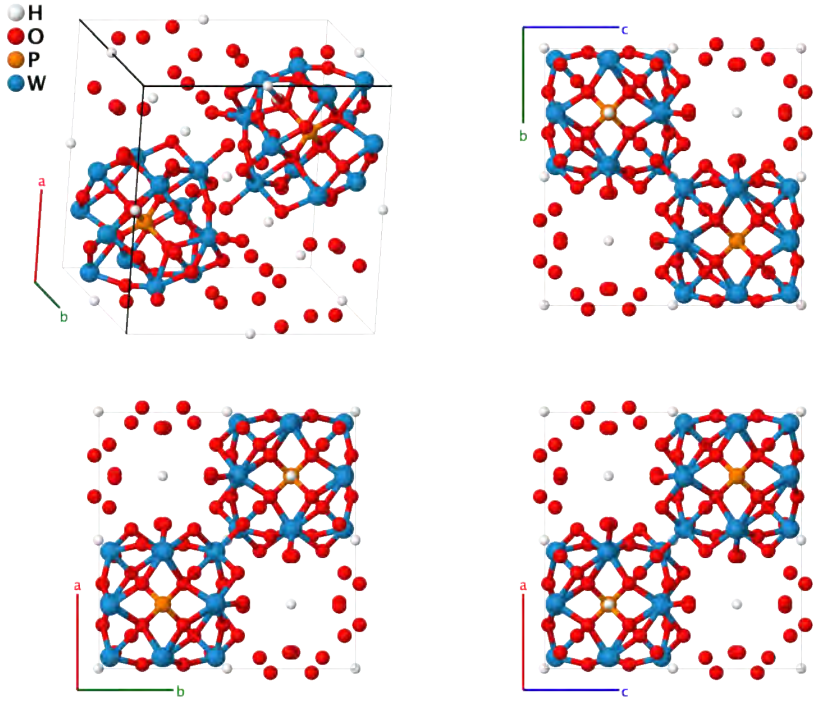
- P. Villars (Chief Editor), *$H_3PW_{12}O_{40}{\cdot}3H_2O$ Crystal Structure*, http://materials.springer.com/isp/crystallographic/docs/sd_1408693 (2016). In: Inorganic Solid Phases, SpringerMaterials (online database).

Geometry files:

- CIF: pp. 1808
- POSCAR: pp. 1809

12-phosphotungstic acid [$H_3PW_{12}O_{40}{\cdot}5H_2O$ ($H4_{16}$)] Structure: A5B40CD12_cP116_224_cd_e3k_a_k

http://aflow.org/prototype-encyclopedia/A5B40CD12_cP116_224_cd_e3k_a_k



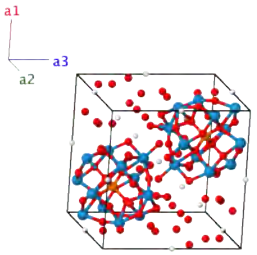
Prototype	:	$(H_{2.6}O)_5O_{40}PW_{12}$
AFLOW prototype label	:	A5B40CD12_cP116_224_cd_e3k_a_k
***Strukturbericht* designation**	:	$H4_{16}$
Pearson symbol	:	cP116
Space group number	:	224
Space group symbol	:	$Pn\bar{3}m$
AFLOW prototype command	:	`aflow --proto=A5B40CD12_cP116_224_cd_e3k_a_k --params=`$a, x_4, x_5, z_5, x_6, z_6, x_7, z_7, x_8, z_8$

- This structure is a bit problematic. (Brown, 1977) offers evidence that the true structure of the hydrated phosphotungstic acid is $H_3PW_{12}O_{40}{\cdot}6H_2O$, rather than $5H_2O$ as deduced by (Keggin, 1934). The presence of the extra water molecule does change the structure somewhat, and Keggin does not give the positions of the hydrogen atoms not attached to water molecules. Ordinarily we would mark this structure obsolete, but it is possible to remove water molecules from this structure and still have a molecule that is recognizably related to the acid, as in $H_3PW_{12}O_{40}{\cdot}3H_2O$. In addition, the lattice constant reported by Keggin is 12.141 Å compared to Brown's 12.506 Å, a change consistent with the loss of a water molecule. Given this, we will not deprecate this five-water molecule structure.
- (Gottfried, 1937) does not give the positions of the water molecules, and they reverse the x and z coordinates for O-IV and W.
- This structure is a partially dehydrated form of $H_3PW_{12}O_{40}{\cdot}29H_2O$ ($H4_{21}$). Further dehydration produces the $H_3PW_{12}O_{40}{\cdot}3H_2O$ structure.

- The positions of the three hydrogen atoms are unknown. If we follow (Marosi, 2000), then the 'free' hydrogens are likely to be bound to some of the water molecules, giving the unusual stoichiometry in the prototype.

Simple Cubic primitive vectors:

$$\mathbf{a}_1 = a\,\hat{\mathbf{x}}$$
$$\mathbf{a}_2 = a\,\hat{\mathbf{y}}$$
$$\mathbf{a}_3 = a\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(2*a*)	P
$\mathbf{B}_2$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(2*a*)	P
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(4*c*)	H_2O I
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{z}}$	(4*c*)	H_2O I
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{y}}$	(4*c*)	H_2O I
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$\frac{1}{2}a\,\hat{\mathbf{x}}$	(4*c*)	H_2O I
$\mathbf{B}_7$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(6*d*)	H_2O II
$\mathbf{B}_8$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(6*d*)	H_2O II
$\mathbf{B}_9$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(6*d*)	H_2O II
$\mathbf{B}_{10}$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(6*d*)	H_2O II
$\mathbf{B}_{11}$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(6*d*)	H_2O II
$\mathbf{B}_{12}$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(6*d*)	H_2O II
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{15}$	=	$\left(\frac{1}{2} - x_4\right)\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{16}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_4\right)\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_4\right)a\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{17}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_4\right)\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{y}} - x_4a\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{18}$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} - x_4a\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{19}$	=	$\left(\frac{1}{2} + x_4\right)\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \left(\frac{1}{2} + x_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{20}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + x_4\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_4\right)\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_4\right)a\,\hat{\mathbf{z}}$	(8*e*)	O I
$\mathbf{B}_{21}$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}} + x_5a\,\hat{\mathbf{y}} + z_5a\,\hat{\mathbf{z}}$	(24*k*)	O II
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_5\right)\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{y}} + z_5a\,\hat{\mathbf{z}}$	(24*k*)	O II
$\mathbf{B}_{23}$	=	$\left(\frac{1}{2} - x_5\right)\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{x}} + x_5a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)a\,\hat{\mathbf{z}}$	(24*k*)	O II
$\mathbf{B}_{24}$	=	$x_5\,\mathbf{a}_1 + \left(\frac{1}{2} - x_5\right)\mathbf{a}_2 + \left(\frac{1}{2} - z_5\right)\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - z_5\right)a\,\hat{\mathbf{z}}$	(24*k*)	O II
$\mathbf{B}_{25}$	=	$z_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$z_5a\,\hat{\mathbf{x}} + x_5a\,\hat{\mathbf{y}} + x_5a\,\hat{\mathbf{z}}$	(24*k*)	O II

$$
\begin{aligned}
\mathbf{B}_{26} &= z_5\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_5\right)\mathbf{a}_3 &&= z_5 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_5\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_5\right) a\,\hat{\mathbf{z}} && (24k) && \text{O II}\\
\mathbf{B}_{27} &= \left(\tfrac{1}{2} - z_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_5\right)\mathbf{a}_2 + x_5\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} - z_5\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_5\right) a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}} && (24k) && \text{O II}\\
\mathbf{B}_{28} &= \left(\tfrac{1}{2} - z_5\right)\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_5\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - z_5\right) a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_5\right) a\,\hat{\mathbf{z}} && (24k) && \text{O II}\\
\mathbf{B}_{29} &= x_5\,\mathbf{a}_1 + z_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3 &&= x_5 a\,\hat{\mathbf{x}} + z_5 a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}} && (24k) && \text{O II}\\
\mathbf{B}_{30} &= \left(\tfrac{1}{2} - x_5\right)\mathbf{a}_1 + z_5\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_5\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_5\right) a\,\hat{\mathbf{x}} + z_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_5\right) a\,\hat{\mathbf{z}} && (24k) && \text{O II}\\
\mathbf{B}_{31} &= x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2} - z_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_5\right)\mathbf{a}_3 &&= x_5 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - z_5\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_5\right) a\,\hat{\mathbf{z}} && (24k) && \text{O II}\\
\mathbf{B}_{32} &= \left(\tfrac{1}{2} - x_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - z_5\right)\mathbf{a}_2 + x_5\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_5\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - z_5\right) a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}} && (24k) && \text{O II}\\
\mathbf{B}_{33} &= \left(\tfrac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_5\right) a\,\hat{\mathbf{y}} - z_5 a\,\hat{\mathbf{z}} && (24k) && \text{O II}\\
\mathbf{B}_{34} &= -x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3 &&= -x_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} - z_5 a\,\hat{\mathbf{z}} && (24k) && \text{O II}\\
\mathbf{B}_{35} &= \left(\tfrac{1}{2} + x_5\right)\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_5\right) a\,\hat{\mathbf{z}} && (24k) && \text{O II}\\
\mathbf{B}_{36} &= -x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_3 &&= -x_5 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_5\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_5\right) a\,\hat{\mathbf{z}} && (24k) && \text{O II}\\
\mathbf{B}_{37} &= \left(\tfrac{1}{2} + x_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_2 - x_5\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + z_5\right) a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}} && (24k) && \text{O II}\\
\mathbf{B}_{38} &= -x_5\,\mathbf{a}_1 + \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_5\right)\mathbf{a}_3 &&= -x_5 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + z_5\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_5\right) a\,\hat{\mathbf{z}} && (24k) && \text{O II}\\
\mathbf{B}_{39} &= -x_5\,\mathbf{a}_1 - z_5\,\mathbf{a}_2 - x_5\,\mathbf{a}_3 &&= -x_5 a\,\hat{\mathbf{x}} - z_5 a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}} && (24k) && \text{O II}\\
\mathbf{B}_{40} &= \left(\tfrac{1}{2} + x_5\right)\mathbf{a}_1 - z_5\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_5\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_5\right) a\,\hat{\mathbf{x}} - z_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_5\right) a\,\hat{\mathbf{z}} && (24k) && \text{O II}\\
\mathbf{B}_{41} &= \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_5\right)\mathbf{a}_2 - x_5\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} + z_5\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_5\right) a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}} && (24k) && \text{O II}\\
\mathbf{B}_{42} &= \left(\tfrac{1}{2} + z_5\right)\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_5\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} + z_5\right) a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_5\right) a\,\hat{\mathbf{z}} && (24k) && \text{O II}\\
\mathbf{B}_{43} &= -z_5\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_5\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_5\right)\mathbf{a}_3 &&= -z_5 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_5\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_5\right) a\,\hat{\mathbf{z}} && (24k) && \text{O II}\\
\mathbf{B}_{44} &= -z_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - x_5\,\mathbf{a}_3 &&= -z_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}} && (24k) && \text{O II}\\
\mathbf{B}_{45} &= x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + z_6\,\mathbf{a}_3 &&= x_6 a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} + z_6 a\,\hat{\mathbf{z}} && (24k) && \text{O III}\\
\mathbf{B}_{46} &= \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_2 + z_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_6\right) a\,\hat{\mathbf{y}} + z_6 a\,\hat{\mathbf{z}} && (24k) && \text{O III}\\
\mathbf{B}_{47} &= \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_1 + x_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_6\right) a\,\hat{\mathbf{z}} && (24k) && \text{O III}\\
\mathbf{B}_{48} &= x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_3 &&= x_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_6\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_6\right) a\,\hat{\mathbf{z}} && (24k) && \text{O III}\\
\mathbf{B}_{49} &= z_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + x_6\,\mathbf{a}_3 &&= z_6 a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} + x_6 a\,\hat{\mathbf{z}} && (24k) && \text{O III}\\
\mathbf{B}_{50} &= z_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_3 &&= z_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_6\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_6\right) a\,\hat{\mathbf{z}} && (24k) && \text{O III}\\
\mathbf{B}_{51} &= \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_2 + x_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} - z_6\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_6\right) a\,\hat{\mathbf{y}} + x_6 a\,\hat{\mathbf{z}} && (24k) && \text{O III}\\
\mathbf{B}_{52} &= \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_1 + x_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - z_6\right) a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_6\right) a\,\hat{\mathbf{z}} && (24k) && \text{O III}\\
\mathbf{B}_{53} &= x_6\,\mathbf{a}_1 + z_6\,\mathbf{a}_2 + x_6\,\mathbf{a}_3 &&= x_6 a\,\hat{\mathbf{x}} + z_6 a\,\hat{\mathbf{y}} + x_6 a\,\hat{\mathbf{z}} && (24k) && \text{O III}\\
\mathbf{B}_{54} &= \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_1 + z_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} + z_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_6\right) a\,\hat{\mathbf{z}} && (24k) && \text{O III}\\
\mathbf{B}_{55} &= x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_3 &&= x_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - z_6\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_6\right) a\,\hat{\mathbf{z}} && (24k) && \text{O III}\\
\mathbf{B}_{56} &= \left(\tfrac{1}{2} - x_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - z_6\right)\mathbf{a}_2 + x_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} - x_6\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - z_6\right) a\,\hat{\mathbf{y}} + x_6 a\,\hat{\mathbf{z}} && (24k) && \text{O III}\\
\mathbf{B}_{57} &= \left(\tfrac{1}{2} + x_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_6\right)\mathbf{a}_2 - z_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_6\right) a\,\hat{\mathbf{y}} - z_6 a\,\hat{\mathbf{z}} && (24k) && \text{O III}\\
\mathbf{B}_{58} &= -x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3 &&= -x_6 a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} - z_6 a\,\hat{\mathbf{z}} && (24k) && \text{O III}\\
\mathbf{B}_{59} &= \left(\tfrac{1}{2} + x_6\right)\mathbf{a}_1 - x_6\,\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right) a\,\hat{\mathbf{z}} && (24k) && \text{O III}\\
\mathbf{B}_{60} &= -x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_3 &&= -x_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_6\right) a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_6\right) a\,\hat{\mathbf{z}} && (24k) && \text{O III}\\
\mathbf{B}_{61} &= \left(\tfrac{1}{2} + x_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + z_6\right)\mathbf{a}_2 - x_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2} + x_6\right) a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + z_6\right) a\,\hat{\mathbf{y}} - x_6 a\,\hat{\mathbf{z}} && (24k) && \text{O III}
\end{aligned}
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\begin{aligned}
\mathbf{B}_{62} &= -x_6\,\mathbf{a}_1 + \left(\tfrac{1}{2}+z_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_3 &&= -x_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+z_6\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{z}} && (24k) && \text{O III}\\
\mathbf{B}_{63} &= -x_6\,\mathbf{a}_1 - z_6\,\mathbf{a}_2 - x_6\,\mathbf{a}_3 &&= -x_6 a\,\hat{\mathbf{x}} - z_6 a\,\hat{\mathbf{y}} - x_6 a\,\hat{\mathbf{z}} && (24k) && \text{O III}\\
\mathbf{B}_{64} &= \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_1 - z_6\,\mathbf{a}_2 + \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{x}} - z_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{z}} && (24k) && \text{O III}\\
\mathbf{B}_{65} &= \left(\tfrac{1}{2}+z_6\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_2 - x_6\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+z_6\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{y}} - x_6 a\,\hat{\mathbf{z}} && (24k) && \text{O III}\\
\mathbf{B}_{66} &= \left(\tfrac{1}{2}+z_6\right)\mathbf{a}_1 - x_6\,\mathbf{a}_2 + \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+z_6\right)a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{z}} && (24k) && \text{O III}\\
\mathbf{B}_{67} &= -z_6\,\mathbf{a}_1 + \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+x_6\right)\mathbf{a}_3 &&= -z_6 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_6\right)a\,\hat{\mathbf{z}} && (24k) && \text{O III}\\
\mathbf{B}_{68} &= -z_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 - x_6\,\mathbf{a}_3 &&= -z_6 a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} - x_6 a\,\hat{\mathbf{z}} && (24k) && \text{O III}\\
\mathbf{B}_{69} &= x_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 &&= x_7 a\,\hat{\mathbf{x}} + x_7 a\,\hat{\mathbf{y}} + z_7 a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{70} &= \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_2 + z_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{y}} + z_7 a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{71} &= \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_1 + x_7\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{x}} + x_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_7\right)a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{72} &= x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_3 &&= x_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_7\right)a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{73} &= z_7\,\mathbf{a}_1 + x_7\,\mathbf{a}_2 + x_7\,\mathbf{a}_3 &&= z_7 a\,\hat{\mathbf{x}} + x_7 a\,\hat{\mathbf{y}} + x_7 a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{74} &= z_7\,\mathbf{a}_1 + \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_3 &&= z_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{75} &= \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_2 + x_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-z_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{y}} + x_7 a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{76} &= \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_1 + x_7\,\mathbf{a}_2 + \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-z_7\right)a\,\hat{\mathbf{x}} + x_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{77} &= x_7\,\mathbf{a}_1 + z_7\,\mathbf{a}_2 + x_7\,\mathbf{a}_3 &&= x_7 a\,\hat{\mathbf{x}} + z_7 a\,\hat{\mathbf{y}} + x_7 a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{78} &= \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_1 + z_7\,\mathbf{a}_2 + \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{x}} + z_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{79} &= x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_3 &&= x_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-z_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{80} &= \left(\tfrac{1}{2}-x_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-z_7\right)\mathbf{a}_2 + x_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-z_7\right)a\,\hat{\mathbf{y}} + x_7 a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{81} &= \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_2 - z_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{y}} - z_7 a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{82} &= -x_7\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3 &&= -x_7 a\,\hat{\mathbf{x}} - x_7 a\,\hat{\mathbf{y}} - z_7 a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{83} &= \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1 - x_7\,\mathbf{a}_2 + \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{x}} - x_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{84} &= -x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_3 &&= -x_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{85} &= \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_2 - x_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{y}} - x_7 a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{86} &= -x_7\,\mathbf{a}_1 + \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_3 &&= -x_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{87} &= -x_7\,\mathbf{a}_1 - z_7\,\mathbf{a}_2 - x_7\,\mathbf{a}_3 &&= -x_7 a\,\hat{\mathbf{x}} - z_7 a\,\hat{\mathbf{y}} - x_7 a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{88} &= \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_1 - z_7\,\mathbf{a}_2 + \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{x}} - z_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{89} &= \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_1 + \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_2 - x_7\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{y}} - x_7 a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{90} &= \left(\tfrac{1}{2}+z_7\right)\mathbf{a}_1 - x_7\,\mathbf{a}_2 + \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}+z_7\right)a\,\hat{\mathbf{x}} - x_7 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{91} &= -z_7\,\mathbf{a}_1 + \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+x_7\right)\mathbf{a}_3 &&= -z_7 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}+x_7\right)a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{92} &= -z_7\,\mathbf{a}_1 - x_7\,\mathbf{a}_2 - x_7\,\mathbf{a}_3 &&= -z_7 a\,\hat{\mathbf{x}} - x_7 a\,\hat{\mathbf{y}} - x_7 a\,\hat{\mathbf{z}} && (24k) && \text{O IV}\\
\mathbf{B}_{93} &= x_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3 &&= x_8 a\,\hat{\mathbf{x}} + x_8 a\,\hat{\mathbf{y}} + z_8 a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{94} &= \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_2 + z_8\,\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{y}} + z_8 a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{95} &= \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_1 + x_8\,\mathbf{a}_2 + \left(\tfrac{1}{2}-z_8\right)\mathbf{a}_3 &&= \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{x}} + x_8 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_8\right)a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{96} &= x_8\,\mathbf{a}_1 + \left(\tfrac{1}{2}-x_8\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-z_8\right)\mathbf{a}_3 &&= x_8 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_8\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-z_8\right)a\,\hat{\mathbf{z}} && (24k) && \text{W}\\
\mathbf{B}_{97} &= z_8\,\mathbf{a}_1 + x_8\,\mathbf{a}_2 + x_8\,\mathbf{a}_3 &&= z_8 a\,\hat{\mathbf{x}} + x_8 a\,\hat{\mathbf{y}} + x_8 a\,\hat{\mathbf{z}} && (24k) && \text{W}
\end{aligned}
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\begin{array}{lcccccc}
\mathbf{B}_{98} & = & z_8\,\mathbf{a}_1 + \left(\frac{1}{2} - x_8\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_8\right)\mathbf{a}_3 & = & z_8 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_8\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_8\right) a\,\hat{\mathbf{z}} & (24k) & \mathrm{W} \\
\mathbf{B}_{99} & = & \left(\frac{1}{2} - z_8\right)\mathbf{a}_1 + \left(\frac{1}{2} - x_8\right)\mathbf{a}_2 + x_8\,\mathbf{a}_3 & = & \left(\frac{1}{2} - z_8\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - x_8\right) a\,\hat{\mathbf{y}} + x_8 a\,\hat{\mathbf{z}} & (24k) & \mathrm{W} \\
\mathbf{B}_{100} & = & \left(\frac{1}{2} - z_8\right)\mathbf{a}_1 + x_8\,\mathbf{a}_2 + \left(\frac{1}{2} - x_8\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - z_8\right) a\,\hat{\mathbf{x}} + x_8 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_8\right) a\,\hat{\mathbf{z}} & (24k) & \mathrm{W} \\
\mathbf{B}_{101} & = & x_8\,\mathbf{a}_1 + z_8\,\mathbf{a}_2 + x_8\,\mathbf{a}_3 & = & x_8 a\,\hat{\mathbf{x}} + z_8 a\,\hat{\mathbf{y}} + x_8 a\,\hat{\mathbf{z}} & (24k) & \mathrm{W} \\
\mathbf{B}_{102} & = & \left(\frac{1}{2} - x_8\right)\mathbf{a}_1 + z_8\,\mathbf{a}_2 + \left(\frac{1}{2} - x_8\right)\mathbf{a}_3 & = & \left(\frac{1}{2} - x_8\right) a\,\hat{\mathbf{x}} + z_8 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_8\right) a\,\hat{\mathbf{z}} & (24k) & \mathrm{W} \\
\mathbf{B}_{103} & = & x_8\,\mathbf{a}_1 + \left(\frac{1}{2} - z_8\right)\mathbf{a}_2 + \left(\frac{1}{2} - x_8\right)\mathbf{a}_3 & = & x_8 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - z_8\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{2} - x_8\right) a\,\hat{\mathbf{z}} & (24k) & \mathrm{W} \\
\mathbf{B}_{104} & = & \left(\frac{1}{2} - x_8\right)\mathbf{a}_1 + \left(\frac{1}{2} - z_8\right)\mathbf{a}_2 + x_8\,\mathbf{a}_3 & = & \left(\frac{1}{2} - x_8\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} - z_8\right) a\,\hat{\mathbf{y}} + x_8 a\,\hat{\mathbf{z}} & (24k) & \mathrm{W} \\
\mathbf{B}_{105} & = & \left(\frac{1}{2} + x_8\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_8\right)\mathbf{a}_2 - z_8\,\mathbf{a}_3 & = & \left(\frac{1}{2} + x_8\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_8\right) a\,\hat{\mathbf{y}} - z_8 a\,\hat{\mathbf{z}} & (24k) & \mathrm{W} \\
\mathbf{B}_{106} & = & -x_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3 & = & -x_8 a\,\hat{\mathbf{x}} - x_8 a\,\hat{\mathbf{y}} - z_8 a\,\hat{\mathbf{z}} & (24k) & \mathrm{W} \\
\mathbf{B}_{107} & = & \left(\frac{1}{2} + x_8\right)\mathbf{a}_1 - x_8\,\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3 & = & \left(\frac{1}{2} + x_8\right) a\,\hat{\mathbf{x}} - x_8 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right) a\,\hat{\mathbf{z}} & (24k) & \mathrm{W} \\
\mathbf{B}_{108} & = & -x_8\,\mathbf{a}_1 + \left(\frac{1}{2} + x_8\right)\mathbf{a}_2 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_3 & = & -x_8 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_8\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + z_8\right) a\,\hat{\mathbf{z}} & (24k) & \mathrm{W} \\
\mathbf{B}_{109} & = & \left(\frac{1}{2} + x_8\right)\mathbf{a}_1 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_2 - x_8\,\mathbf{a}_3 & = & \left(\frac{1}{2} + x_8\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + z_8\right) a\,\hat{\mathbf{y}} - x_8 a\,\hat{\mathbf{z}} & (24k) & \mathrm{W} \\
\mathbf{B}_{110} & = & -x_8\,\mathbf{a}_1 + \left(\frac{1}{2} + z_8\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_8\right)\mathbf{a}_3 & = & -x_8 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + z_8\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_8\right) a\,\hat{\mathbf{z}} & (24k) & \mathrm{W} \\
\mathbf{B}_{111} & = & -x_8\,\mathbf{a}_1 - z_8\,\mathbf{a}_2 - x_8\,\mathbf{a}_3 & = & -x_8 a\,\hat{\mathbf{x}} - z_8 a\,\hat{\mathbf{y}} - x_8 a\,\hat{\mathbf{z}} & (24k) & \mathrm{W} \\
\mathbf{B}_{112} & = & \left(\frac{1}{2} + x_8\right)\mathbf{a}_1 - z_8\,\mathbf{a}_2 + \left(\frac{1}{2} + x_8\right)\mathbf{a}_3 & = & \left(\frac{1}{2} + x_8\right) a\,\hat{\mathbf{x}} - z_8 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_8\right) a\,\hat{\mathbf{z}} & (24k) & \mathrm{W} \\
\mathbf{B}_{113} & = & \left(\frac{1}{2} + z_8\right)\mathbf{a}_1 + \left(\frac{1}{2} + x_8\right)\mathbf{a}_2 - x_8\,\mathbf{a}_3 & = & \left(\frac{1}{2} + z_8\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_8\right) a\,\hat{\mathbf{y}} - x_8 a\,\hat{\mathbf{z}} & (24k) & \mathrm{W} \\
\mathbf{B}_{114} & = & \left(\frac{1}{2} + z_8\right)\mathbf{a}_1 - x_8\,\mathbf{a}_2 + \left(\frac{1}{2} + x_8\right)\mathbf{a}_3 & = & \left(\frac{1}{2} + z_8\right) a\,\hat{\mathbf{x}} - x_8 a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_8\right) a\,\hat{\mathbf{z}} & (24k) & \mathrm{W} \\
\mathbf{B}_{115} & = & -z_8\,\mathbf{a}_1 + \left(\frac{1}{2} + x_8\right)\mathbf{a}_2 + \left(\frac{1}{2} + x_8\right)\mathbf{a}_3 & = & -z_8 a\,\hat{\mathbf{x}} + \left(\frac{1}{2} + x_8\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{2} + x_8\right) a\,\hat{\mathbf{z}} & (24k) & \mathrm{W} \\
\mathbf{B}_{116} & = & -z_8\,\mathbf{a}_1 - x_8\,\mathbf{a}_2 - x_8\,\mathbf{a}_3 & = & -z_8 a\,\hat{\mathbf{x}} - x_8 a\,\hat{\mathbf{y}} - x_8 a\,\hat{\mathbf{z}} & (24k) & \mathrm{W}
\end{array}
$$

References:

- J. F. Keggin, *The structure and formula of 12-phosphotungstic acid*, Proc. Roy. Soc. Lond. A **144**, 75–100 (1934), doi:10.1098/rspa.1934.0035.
- C. Gottfried and F. Schossberger, eds., *Strukturbericht Band III 1933-1935* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).
- L. Marosi, E. E. Platero, J. Cifre, and C. O. Areán, *Thermal dehydration of $H_{3+x}PV_xM_{12-x}O_{40}{\cdot}yH_2O$ Keggin type heteropolyacids; formation, thermal stability and structure of the anhydrous acids $H_3PM_{12}O_{40}$, of the corresponding anhydrides $PM_{12}O_{38.5}$ and of a novel trihydrate $H_3PW_{12}O_{40}{\cdot}3H_2O$*, J. Mater. Chem. **10**, 1949–1955 (2000), doi:10.1039/b001476l.

Found in:

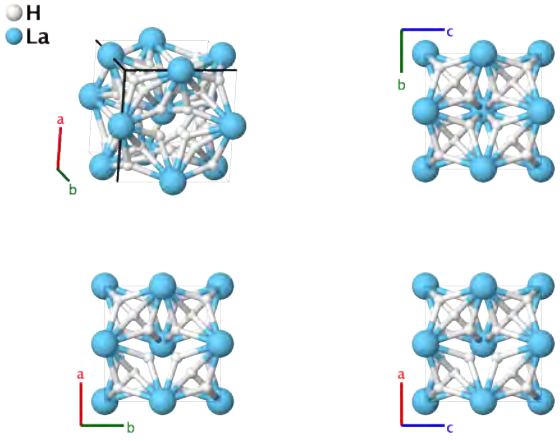
- G. M. Brown, M.-R. Noe-Spirlet, W. R. Busing, and H. A. Levy, *Dodecatungstophosphoric acid hexahydrate, $(H_5O_2^+)_3(PW_{12}O_{40}^{3-})$. The true structure of Keggin's 'pentahydrate' from single-crystal X-ray and neutron diffraction data*, Acta Crystallogr. Sect. B Struct. Sci. **33**, 1038–1046 (1977), doi:10.1107/S0567740877005330.

Geometry files:

- CIF: pp. 1810
- POSCAR: pp. 1810

LaH_{10} High-T_c Superconductor Structure: A10B_cF44_225_cf_b

http://aflow.org/prototype-encyclopedia/A10B_cF44_225_cf_b

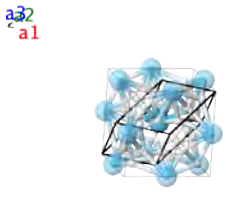


Prototype	:	$H_{10}La$
AFLOW prototype label	:	A10B_cF44_225_cf_b
***Strukturbericht* designation**	:	None
Pearson symbol	:	cF44
Space group number	:	225
Space group symbol	:	$Fm\bar{3}m$
AFLOW prototype command	:	`aflow --proto=A10B_cF44_225_cf_b --params=`a, x_3

- This structure was predicted by (Liu, 2017) as a possible near-room-temperature superconductor at pressures above 200 GPa. Its existence was confirmed by (Geballe, 2018).
- The lattice constant a = 4.78 Å is taken from (Liu, 2017) prediction at 300 GPa. (Geballe, 2018) found a = 5.09 Å at 172 GPa.
- (Liu, 2017) give a list of hydrogen positions which are not compatible with space group $Fm\bar{3}m$ #225. The positions we assigned are consistent with the work of (Geballe, 2018).

Face-centered Cubic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$$



Basis vectors:

Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type

$\mathbf{B}_1$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(4*b*)	La
$\mathbf{B}_2$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(8*c*)	H I
$\mathbf{B}_3$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(8*c*)	H I
$\mathbf{B}_4$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(32*f*)	H II
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 - 3x_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(32*f*)	H II
$\mathbf{B}_6$	=	$x_3\,\mathbf{a}_1 - 3x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(32*f*)	H II
$\mathbf{B}_7$	=	$-3x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(32*f*)	H II
$\mathbf{B}_8$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + 3x_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(32*f*)	H II
$\mathbf{B}_9$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(32*f*)	H II
$\mathbf{B}_{10}$	=	$-x_3\,\mathbf{a}_1 + 3x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(32*f*)	H II
$\mathbf{B}_{11}$	=	$3x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(32*f*)	H II

References:

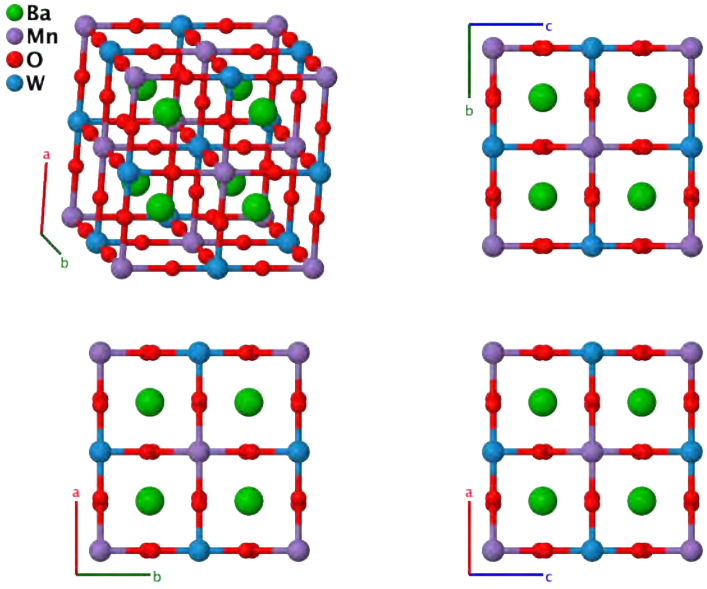
- H. Liu, I. I. Naumov, R. Hoffmann, N. W. Ashcroft, and R. J. Hemley, *Potential high-T_c superconducting lanthanum and yttrium hydrides at high pressure*, Proc. Natl. Acad. Sci. **114**, 6990–6995 (2017), doi:10.1073/pnas.1704505114.
- Z. M. Geballe, H. Liu, A. K. Mishra, M. Ahart, M. Somayazulu, Y. Meng, M. Baldini, and R. J. Hemley, *Synthesis and Stability of Lanthanum Superhydrides*, Angew. Chem. Int. Ed. **57**, 688–692 (2018), doi:10.1002/anie.201709970.

Geometry files:

- CIF: pp. 1811
- POSCAR: pp. 1812

Double Perovskite (Ba_2MnWO_6) Structure: A2BC6D_cF40_225_c_a_e_b

http://aflow.org/prototype-encyclopedia/A2BC6D_cF40_225_c_a_e_b



Prototype	:	Ba_2MnO_6W
AFLOW prototype label	:	A2BC6D_cF40_225_c_a_e_b
***Strukturbericht* designation**	:	None
Pearson symbol	:	cF40
Space group number	:	225
Space group symbol	:	$Fm\bar{3}m$
AFLOW prototype command	:	`aflow --proto=A2BC6D_cF40_225_c_a_e_b --params=`a, x_4

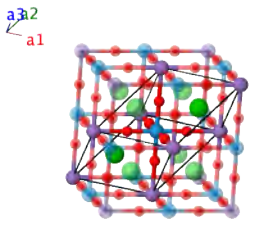
Other compounds with this structure

- Ba_2CaWO_6, Ba_2LiOsO_6, Ba_2NaOsO_6, Bi_2FeCrO_6, Ca_2MnReO_6, Cu_2TiSiO_6, Mn_2FeSbO_6, Sr_2FeMoO_6, Sr_2ReMoO_6, $(La_{0.5}Sc_{0.5})_2MnCoO_6$, and K_2NaAlF_6 (elpasolite)

- We use the data taken at room temperature.

Face-centered Cubic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(4*a*)	Mn
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(4*b*)	W
$\mathbf{B}_3$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(8*c*)	Ba
$\mathbf{B}_4$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(8*c*)	Ba
$\mathbf{B}_5$	=	$-x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}}$	(24*e*)	O
$\mathbf{B}_6$	=	$x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}}$	(24*e*)	O
$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{y}}$	(24*e*)	O
$\mathbf{B}_8$	=	$-x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{y}}$	(24*e*)	O
$\mathbf{B}_9$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{z}}$	(24*e*)	O
$\mathbf{B}_{10}$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{z}}$	(24*e*)	O

References:

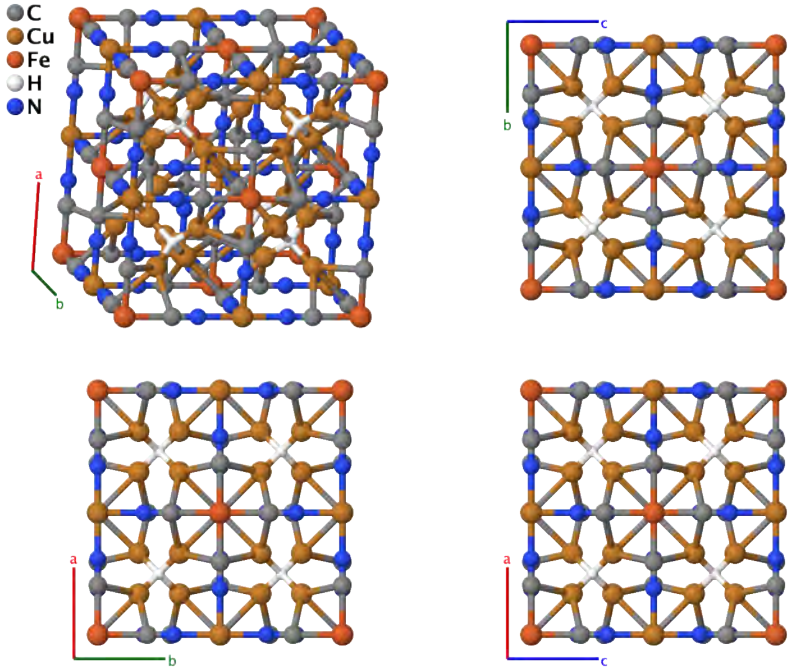
- A. K. Azad, S. A. Ivanov, S.-G. Eriksson, J. Eriksen, H. Rundlöf, R. Mathieu, and P. Svedlindh, *Synthesis, crystal structure, and magnetic characterization of the double perovskite Ba_2MnWO_6*, Mater. Res. Bull. **36**, 2215–2228 (2001), doi:10.1016/S0025-5408(01)00707-3.

Geometry files:

- CIF: pp. 1812
- POSCAR: pp. 1813

$Cu_3[Fe(CN)_6]_2{\cdot}xH_2O$ ($J2_5$, $x \approx 3$) Structure: A6B9CD2E6_cF96_225_e_bf_a_c_e

http://aflow.org/prototype-encyclopedia/A6B9CD2E6_cF96_225_e_bf_a_c_e



Prototype	:	$C_{12}Cu_3Fe_2(H_2O)_xN_{12}$
AFLOW prototype label	:	A6B9CD2E6_cF96_225_e_bf_a_c_e
***Strukturbericht* designation**	:	$J2_5$
Pearson symbol	:	cF96
Space group number	:	225
Space group symbol	:	$Fm\bar{3}m$
AFLOW prototype command	:	`aflow --proto=A6B9CD2E6_cF96_225_e_bf_a_c_e --params=`a, x_4, x_5, x_6

Other compounds with this structure

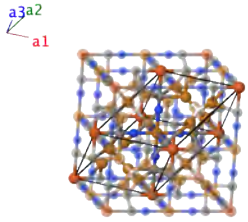
- $Cd_3[Co(CN)_6]_2$, $Co_3[Co(CN)_6]_2$, $Cu_3[Fe(CN)_6]_2$, $Fe_3[Fe(CN)_6]_2$, $Td_3[Fe(CN)_6]_2$, and $Zn_3[Fe(CN)_6]_2$

- These compounds form a class called "Prussian Blue Analogs," where Prussian Blue is $Fe_3[Fe(CN)_6]_2$.
- (van Bever, 1938) studied what he believed to be the hydrated form of this structure, with $x \approx 3$. In that case, the water molecules occupy the ($8c$) sites, but each site is only occupied 75% of the time. The water sites are surrounded by a tetrahedron of copper ($32f$) sites, but only 6.25% of these sites are occupied.
- (Weiser, 1942) studied the anhydrous form. They found that in this case the copper atoms that were on the ($32f$) sites replace the water molecules on the ($8c$) site, and this site is now fully occupied with copper.
- To convert from the hydrated to anhydrous structure, remove the copper atoms from the ($32f$) sites in the CIF or POSCAR file, and relabel the ($8c$) site as copper.

- For a picture of the resulting structure see (Jiao, 2017).
- The AFLOW label models the structure as if the sites were fully occupied.

Face-centered Cubic primitive vectors:

$$\mathbf{a}_1 = \tfrac{1}{2}\,a\,\hat{\mathbf{y}} + \tfrac{1}{2}\,a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{y}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(4a)$	Fe
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	$(4b)$	Cu I
$\mathbf{B}_3$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	$(8c)$	H_2O
$\mathbf{B}_4$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	$(8c)$	H_2O
$\mathbf{B}_5$	=	$-x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}}$	$(24e)$	C
$\mathbf{B}_6$	=	$x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}}$	$(24e)$	C
$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{y}}$	$(24e)$	C
$\mathbf{B}_8$	=	$-x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{y}}$	$(24e)$	C
$\mathbf{B}_9$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{z}}$	$(24e)$	C
$\mathbf{B}_{10}$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{z}}$	$(24e)$	C
$\mathbf{B}_{11}$	=	$-x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}}$	$(24e)$	N
$\mathbf{B}_{12}$	=	$x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - x_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}}$	$(24e)$	N
$\mathbf{B}_{13}$	=	$x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{y}}$	$(24e)$	N
$\mathbf{B}_{14}$	=	$-x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 - x_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{y}}$	$(24e)$	N
$\mathbf{B}_{15}$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 - x_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{z}}$	$(24e)$	N
$\mathbf{B}_{16}$	=	$-x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{z}}$	$(24e)$	N
$\mathbf{B}_{17}$	=	$x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + x_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} + x_6 a\,\hat{\mathbf{z}}$	$(32f)$	Cu II
$\mathbf{B}_{18}$	=	$x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 - 3x_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} + x_6 a\,\hat{\mathbf{z}}$	$(32f)$	Cu II
$\mathbf{B}_{19}$	=	$x_6\,\mathbf{a}_1 - 3x_6\,\mathbf{a}_2 + x_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} - x_6 a\,\hat{\mathbf{z}}$	$(32f)$	Cu II
$\mathbf{B}_{20}$	=	$-3x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + x_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} - x_6 a\,\hat{\mathbf{z}}$	$(32f)$	Cu II
$\mathbf{B}_{21}$	=	$-x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 + 3x_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} - x_6 a\,\hat{\mathbf{z}}$	$(32f)$	Cu II
$\mathbf{B}_{22}$	=	$-x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 - x_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} - x_6 a\,\hat{\mathbf{z}}$	$(32f)$	Cu II
$\mathbf{B}_{23}$	=	$-x_6\,\mathbf{a}_1 + 3x_6\,\mathbf{a}_2 - x_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} + x_6 a\,\hat{\mathbf{z}}$	$(32f)$	Cu II
$\mathbf{B}_{24}$	=	$3x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 - x_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} + x_6 a\,\hat{\mathbf{z}}$	$(32f)$	Cu II

References:

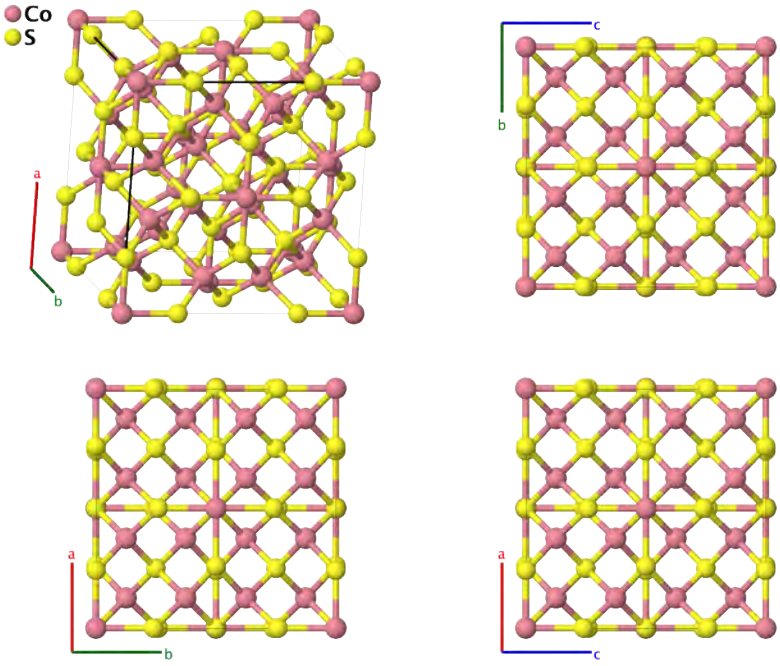
- A. K. van Bever, *The Crystal Structure of Some Ferricyanides with Bivalent Kations*, Rec. Trav. Chim. Pays-Bas **57**, 1259–1268 (1938), doi:10.1002/recl.19380571108.
- H. B. Weiser, W. O. Milligan, and J. B. Bates, *X-ray Diffraction Studies on Heavy-metal Iron-cyanides*, J. Phys. Chem. **46**, 99–111 (1942), doi:10.1021/j150415a013.
- S. Jiao, J. Tuo, H. Xie, Z. Cai, S. Wang, and J. Zhu, *The electrochemical performance of $Cu_3[Fe(CN)_6]_2$ as a cathode material for sodium-ion batteries*, Mater. Res. Bull. **86**, 194–200 (2017), doi:10.1016/j.materresbull.2016.10.019.

Geometry files:

- CIF: pp. 1813
- POSCAR: pp. 1814

Co_9S_8 ($D8_9$) Structure: A9B8_cF68_225_af_ce

http://aflow.org/prototype-encyclopedia/A9B8_cF68_225_af_ce



Prototype	:	Co_9S_8
AFLOW prototype label	:	A9B8_cF68_225_af_ce
***Strukturbericht* designation**	:	$D8_9$
Pearson symbol	:	cF68
Space group number	:	225
Space group symbol	:	$Fm\bar{3}m$
AFLOW prototype command	:	`aflow --proto=A9B8_cF68_225_af_ce --params=`a, x_3, x_4

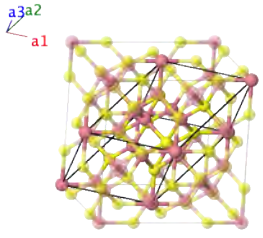
Other compounds with this structure

- $(Fe,Ni)_9S_8$ (pentlandite)

- (Geller, 1962) placed the first Co atom at the (4*b*) Wyckoff position. We have shifted this to the origin, the (4*a*) Wyckoff position.

Face-centered Cubic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(4*a*)	Co I
$\mathbf{B}_2$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(8*c*)	S I
$\mathbf{B}_3$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(8*c*)	S I
$\mathbf{B}_4$	=	$-x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}}$	(24*e*)	S II
$\mathbf{B}_5$	=	$x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}}$	(24*e*)	S II
$\mathbf{B}_6$	=	$x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{y}}$	(24*e*)	S II
$\mathbf{B}_7$	=	$-x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{y}}$	(24*e*)	S II
$\mathbf{B}_8$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{z}}$	(24*e*)	S II
$\mathbf{B}_9$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{z}}$	(24*e*)	S II
$\mathbf{B}_{10}$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	(32*f*)	Co II
$\mathbf{B}_{11}$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 - 3x_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	(32*f*)	Co II
$\mathbf{B}_{12}$	=	$x_4\,\mathbf{a}_1 - 3x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} - x_4a\,\hat{\mathbf{z}}$	(32*f*)	Co II
$\mathbf{B}_{13}$	=	$-3x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} - x_4a\,\hat{\mathbf{z}}$	(32*f*)	Co II
$\mathbf{B}_{14}$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + 3x_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} - x_4a\,\hat{\mathbf{z}}$	(32*f*)	Co II
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} - x_4a\,\hat{\mathbf{z}}$	(32*f*)	Co II
$\mathbf{B}_{16}$	=	$-x_4\,\mathbf{a}_1 + 3x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	(32*f*)	Co II
$\mathbf{B}_{17}$	=	$3x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	(32*f*)	Co II

References:

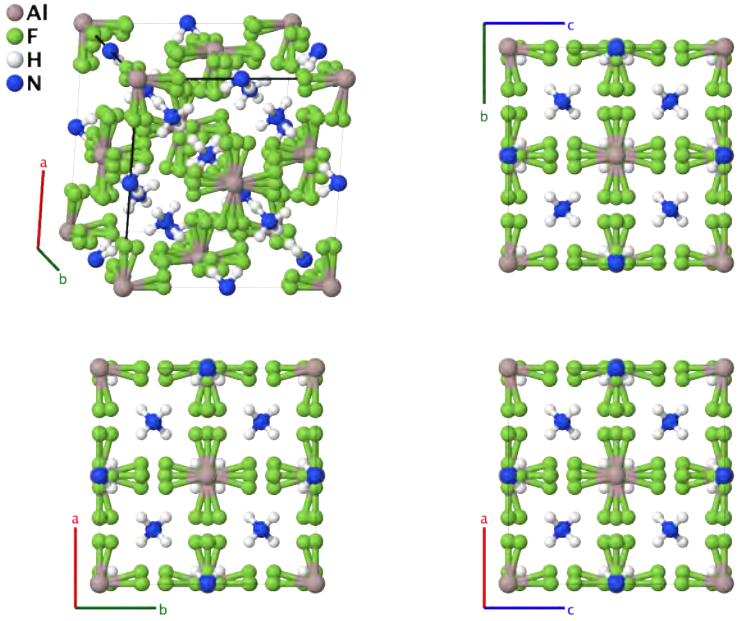
- S. Geller, *Refinement of the crystal structure of Co_9S_8*, Acta Cryst. **15**, 1195–1198 (1962), doi:10.1107/S0365110X62003187.

Geometry files:

- CIF: pp. 1815
- POSCAR: pp. 1816

$(NH_4)_3AlF_6$ ($J2_1$) Structure: AB30C16D3_cF200_225_a_ej_2f_bc

http://aflow.org/prototype-encyclopedia/AB30C16D3_cF200_225_a_ej_2f_bc



Prototype	:	$AlF_6H_{12}N_3$
AFLOW prototype label	:	AB30C16D3_cF200_225_a_ej_2f_bc
***Strukturbericht* designation**	:	$J2_1$
Pearson symbol	:	cF200
Space group number	:	225
Space group symbol	:	$Fm\bar{3}m$
AFLOW prototype command	:	`aflow --proto=AB30C16D3_cF200_225_a_ej_2f_bc --params=`$a, x_4, x_5, x_6, y_7, z_7$

Other compounds with this structure

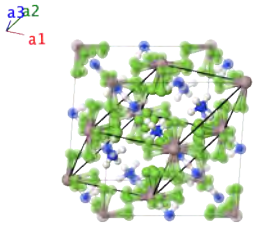
- $(NH_4)_3FeF_6$, $(NH_4)_3TiOF_5$, $(NH_4)_3Fe(NO_2)_6$, $K_3Ir(NO_2)_6$, $Cs_3Ir(NO_2)_6$, $Rb_3Ir(NO_2)_6$, and $Tl_3Ir(NO_2)_6$

- Early determinations of this structure placed all of the fluorine atoms on the (24*e*) site and were not able to determine the positions of the hydrogen atoms in the ammonium ion. This structure was designated $H71$ ($H7_1$) by (Ewald, 1931), renamed $I2_1$ by (Hermann, 1937) and finally given the label $J2_1$ by (Gottfried, 1937).

- The structure determined by (Udovenko, 2003) found that the fluorine atoms are split onto two sites, F-I, on Wyckoff position (24*e*) is 1/3 filled, and F-II, on (96*j*) is 1/6 filled. The positions are so close, however, that a reasonable approximation can be made by eliminating the (96*j*) site and fully occupying the (24*e*) site. The H-I (32*f*) site is fully occupied, while the H-II (32*f*) site is 50% occupied.

Face-centered Cubic primitive vectors:

$$\mathbf{a}_1 = \tfrac{1}{2}\,a\,\hat{\mathbf{y}} + \tfrac{1}{2}\,a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{y}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(4*a*)	Al
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(4*b*)	N I
$\mathbf{B}_3$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(8*c*)	N II
$\mathbf{B}_4$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(8*c*)	N II
$\mathbf{B}_5$	=	$-x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}}$	(24*e*)	F I
$\mathbf{B}_6$	=	$x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}}$	(24*e*)	F I
$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{y}}$	(24*e*)	F I
$\mathbf{B}_8$	=	$-x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{y}}$	(24*e*)	F I
$\mathbf{B}_9$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{z}}$	(24*e*)	F I
$\mathbf{B}_{10}$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{z}}$	(24*e*)	F I
$\mathbf{B}_{11}$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}}$	(32*f*)	H I
$\mathbf{B}_{12}$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 - 3x_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}}$	(32*f*)	H I
$\mathbf{B}_{13}$	=	$x_5\,\mathbf{a}_1 - 3x_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}}$	(32*f*)	H I
$\mathbf{B}_{14}$	=	$-3x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}}$	(32*f*)	H I
$\mathbf{B}_{15}$	=	$-x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + 3x_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}}$	(32*f*)	H I
$\mathbf{B}_{16}$	=	$-x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - x_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}}$	(32*f*)	H I
$\mathbf{B}_{17}$	=	$-x_5\,\mathbf{a}_1 + 3x_5\,\mathbf{a}_2 - x_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}}$	(32*f*)	H I
$\mathbf{B}_{18}$	=	$3x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - x_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}}$	(32*f*)	H I
$\mathbf{B}_{19}$	=	$x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + x_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} + x_6 a\,\hat{\mathbf{z}}$	(32*f*)	H II
$\mathbf{B}_{20}$	=	$x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 - 3x_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} + x_6 a\,\hat{\mathbf{z}}$	(32*f*)	H II
$\mathbf{B}_{21}$	=	$x_6\,\mathbf{a}_1 - 3x_6\,\mathbf{a}_2 + x_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} - x_6 a\,\hat{\mathbf{z}}$	(32*f*)	H II
$\mathbf{B}_{22}$	=	$-3x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + x_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} - x_6 a\,\hat{\mathbf{z}}$	(32*f*)	H II
$\mathbf{B}_{23}$	=	$-x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 + 3x_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} - x_6 a\,\hat{\mathbf{z}}$	(32*f*)	H II
$\mathbf{B}_{24}$	=	$-x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 - x_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} - x_6 a\,\hat{\mathbf{z}}$	(32*f*)	H II
$\mathbf{B}_{25}$	=	$-x_6\,\mathbf{a}_1 + 3x_6\,\mathbf{a}_2 - x_6\,\mathbf{a}_3$	=	$x_6 a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} + x_6 a\,\hat{\mathbf{z}}$	(32*f*)	H II
$\mathbf{B}_{26}$	=	$3x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 - x_6\,\mathbf{a}_3$	=	$-x_6 a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} + x_6 a\,\hat{\mathbf{z}}$	(32*f*)	H II
$\mathbf{B}_{27}$	=	$(y_7 + z_7)\,\mathbf{a}_1 + (-y_7 + z_7)\,\mathbf{a}_2 + (y_7 - z_7)\,\mathbf{a}_3$	=	$y_7 a\,\hat{\mathbf{y}} + z_7 a\,\hat{\mathbf{z}}$	(96*j*)	F II

$\mathbf{B}_{28}$	=	$(-y_7 + z_7)\,\mathbf{a}_1 + (y_7 + z_7)\,\mathbf{a}_2 + (-y_7 - z_7)\,\mathbf{a}_3$	=	$-y_7 a\,\hat{\mathbf{y}} + z_7 a\,\hat{\mathbf{z}}$	(96*j*)	F II
$\mathbf{B}_{29}$	=	$(y_7 - z_7)\,\mathbf{a}_1 + (-y_7 - z_7)\,\mathbf{a}_2 + (y_7 + z_7)\,\mathbf{a}_3$	=	$y_7 a\,\hat{\mathbf{y}} - z_7 a\,\hat{\mathbf{z}}$	(96*j*)	F II
$\mathbf{B}_{30}$	=	$(-y_7 - z_7)\,\mathbf{a}_1 + (y_7 - z_7)\,\mathbf{a}_2 + (-y_7 + z_7)\,\mathbf{a}_3$	=	$-y_7 a\,\hat{\mathbf{y}} - z_7 a\,\hat{\mathbf{z}}$	(96*j*)	F II
$\mathbf{B}_{31}$	=	$(y_7 - z_7)\,\mathbf{a}_1 + (y_7 + z_7)\,\mathbf{a}_2 + (-y_7 + z_7)\,\mathbf{a}_3$	=	$z_7 a\,\hat{\mathbf{x}} + y_7 a\,\hat{\mathbf{z}}$	(96*j*)	F II
$\mathbf{B}_{32}$	=	$(-y_7 - z_7)\,\mathbf{a}_1 + (-y_7 + z_7)\,\mathbf{a}_2 + (y_7 + z_7)\,\mathbf{a}_3$	=	$z_7 a\,\hat{\mathbf{x}} - y_7 a\,\hat{\mathbf{z}}$	(96*j*)	F II
$\mathbf{B}_{33}$	=	$(y_7 + z_7)\,\mathbf{a}_1 + (y_7 - z_7)\,\mathbf{a}_2 + (-y_7 - z_7)\,\mathbf{a}_3$	=	$-z_7 a\,\hat{\mathbf{x}} + y_7 a\,\hat{\mathbf{z}}$	(96*j*)	F II
$\mathbf{B}_{34}$	=	$(-y_7 + z_7)\,\mathbf{a}_1 + (-y_7 - z_7)\,\mathbf{a}_2 + (y_7 - z_7)\,\mathbf{a}_3$	=	$-z_7 a\,\hat{\mathbf{x}} - y_7 a\,\hat{\mathbf{z}}$	(96*j*)	F II
$\mathbf{B}_{35}$	=	$(-y_7 + z_7)\,\mathbf{a}_1 + (y_7 - z_7)\,\mathbf{a}_2 + (y_7 + z_7)\,\mathbf{a}_3$	=	$y_7 a\,\hat{\mathbf{x}} + z_7 a\,\hat{\mathbf{y}}$	(96*j*)	F II
$\mathbf{B}_{36}$	=	$(y_7 + z_7)\,\mathbf{a}_1 + (-y_7 - z_7)\,\mathbf{a}_2 + (-y_7 + z_7)\,\mathbf{a}_3$	=	$-y_7 a\,\hat{\mathbf{x}} + z_7 a\,\hat{\mathbf{y}}$	(96*j*)	F II
$\mathbf{B}_{37}$	=	$(-y_7 - z_7)\,\mathbf{a}_1 + (y_7 + z_7)\,\mathbf{a}_2 + (y_7 - z_7)\,\mathbf{a}_3$	=	$y_7 a\,\hat{\mathbf{x}} - z_7 a\,\hat{\mathbf{y}}$	(96*j*)	F II
$\mathbf{B}_{38}$	=	$(y_7 - z_7)\,\mathbf{a}_1 + (-y_7 + z_7)\,\mathbf{a}_2 + (-y_7 - z_7)\,\mathbf{a}_3$	=	$-y_7 a\,\hat{\mathbf{x}} - z_7 a\,\hat{\mathbf{y}}$	(96*j*)	F II
$\mathbf{B}_{39}$	=	$(-y_7 - z_7)\,\mathbf{a}_1 + (y_7 - z_7)\,\mathbf{a}_2 + (y_7 + z_7)\,\mathbf{a}_3$	=	$y_7 a\,\hat{\mathbf{x}} - z_7 a\,\hat{\mathbf{z}}$	(96*j*)	F II
$\mathbf{B}_{40}$	=	$(y_7 - z_7)\,\mathbf{a}_1 + (-y_7 - z_7)\,\mathbf{a}_2 + (-y_7 + z_7)\,\mathbf{a}_3$	=	$-y_7 a\,\hat{\mathbf{x}} - z_7 a\,\hat{\mathbf{z}}$	(96*j*)	F II
$\mathbf{B}_{41}$	=	$(-y_7 + z_7)\,\mathbf{a}_1 + (y_7 + z_7)\,\mathbf{a}_2 + (y_7 - z_7)\,\mathbf{a}_3$	=	$y_7 a\,\hat{\mathbf{x}} + z_7 a\,\hat{\mathbf{z}}$	(96*j*)	F II
$\mathbf{B}_{42}$	=	$(y_7 + z_7)\,\mathbf{a}_1 + (-y_7 + z_7)\,\mathbf{a}_2 + (-y_7 - z_7)\,\mathbf{a}_3$	=	$-y_7 a\,\hat{\mathbf{x}} + z_7 a\,\hat{\mathbf{z}}$	(96*j*)	F II
$\mathbf{B}_{43}$	=	$(-y_7 + z_7)\,\mathbf{a}_1 + (-y_7 - z_7)\,\mathbf{a}_2 + (y_7 + z_7)\,\mathbf{a}_3$	=	$z_7 a\,\hat{\mathbf{y}} - y_7 a\,\hat{\mathbf{z}}$	(96*j*)	F II
$\mathbf{B}_{44}$	=	$(y_7 + z_7)\,\mathbf{a}_1 + (y_7 - z_7)\,\mathbf{a}_2 + (-y_7 + z_7)\,\mathbf{a}_3$	=	$z_7 a\,\hat{\mathbf{y}} + y_7 a\,\hat{\mathbf{z}}$	(96*j*)	F II
$\mathbf{B}_{45}$	=	$(-y_7 - z_7)\,\mathbf{a}_1 + (-y_7 + z_7)\,\mathbf{a}_2 + (y_7 - z_7)\,\mathbf{a}_3$	=	$-z_7 a\,\hat{\mathbf{y}} - y_7 a\,\hat{\mathbf{z}}$	(96*j*)	F II
$\mathbf{B}_{46}$	=	$(y_7 - z_7)\,\mathbf{a}_1 + (y_7 + z_7)\,\mathbf{a}_2 + (-y_7 - z_7)\,\mathbf{a}_3$	=	$-z_7 a\,\hat{\mathbf{y}} + y_7 a\,\hat{\mathbf{z}}$	(96*j*)	F II
$\mathbf{B}_{47}$	=	$(y_7 - z_7)\,\mathbf{a}_1 + (-y_7 + z_7)\,\mathbf{a}_2 + (y_7 + z_7)\,\mathbf{a}_3$	=	$z_7 a\,\hat{\mathbf{x}} + y_7 a\,\hat{\mathbf{y}}$	(96*j*)	F II
$\mathbf{B}_{48}$	=	$(-y_7 - z_7)\,\mathbf{a}_1 + (y_7 + z_7)\,\mathbf{a}_2 + (-y_7 + z_7)\,\mathbf{a}_3$	=	$z_7 a\,\hat{\mathbf{x}} - y_7 a\,\hat{\mathbf{y}}$	(96*j*)	F II
$\mathbf{B}_{49}$	=	$(y_7 + z_7)\,\mathbf{a}_1 + (-y_7 - z_7)\,\mathbf{a}_2 + (y_7 - z_7)\,\mathbf{a}_3$	=	$-z_7 a\,\hat{\mathbf{x}} + y_7 a\,\hat{\mathbf{y}}$	(96*j*)	F II
$\mathbf{B}_{50}$	=	$(-y_7 + z_7)\,\mathbf{a}_1 + (y_7 - z_7)\,\mathbf{a}_2 + (-y_7 - z_7)\,\mathbf{a}_3$	=	$-z_7 a\,\hat{\mathbf{x}} - y_7 a\,\hat{\mathbf{y}}$	(96*j*)	F II

References:

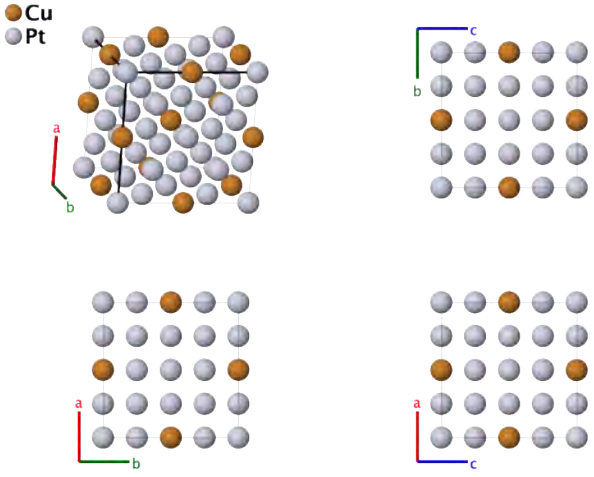
- A. A. Udovenko, N. M. Laptash, and I. G. Maslennikova, *Orientation disorder in ammonium elpasolites: Crystal structures of $(NH_4)_3AlF_6$, $(NH_4)_3TiOF_5$ and $(NH_4)_3FeF_6$*, J. Fluor. Chem. **124**, 5–15 (2003), doi:10.1016/S0022-1139(03)00166-0.
- P. P. Ewald and C. Hermann, eds., *Strukturbericht 1913-1928* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1931).
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).
- C. Gottfried and F. Schossberger, eds., *Strukturbericht Band III 1933-1935* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1816
- POSCAR: pp. 1817

$L1_a$ (disputed $CuPt_3$) Structure: AB7_cF32_225_b_ad

http://aflow.org/prototype-encyclopedia/AB7_cF32_225_b_ad

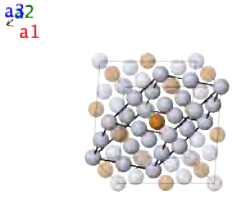


Prototype	:	$CuPt_3$
AFLOW prototype label	:	AB7_cF32_225_b_ad
***Strukturbericht* designation**	:	$L1_a$
Pearson symbol	:	cF32
Space group number	:	225
Space group symbol	:	$Fm\bar{3}m$
AFLOW prototype command	:	`aflow --proto=AB7_cF32_225_b_ad --params=`a

- According to (Tang, 1951), the (24*d*) sites have the composition $Pt_{0.8}Cu_{0.2}$ in stoichiometric $CuPt_3$. Here we use "Pt" to specify the atoms on this site.
- (Tang, 1951) states that the crystal structure of $CuPt_3$ must be cubic, but (Mshumi, 2014) argue that it is orthorhombic, and in fact the $L1_3$ structure.
- (Smithells, 1955) gave this structure the $L1_a$ designation as part of his extension of the original *Strukturbericht* labels. He does note that an alternative orthorhombic structure had been proposed.
- (Smithells, 1955) assigns this structure to space group $F432$ #209, but the positions given by (Tang, 1951) are also consistent with $Fm\bar{3}m$ #225, so we assign this structure to the higher symmetry space group.
- (Tang, 1951) does not give the lattice constant, so we use the value estimated by (Smithells, 1955).
- The Wyckoff positions are identical to those of the Ca_7Ge structure.

Face-centered Cubic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{y}} + \frac{1}{2}\,a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{1}{2}\,a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{1}{2}\,a\,\hat{\mathbf{y}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(4*a*)	Pt I
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(4*b*)	Cu
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$\frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(24*d*)	Pt II
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(24*d*)	Pt II
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(24*d*)	Pt II
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(24*d*)	Pt II
$\mathbf{B}_7$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}}$	(24*d*)	Pt II
$\mathbf{B}_8$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(24*d*)	Pt II

References:

- Y.-C. Tang, *A cubic structure for the phase Pt_3Cu*, Acta Cryst. **4**, 377–378 (1951), doi:10.1107/S0365110X51001185.
- C. J. Smithells, *Metals Reference Book* (Butterworths Scientific, London, 1955), second edn.

Found in:

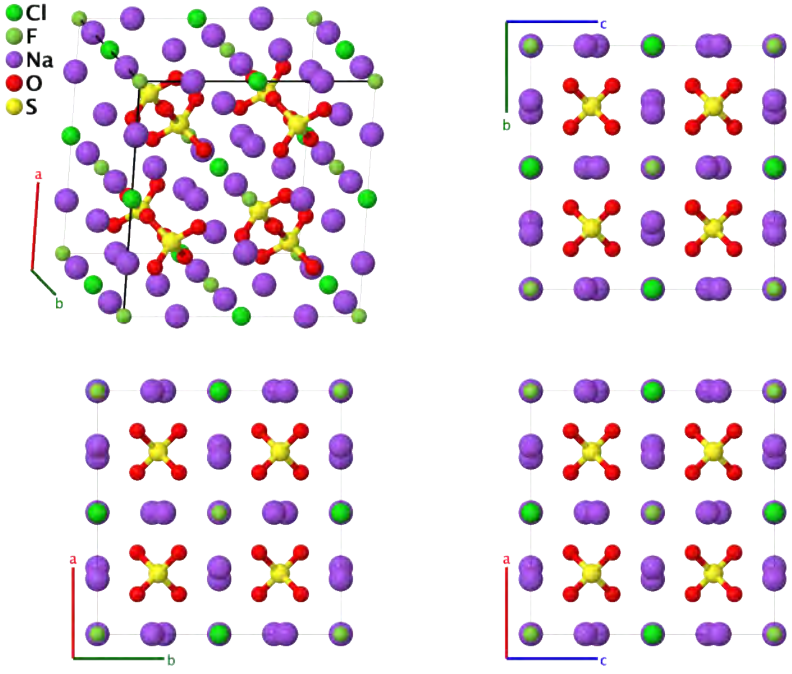
- C. Mshumi, C. I. Lang, L. R. Richey, K. C. Erb, C. W. Rosenbrock, L. J. Nelson, R. R. Vanfleet, H. T. Stokes, B. J. Campbell, and G. L. W. Hart, *Revisiting the $CuPt_3$ prototype and the $L1_3$ structure*, Acta Mater. **73**, 326–336 (2014), doi:10.1016/j.actamat.2014.03.029.

Geometry files:

- CIF: pp. 1818
- POSCAR: pp. 1819

Sulphohalite [$Na_6ClF(SO_4)_2$, $H5_8$] Structure: ABC6D8E2_cF72_225_b_a_e_f_c

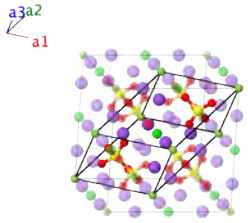
http://aflow.org/prototype-encyclopedia/ABC6D8E2_cF72_225_b_a_e_f_c



Prototype	:	$ClFNa_6O_8S_2$
AFLOW prototype label	:	ABC6D8E2_cF72_225_b_a_e_f_c
***Strukturbericht* designation**	:	$H5_8$
Pearson symbol	:	cF72
Space group number	:	225
Space group symbol	:	$Fm\bar{3}m$
AFLOW prototype command	:	`aflow --proto=ABC6D8E2_cF72_225_b_a_e_f_c --params=`a, x_4, x_5

Face-centered Cubic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	$(4a)$	F

$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(4*b*)	Cl
$\mathbf{B}_3$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(8*c*)	S
$\mathbf{B}_4$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{3}{4}a\,\hat{\mathbf{x}} + \frac{3}{4}a\,\hat{\mathbf{y}} + \frac{3}{4}a\,\hat{\mathbf{z}}$	(8*c*)	S
$\mathbf{B}_5$	=	$-x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}}$	(24*e*)	Na
$\mathbf{B}_6$	=	$x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}}$	(24*e*)	Na
$\mathbf{B}_7$	=	$x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{y}}$	(24*e*)	Na
$\mathbf{B}_8$	=	$-x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{y}}$	(24*e*)	Na
$\mathbf{B}_9$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{z}}$	(24*e*)	Na
$\mathbf{B}_{10}$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{z}}$	(24*e*)	Na
$\mathbf{B}_{11}$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}}$	(32*f*)	O
$\mathbf{B}_{12}$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 - 3x_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}}$	(32*f*)	O
$\mathbf{B}_{13}$	=	$x_5\,\mathbf{a}_1 - 3x_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}}$	(32*f*)	O
$\mathbf{B}_{14}$	=	$-3x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}}$	(32*f*)	O
$\mathbf{B}_{15}$	=	$-x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + 3x_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}}$	(32*f*)	O
$\mathbf{B}_{16}$	=	$-x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - x_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}}$	(32*f*)	O
$\mathbf{B}_{17}$	=	$-x_5\,\mathbf{a}_1 + 3x_5\,\mathbf{a}_2 - x_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}}$	(32*f*)	O
$\mathbf{B}_{18}$	=	$3x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 - x_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}}$	(32*f*)	O

References:

- A. Pabst, *The Crystal Structure of Sulphohalite*, Zeitschrift für Kristallographie - Crystalline Materials **89**, 514–517 (1934), doi:10.1524/zkri.1934.89.1.514.

Found in:

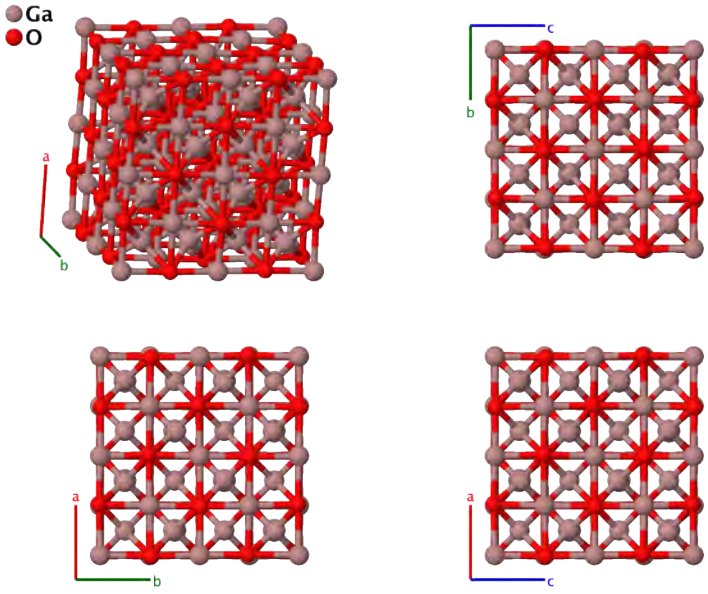
- C. Gottfried and F. Schossberger, eds., *Strukturbericht Band III 1933-1935* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1819
- POSCAR: pp. 1820

γ-Ga_2O_3 Structure: A11B4_cF120_227_acdf_e

http://aflow.org/prototype-encyclopedia/A11B4_cF120_227_acdf_e



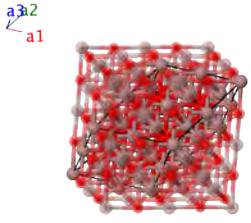
Prototype	:	Ga_2O_3
AFLOW prototype label	:	A11B4_cF120_227_acdf_e
***Strukturbericht* designation**	:	None
Pearson symbol	:	cF120
Space group number	:	227
Space group symbol	:	$Fd\bar{3}m$
AFLOW prototype command	:	`aflow --proto=A11B4_cF120_227_acdf_e --params=`a, x_4, x_5

- Ga_2O_3 takes on a variety of structures:
  - α-Ga_2O_3, which has the corundum ($D5_1$) structure,
  - β-Ga_2O_3,
  - γ-Ga_2O_3, this structure, and
  - ϵ-Ga_2O_3, a structure with many vacancies which can be approximated by the κ-alumina structure.

  Note that all of the gallium sites are only partially occupied.

Face-centered Cubic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{8}\,\mathbf{a}_1 + \frac{1}{8}\,\mathbf{a}_2 + \frac{1}{8}\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(8*a*)	Ga I
$\mathbf{B}_2$	=	$\frac{7}{8}\,\mathbf{a}_1 + \frac{7}{8}\,\mathbf{a}_2 + \frac{7}{8}\,\mathbf{a}_3$	=	$\frac{7}{8}a\,\hat{\mathbf{x}} + \frac{7}{8}a\,\hat{\mathbf{y}} + \frac{7}{8}a\,\hat{\mathbf{z}}$	(8*a*)	Ga I
$\mathbf{B}_3$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(16*c*)	Ga II
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}}$	(16*c*)	Ga II
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(16*c*)	Ga II
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$\frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(16*c*)	Ga II
$\mathbf{B}_7$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(16*d*)	Ga III
$\mathbf{B}_8$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(16*d*)	Ga III
$\mathbf{B}_9$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(16*d*)	Ga III
$\mathbf{B}_{10}$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(16*d*)	Ga III
$\mathbf{B}_{11}$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	(32*e*)	O
$\mathbf{B}_{12}$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \left(\frac{1}{2} - 3x_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	(32*e*)	O
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2} - 3x_4\right)\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{z}}$	(32*e*)	O
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} - 3x_4\right)\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{z}}$	(32*e*)	O
$\mathbf{B}_{15}$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \left(\frac{1}{2} + 3x_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_4\right)a\,\hat{\mathbf{y}} - x_4a\,\hat{\mathbf{z}}$	(32*e*)	O
$\mathbf{B}_{16}$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} - x_4a\,\hat{\mathbf{z}}$	(32*e*)	O
$\mathbf{B}_{17}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + 3x_4\right)\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_4\right)a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + x_4\right)a\,\hat{\mathbf{z}}$	(32*e*)	O
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} + 3x_4\right)\mathbf{a}_1 - x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + x_4\right)a\,\hat{\mathbf{z}}$	(32*e*)	O
$\mathbf{B}_{19}$	=	$\left(\frac{1}{4} - x_5\right)\mathbf{a}_1 + x_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Ga IV
$\mathbf{B}_{20}$	=	$x_5\,\mathbf{a}_1 + \left(\frac{1}{4} - x_5\right)\mathbf{a}_2 + \left(\frac{1}{4} - x_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_5\right)a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Ga IV
$\mathbf{B}_{21}$	=	$x_5\,\mathbf{a}_1 + \left(\frac{1}{4} - x_5\right)\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + x_5a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Ga IV
$\mathbf{B}_{22}$	=	$\left(\frac{1}{4} - x_5\right)\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \left(\frac{1}{4} - x_5\right)\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_5\right)a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Ga IV
$\mathbf{B}_{23}$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \left(\frac{1}{4} - x_5\right)\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + x_5a\,\hat{\mathbf{z}}$	(48*f*)	Ga IV
$\mathbf{B}_{24}$	=	$\left(\frac{1}{4} - x_5\right)\mathbf{a}_1 + \left(\frac{1}{4} - x_5\right)\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_5\right)a\,\hat{\mathbf{z}}$	(48*f*)	Ga IV
$\mathbf{B}_{25}$	=	$\left(\frac{3}{4} + x_5\right)\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \left(\frac{3}{4} + x_5\right)\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} + \left(\frac{3}{4} + x_5\right)a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Ga IV
$\mathbf{B}_{26}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{3}{4} + x_5\right)\mathbf{a}_2 - x_5\,\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} - x_5a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Ga IV
$\mathbf{B}_{27}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{3}{4} + x_5\right)\mathbf{a}_2 + \left(\frac{3}{4} + x_5\right)\mathbf{a}_3$	=	$\left(\frac{3}{4} + x_5\right)a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Ga IV
$\mathbf{B}_{28}$	=	$\left(\frac{3}{4} + x_5\right)\mathbf{a}_1 - x_5\,\mathbf{a}_2 - x_5\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Ga IV
$\mathbf{B}_{29}$	=	$-x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \left(\frac{3}{4} + x_5\right)\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} - x_5a\,\hat{\mathbf{z}}$	(48*f*)	Ga IV
$\mathbf{B}_{30}$	=	$\left(\frac{3}{4} + x_5\right)\mathbf{a}_1 + \left(\frac{3}{4} + x_5\right)\mathbf{a}_2 - x_5\,\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} + \left(\frac{3}{4} + x_5\right)a\,\hat{\mathbf{z}}$	(48*f*)	Ga IV

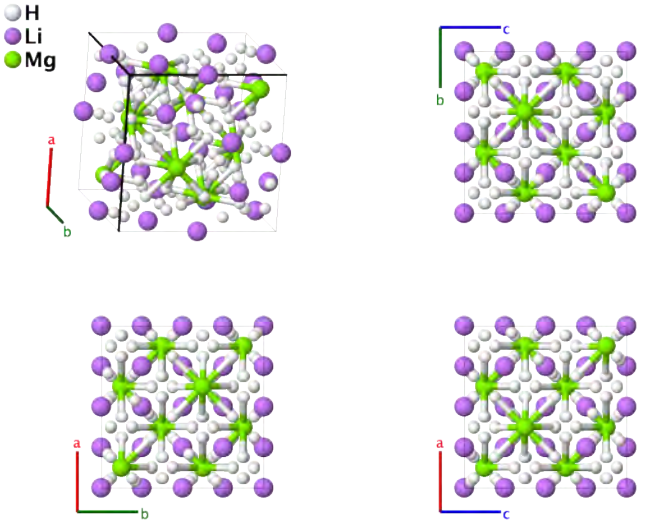
References:

- H. Y. Playford, A. C. Hannon, E. R. Barney, and R. I. Walton, *Structures of Uncharacterised Polymorphs of Gallium Oxide from Total Neutron Diffraction*, Chem.: Eur. J. **19**, 2803–2813 (2013), doi:10.1002/chem.201203359.

Geometry files:
- CIF: pp. 1820
- POSCAR: pp. 1821

Predicted Li_2MgH_{16} High-Temperature Superconductor (250 GPa) Structure: A16B2C_cF152_227_eg_d_a

http://aflow.org/prototype-encyclopedia/A16B2C_cF152_227_eg_d_a

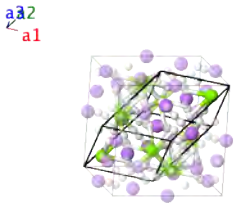


Prototype	:	$H_{16}Li_2Mg$
AFLOW prototype label	:	A16B2C_cF152_227_eg_d_a
***Strukturbericht* designation**	:	None
Pearson symbol	:	cF152
Space group number	:	227
Space group symbol	:	$Fd\bar{3}m$
AFLOW prototype command	:	`aflow --proto=A16B2C_cF152_227_eg_d_a --params=`a, x_3, x_4, z_4

- This structure was predicted by (Sun, 2019) as a metastable state of Li_2MgH_{16} at 250 GPa and $T = 0$ K. If it is possible to construct this compound, or if it becomes stable due to thermodynamic considerations, it is predicted to have a superconducting transition T_c between 430 and 473 K.
- The predicted $T = 0$ K ground state at 300 GPa is a $P\bar{3}m1$ #164 structure with molecular hydrogen.
- (Sun, 2019) give the Wyckoff positions in setting 1 of space group $Fd\bar{3}m$ #227. (They list the Wyckoff positions of the magnesium and lithium atoms as (8*b*) and (16*c*), respectively. They are actually at (8*a*) and (16*d*), as in their 300 GPa data.) We used FINDSYM to shift this to our standard setting 2.

Face-centered Cubic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{y}} + \tfrac{1}{2}\,a\,\hat{\mathbf{z}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{z}}\\ \mathbf{a}_3 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{y}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{8}\,\mathbf{a}_1 + \frac{1}{8}\,\mathbf{a}_2 + \frac{1}{8}\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(8*a*)	Mg
$\mathbf{B}_2$	=	$\frac{7}{8}\,\mathbf{a}_1 + \frac{7}{8}\,\mathbf{a}_2 + \frac{7}{8}\,\mathbf{a}_3$	=	$\frac{7}{8}a\,\hat{\mathbf{x}} + \frac{7}{8}a\,\hat{\mathbf{y}} + \frac{7}{8}a\,\hat{\mathbf{z}}$	(8*a*)	Mg
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(16*d*)	Li
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(16*d*)	Li
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(16*d*)	Li
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(16*d*)	Li
$\mathbf{B}_7$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(32*e*)	H I
$\mathbf{B}_8$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \left(\frac{1}{2} - 3x_3\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(32*e*)	H I
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - 3x_3\right)\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{z}}$	(32*e*)	H I
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - 3x_3\right)\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{z}}$	(32*e*)	H I
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\frac{1}{2} + 3x_3\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(32*e*)	H I
$\mathbf{B}_{12}$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(32*e*)	H I
$\mathbf{B}_{13}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + 3x_3\right)\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{z}}$	(32*e*)	H I
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} + 3x_3\right)\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{z}}$	(32*e*)	H I
$\mathbf{B}_{15}$	=	$z_4\,\mathbf{a}_1 + z_4\,\mathbf{a}_2 + (2x_4 - z_4)\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + z_4a\,\hat{\mathbf{z}}$	(96*g*)	H II
$\mathbf{B}_{16}$	=	$z_4\,\mathbf{a}_1 + z_4\,\mathbf{a}_2 + \left(\frac{1}{2} - 2x_4 - z_4\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{y}} + z_4a\,\hat{\mathbf{z}}$	(96*g*)	H II
$\mathbf{B}_{17}$	=	$(2x_4 - z_4)\,\mathbf{a}_1 + \left(\frac{1}{2} - 2x_4 - z_4\right)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_4\right)a\,\hat{\mathbf{z}}$	(96*g*)	H II
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} - 2x_4 - z_4\right)\,\mathbf{a}_1 + (2x_4 - z_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_4\right)a\,\hat{\mathbf{z}}$	(96*g*)	H II
$\mathbf{B}_{19}$	=	$(2x_4 - z_4)\,\mathbf{a}_1 + z_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$z_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	(96*g*)	H II
$\mathbf{B}_{20}$	=	$\left(\frac{1}{2} - 2x_4 - z_4\right)\,\mathbf{a}_1 + z_4\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$z_4a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{z}}$	(96*g*)	H II
$\mathbf{B}_{21}$	=	$z_4\,\mathbf{a}_1 + (2x_4 - z_4)\,\mathbf{a}_2 + \left(\frac{1}{2} - 2x_4 - z_4\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - z_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	(96*g*)	H II
$\mathbf{B}_{22}$	=	$z_4\,\mathbf{a}_1 + \left(\frac{1}{2} - 2x_4 - z_4\right)\,\mathbf{a}_2 + (2x_4 - z_4)\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - z_4\right)a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{z}}$	(96*g*)	H II
$\mathbf{B}_{23}$	=	$z_4\,\mathbf{a}_1 + (2x_4 - z_4)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + z_4a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	(96*g*)	H II
$\mathbf{B}_{24}$	=	$z_4\,\mathbf{a}_1 + \left(\frac{1}{2} - 2x_4 - z_4\right)\,\mathbf{a}_2 + z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{x}} + z_4a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{z}}$	(96*g*)	H II

$\mathbf{B}_{25}$	=	$\left(\frac{1}{2}-2x_4-z_4\right)\mathbf{a}_1+z_4\,\mathbf{a}_2+(2x_4-z_4)\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-z_4\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-x_4\right)a\,\hat{\mathbf{z}}$	(96*g*)	H II
$\mathbf{B}_{26}$	=	$(2x_4-z_4)\,\mathbf{a}_1+z_4\,\mathbf{a}_2+\left(\frac{1}{2}-2x_4-z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}-x_4\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-z_4\right)a\,\hat{\mathbf{y}}+x_4a\,\hat{\mathbf{z}}$	(96*g*)	H II
$\mathbf{B}_{27}$	=	$-z_4\,\mathbf{a}_1-z_4\,\mathbf{a}_2+\left(\frac{1}{2}+2x_4+z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}+x_4\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+x_4\right)a\,\hat{\mathbf{y}}-z_4a\,\hat{\mathbf{z}}$	(96*g*)	H II
$\mathbf{B}_{28}$	=	$-z_4\,\mathbf{a}_1-z_4\,\mathbf{a}_2+(-2x_4+z_4)\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}}-x_4a\,\hat{\mathbf{y}}-z_4a\,\hat{\mathbf{z}}$	(96*g*)	H II
$\mathbf{B}_{29}$	=	$(-2x_4+z_4)\,\mathbf{a}_1+\left(\frac{1}{2}+2x_4+z_4\right)\mathbf{a}_2-z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{4}+x_4\right)a\,\hat{\mathbf{x}}-x_4a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+z_4\right)a\,\hat{\mathbf{z}}$	(96*g*)	H II
$\mathbf{B}_{30}$	=	$\left(\frac{1}{2}+2x_4+z_4\right)\mathbf{a}_1+(-2x_4+z_4)\,\mathbf{a}_2-z_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+x_4\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+z_4\right)a\,\hat{\mathbf{z}}$	(96*g*)	H II
$\mathbf{B}_{31}$	=	$(-2x_4+z_4)\,\mathbf{a}_1-z_4\,\mathbf{a}_2+\left(\frac{1}{2}+2x_4+z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}+x_4\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+z_4\right)a\,\hat{\mathbf{y}}-x_4a\,\hat{\mathbf{z}}$	(96*g*)	H II
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2}+2x_4+z_4\right)\mathbf{a}_1-z_4\,\mathbf{a}_2+(-2x_4+z_4)\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+z_4\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+x_4\right)a\,\hat{\mathbf{z}}$	(96*g*)	H II
$\mathbf{B}_{33}$	=	$-z_4\,\mathbf{a}_1+(-2x_4+z_4)\,\mathbf{a}_2-z_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}}-z_4a\,\hat{\mathbf{y}}-x_4a\,\hat{\mathbf{z}}$	(96*g*)	H II
$\mathbf{B}_{34}$	=	$-z_4\,\mathbf{a}_1+\left(\frac{1}{2}+2x_4+z_4\right)\mathbf{a}_2-z_4\,\mathbf{a}_3$	=	$\left(\frac{1}{4}+x_4\right)a\,\hat{\mathbf{x}}-z_4a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+x_4\right)a\,\hat{\mathbf{z}}$	(96*g*)	H II
$\mathbf{B}_{35}$	=	$-z_4\,\mathbf{a}_1+(-2x_4+z_4)\,\mathbf{a}_2+\left(\frac{1}{2}+2x_4+z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}+z_4\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+x_4\right)a\,\hat{\mathbf{y}}-x_4a\,\hat{\mathbf{z}}$	(96*g*)	H II
$\mathbf{B}_{36}$	=	$-z_4\,\mathbf{a}_1+\left(\frac{1}{2}+2x_4+z_4\right)\mathbf{a}_2+(-2x_4+z_4)\,\mathbf{a}_3$	=	$\left(\frac{1}{4}+z_4\right)a\,\hat{\mathbf{x}}-x_4a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+x_4\right)a\,\hat{\mathbf{z}}$	(96*g*)	H II
$\mathbf{B}_{37}$	=	$\left(\frac{1}{2}+2x_4+z_4\right)\mathbf{a}_1-z_4\,\mathbf{a}_2-z_4\,\mathbf{a}_3$	=	$-z_4a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+x_4\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+x_4\right)a\,\hat{\mathbf{z}}$	(96*g*)	H II
$\mathbf{B}_{38}$	=	$(-2x_4+z_4)\,\mathbf{a}_1-z_4\,\mathbf{a}_2-z_4\,\mathbf{a}_3$	=	$-z_4a\,\hat{\mathbf{x}}-x_4a\,\hat{\mathbf{y}}-x_4a\,\hat{\mathbf{z}}$	(96*g*)	H II

References:

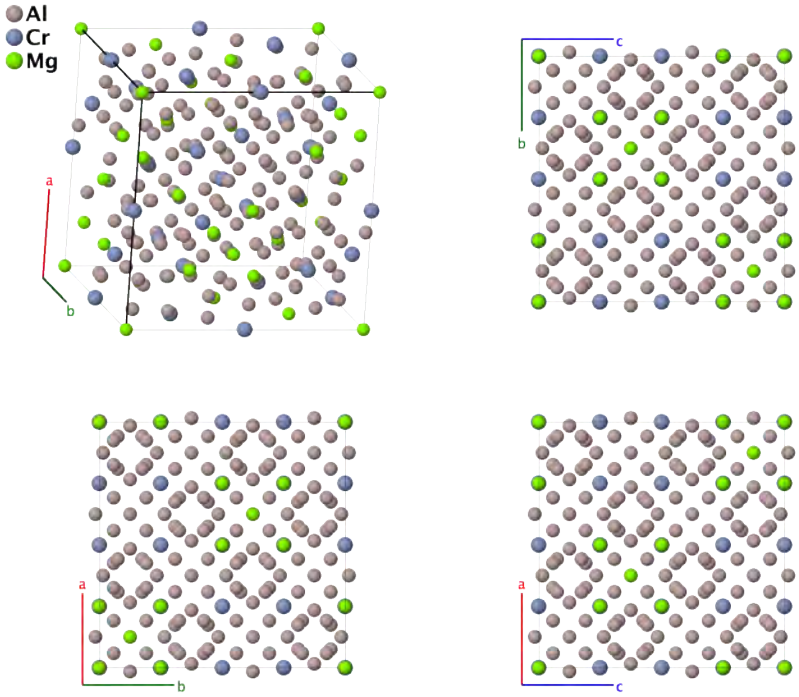
- Y. Sun, J. Lv, Y. Xie, H. Liu, and Y. Ma, *Route to a Superconducting Phase above Room Temperature in Electron-Doped Hydride Compounds under High Pressure*, Phys. Rev. Lett. **123**, 097001 (2019), doi:10.1103/PhysRevLett.123.097001.

Geometry files:

- CIF: pp. 1822
- POSCAR: pp. 1823

$Mg_3Cr_2Al_{18}$ Structure: A18B2C3_cF184_227_fg_d_ac

http://aflow.org/prototype-encyclopedia/A18B2C3_cF184_227_fg_d_ac



Prototype	:	$Al_{18}Cr_2Mg_3$
AFLOW prototype label	:	A18B2C3_cF184_227_fg_d_ac
***Strukturbericht* designation**	:	None
Pearson symbol	:	cF184
Space group number	:	227
Space group symbol	:	$Fd\bar{3}m$
AFLOW prototype command	:	`aflow --proto=A18B2C3_cF184_227_fg_d_ac --params=`a, x_4, x_5, z_5

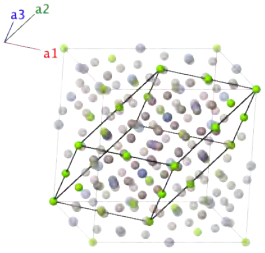
Other compounds with this structure

- $LaCr_2Al_{20}$, $CeCr_2Al_{20}$, $PrCr_2Al_{20}$, $SmCr_2Al_{20}$, $YbCr_2Al_{20}$, $CeTi_2Al_{20}$, $PrTi_2Al_{20}$, $SmTi_2Al_{20}$, $YbTi_2Al_{20}$, CeV_2Al_{20}, GdV_2Al_{20}, LaV_2Al_{20}, PrV_2Al_{20}, SmV_2Al_{20}, $CeNi_2Cd_{20}$, $GdNi_2Cd_{20}$, $LaNi_2Cd_{20}$, $NdNi_2Cd_{20}$, $PrNi_2Cd_{20}$, $SmNi_2Cd_{20}$, YNi_2Cd_{20}, $CePd_2Cd_{20}$, $PrPd_2Cd_{20}$, $SmPd_2Cd_{20}$, and UOs_2Zn_{20}

- (Samson, 1958) gives the atomic coordinates in terms of setting 1 of space group $Fd\bar{3}m$ #227. We have shifted this to the standard setting 2, where the inversion site of the lattice is at the origin.
- If the (8*a*), (16*c*) and (16*d*) sites are occupied by the same type of atom this becomes the $Zn_{22}Zr$ structure.
- In the ternary compounds $LnMX_{20}$, the rare earth (*Ln*) metal occupies the (8*a*) site, the transition metal (*M*) the (16*d*) site, and *X*=Al,Cd,Zn occupies the (16*c*), (48*f*) and (96*g*) sites. These compounds are sometimes listed under the $CeCr_2Al_{20}$ prototype.

Face-centered Cubic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{y}} + \tfrac{1}{2}\,a\,\hat{\mathbf{z}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{z}}\\ \mathbf{a}_3 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{y}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{8}\,\mathbf{a}_1 + \frac{1}{8}\,\mathbf{a}_2 + \frac{1}{8}\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(8*a*)	Mg I
$\mathbf{B}_2$	=	$\frac{7}{8}\,\mathbf{a}_1 + \frac{7}{8}\,\mathbf{a}_2 + \frac{7}{8}\,\mathbf{a}_3$	=	$\frac{7}{8}a\,\hat{\mathbf{x}} + \frac{7}{8}a\,\hat{\mathbf{y}} + \frac{7}{8}a\,\hat{\mathbf{z}}$	(8*a*)	Mg I
$\mathbf{B}_3$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(16*c*)	Mg II
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}}$	(16*c*)	Mg II
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(16*c*)	Mg II
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$\frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(16*c*)	Mg II
$\mathbf{B}_7$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(16*d*)	Cr
$\mathbf{B}_8$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(16*d*)	Cr
$\mathbf{B}_9$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(16*d*)	Cr
$\mathbf{B}_{10}$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(16*d*)	Cr
$\mathbf{B}_{11}$	=	$\left(\frac{1}{4} - x_4\right)\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Al I
$\mathbf{B}_{12}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{4} - x_4\right)\,\mathbf{a}_2 + \left(\frac{1}{4} - x_4\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Al I
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{4} - x_4\right)\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Al I
$\mathbf{B}_{14}$	=	$\left(\frac{1}{4} - x_4\right)\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \left(\frac{1}{4} - x_4\right)\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Al I
$\mathbf{B}_{15}$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \left(\frac{1}{4} - x_4\right)\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + x_4 a\,\hat{\mathbf{z}}$	(48*f*)	Al I
$\mathbf{B}_{16}$	=	$\left(\frac{1}{4} - x_4\right)\,\mathbf{a}_1 + \left(\frac{1}{4} - x_4\right)\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{z}}$	(48*f*)	Al I
$\mathbf{B}_{17}$	=	$\left(\frac{3}{4} + x_4\right)\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \left(\frac{3}{4} + x_4\right)\,\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} + \left(\frac{3}{4} + x_4\right)a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Al I
$\mathbf{B}_{18}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{3}{4} + x_4\right)\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Al I
$\mathbf{B}_{19}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{3}{4} + x_4\right)\,\mathbf{a}_2 + \left(\frac{3}{4} + x_4\right)\,\mathbf{a}_3$	=	$\left(\frac{3}{4} + x_4\right)a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Al I
$\mathbf{B}_{20}$	=	$\left(\frac{3}{4} + x_4\right)\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Al I
$\mathbf{B}_{21}$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \left(\frac{3}{4} + x_4\right)\,\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} - x_4 a\,\hat{\mathbf{z}}$	(48*f*)	Al I
$\mathbf{B}_{22}$	=	$\left(\frac{3}{4} + x_4\right)\,\mathbf{a}_1 + \left(\frac{3}{4} + x_4\right)\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} + \left(\frac{3}{4} + x_4\right)a\,\hat{\mathbf{z}}$	(48*f*)	Al I
$\mathbf{B}_{23}$	=	$z_5\,\mathbf{a}_1 + z_5\,\mathbf{a}_2 + (2x_5 - z_5)\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + z_5 a\,\hat{\mathbf{z}}$	(96*g*)	Al II
$\mathbf{B}_{24}$	=	$z_5\,\mathbf{a}_1 + z_5\,\mathbf{a}_2 + \left(\frac{1}{2} - 2x_5 - z_5\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_5\right)a\,\hat{\mathbf{y}} + z_5 a\,\hat{\mathbf{z}}$	(96*g*)	Al II
$\mathbf{B}_{25}$	=	$(2x_5 - z_5)\,\mathbf{a}_1 + \left(\frac{1}{2} - 2x_5 - z_5\right)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_5\right)a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_5\right)a\,\hat{\mathbf{z}}$	(96*g*)	Al II
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} - 2x_5 - z_5\right)\,\mathbf{a}_1 + (2x_5 - z_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_5\right)a\,\hat{\mathbf{z}}$	(96*g*)	Al II

$\mathbf{B}_{27}$	=	$(2x_5 - z_5)\,\mathbf{a}_1 + z_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$z_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}}$	(96*g*)	Al II
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} - 2x_5 - z_5\right)\mathbf{a}_1 + z_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$z_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_5\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_5\right) a\,\hat{\mathbf{z}}$	(96*g*)	Al II
$\mathbf{B}_{29}$	=	$z_5\,\mathbf{a}_1 + (2x_5 - z_5)\,\mathbf{a}_2 + \left(\frac{1}{2} - 2x_5 - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - z_5\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_5\right) a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}}$	(96*g*)	Al II
$\mathbf{B}_{30}$	=	$z_5\,\mathbf{a}_1 + \left(\frac{1}{2} - 2x_5 - z_5\right)\mathbf{a}_2 + (2x_5 - z_5)\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - z_5\right) a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_5\right) a\,\hat{\mathbf{z}}$	(96*g*)	Al II
$\mathbf{B}_{31}$	=	$z_5\,\mathbf{a}_1 + (2x_5 - z_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + z_5 a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}}$	(96*g*)	Al II
$\mathbf{B}_{32}$	=	$z_5\,\mathbf{a}_1 + \left(\frac{1}{2} - 2x_5 - z_5\right)\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_5\right) a\,\hat{\mathbf{x}} + z_5 a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_5\right) a\,\hat{\mathbf{z}}$	(96*g*)	Al II
$\mathbf{B}_{33}$	=	$\left(\frac{1}{2} - 2x_5 - z_5\right)\mathbf{a}_1 + z_5\,\mathbf{a}_2 + (2x_5 - z_5)\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - z_5\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_5\right) a\,\hat{\mathbf{z}}$	(96*g*)	Al II
$\mathbf{B}_{34}$	=	$(2x_5 - z_5)\,\mathbf{a}_1 + z_5\,\mathbf{a}_2 + \left(\frac{1}{2} - 2x_5 - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_5\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - z_5\right) a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}}$	(96*g*)	Al II
$\mathbf{B}_{35}$	=	$-z_5\,\mathbf{a}_1 - z_5\,\mathbf{a}_2 + \left(\frac{1}{2} + 2x_5 + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_5\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_5\right) a\,\hat{\mathbf{y}} - z_5 a\,\hat{\mathbf{z}}$	(96*g*)	Al II
$\mathbf{B}_{36}$	=	$-z_5\,\mathbf{a}_1 - z_5\,\mathbf{a}_2 + (-2x_5 + z_5)\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} - z_5 a\,\hat{\mathbf{z}}$	(96*g*)	Al II
$\mathbf{B}_{37}$	=	$(-2x_5 + z_5)\,\mathbf{a}_1 + \left(\frac{1}{2} + 2x_5 + z_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_5\right) a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_5\right) a\,\hat{\mathbf{z}}$	(96*g*)	Al II
$\mathbf{B}_{38}$	=	$\left(\frac{1}{2} + 2x_5 + z_5\right)\mathbf{a}_1 + (-2x_5 + z_5)\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_5\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_5\right) a\,\hat{\mathbf{z}}$	(96*g*)	Al II
$\mathbf{B}_{39}$	=	$(-2x_5 + z_5)\,\mathbf{a}_1 - z_5\,\mathbf{a}_2 + \left(\frac{1}{2} + 2x_5 + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_5\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + z_5\right) a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}}$	(96*g*)	Al II
$\mathbf{B}_{40}$	=	$\left(\frac{1}{2} + 2x_5 + z_5\right)\mathbf{a}_1 - z_5\,\mathbf{a}_2 + (-2x_5 + z_5)\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + z_5\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + x_5\right) a\,\hat{\mathbf{z}}$	(96*g*)	Al II
$\mathbf{B}_{41}$	=	$-z_5\,\mathbf{a}_1 + (-2x_5 + z_5)\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - z_5 a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}}$	(96*g*)	Al II
$\mathbf{B}_{42}$	=	$-z_5\,\mathbf{a}_1 + \left(\frac{1}{2} + 2x_5 + z_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_5\right) a\,\hat{\mathbf{x}} - z_5 a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + x_5\right) a\,\hat{\mathbf{z}}$	(96*g*)	Al II
$\mathbf{B}_{43}$	=	$-z_5\,\mathbf{a}_1 + (-2x_5 + z_5)\,\mathbf{a}_2 + \left(\frac{1}{2} + 2x_5 + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + z_5\right) a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_5\right) a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}}$	(96*g*)	Al II
$\mathbf{B}_{44}$	=	$-z_5\,\mathbf{a}_1 + \left(\frac{1}{2} + 2x_5 + z_5\right)\mathbf{a}_2 + (-2x_5 + z_5)\,\mathbf{a}_3$	=	$\left(\frac{1}{4} + z_5\right) a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + x_5\right) a\,\hat{\mathbf{z}}$	(96*g*)	Al II
$\mathbf{B}_{45}$	=	$\left(\frac{1}{2} + 2x_5 + z_5\right)\mathbf{a}_1 - z_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-z_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_5\right) a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + x_5\right) a\,\hat{\mathbf{z}}$	(96*g*)	Al II
$\mathbf{B}_{46}$	=	$(-2x_5 + z_5)\,\mathbf{a}_1 - z_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-z_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}}$	(96*g*)	Al II

References:

- S. Samson, *The Crystal Structure of the Intermetallic Compound $Mg_3Cr_2Al_{18}$*, Acta Cryst. **11**, 851–857 (1958), doi:10.1107/S0365110X58002425.

Found in:

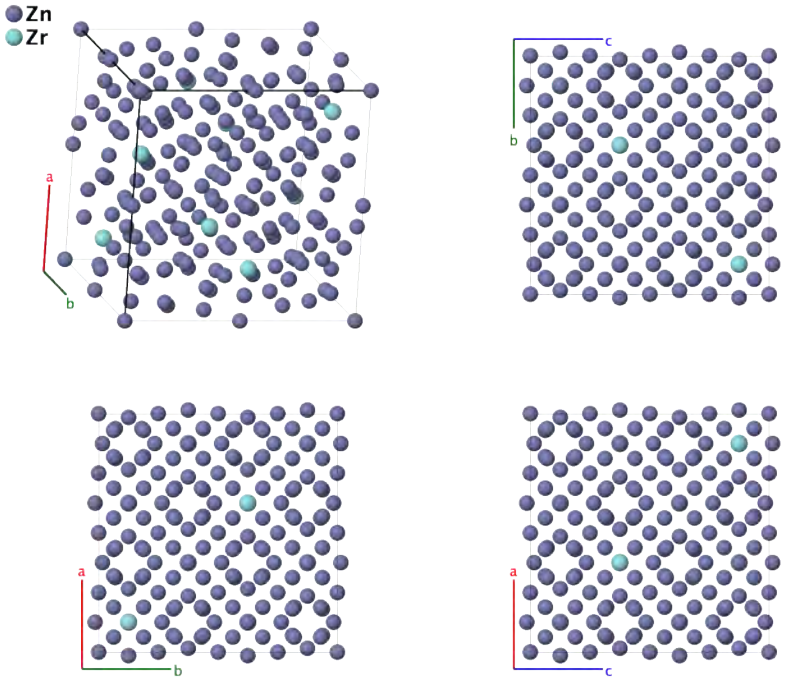
- S. Samson, *The Crystal Structure of the Intermetallic Compound $ZrZn_{22}$*, Acta Cryst. **14**, 1229–1236 (1961), doi:10.1107/S0365110X61003600.

Geometry files:

- CIF: pp. 1823
- POSCAR: pp. 1824

$Zn_{22}Zr$ Structure: A22B_cF184_227_cdfg_a

http://aflow.org/prototype-encyclopedia/A22B_cF184_227_cdfg_a



Prototype	:	$Zn_{22}Zr$
AFLOW prototype label	:	A22B_cF184_227_cdfg_a
***Strukturbericht* designation**	:	None
Pearson symbol	:	cF184
Space group number	:	227
Space group symbol	:	$Fd\bar{3}m$
AFLOW prototype command	:	`aflow --proto=A22B_cF184_227_cdfg_a --params=`a, x_4, x_5, z_5

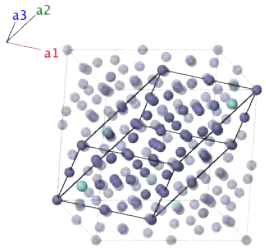
Other compounds with this structure

- $Be_{22}Mo$ and $Be_{22}W$

- (Samson, 1961) gives the atomic coordinates in terms of setting 1 of space group $F\bar{3}m$ #227. We have shifted this to the standard setting 2, where the inversion site of the lattice is at the origin.
- Samson also suggests that the "$Zn_{14}Zr$" structure is created when zirconium atoms replace some of the zinc atoms on the (16*c*) site [the (16*d*) site in Samson's orientation].
- This structure can be derived from the $Mg_3Cr_2Al_{18}$ structure.

Face-centered Cubic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{y}} + \tfrac{1}{2}\,a\,\hat{\mathbf{z}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{z}}\\ \mathbf{a}_3 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{y}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{8}\,\mathbf{a}_1 + \frac{1}{8}\,\mathbf{a}_2 + \frac{1}{8}\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(8*a*)	Zr
$\mathbf{B}_2$	=	$\frac{7}{8}\,\mathbf{a}_1 + \frac{7}{8}\,\mathbf{a}_2 + \frac{7}{8}\,\mathbf{a}_3$	=	$\frac{7}{8}a\,\hat{\mathbf{x}} + \frac{7}{8}a\,\hat{\mathbf{y}} + \frac{7}{8}a\,\hat{\mathbf{z}}$	(8*a*)	Zr
$\mathbf{B}_3$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(16*c*)	Zn I
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}}$	(16*c*)	Zn I
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(16*c*)	Zn I
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$\frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(16*c*)	Zn I
$\mathbf{B}_7$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(16*d*)	Zn II
$\mathbf{B}_8$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(16*d*)	Zn II
$\mathbf{B}_9$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(16*d*)	Zn II
$\mathbf{B}_{10}$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(16*d*)	Zn II
$\mathbf{B}_{11}$	=	$\left(\frac{1}{4} - x_4\right)\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4 a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Zn III
$\mathbf{B}_{12}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{4} - x_4\right)\,\mathbf{a}_2 + \left(\frac{1}{4} - x_4\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Zn III
$\mathbf{B}_{13}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{4} - x_4\right)\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Zn III
$\mathbf{B}_{14}$	=	$\left(\frac{1}{4} - x_4\right)\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \left(\frac{1}{4} - x_4\right)\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Zn III
$\mathbf{B}_{15}$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \left(\frac{1}{4} - x_4\right)\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + x_4 a\,\hat{\mathbf{z}}$	(48*f*)	Zn III
$\mathbf{B}_{16}$	=	$\left(\frac{1}{4} - x_4\right)\,\mathbf{a}_1 + \left(\frac{1}{4} - x_4\right)\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{z}}$	(48*f*)	Zn III
$\mathbf{B}_{17}$	=	$\left(\frac{3}{4} + x_4\right)\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \left(\frac{3}{4} + x_4\right)\,\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} + \left(\frac{3}{4} + x_4\right)a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Zn III
$\mathbf{B}_{18}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{3}{4} + x_4\right)\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Zn III
$\mathbf{B}_{19}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{3}{4} + x_4\right)\,\mathbf{a}_2 + \left(\frac{3}{4} + x_4\right)\,\mathbf{a}_3$	=	$\left(\frac{3}{4} + x_4\right)a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Zn III
$\mathbf{B}_{20}$	=	$\left(\frac{3}{4} + x_4\right)\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$-x_4 a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Zn III
$\mathbf{B}_{21}$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \left(\frac{3}{4} + x_4\right)\,\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} - x_4 a\,\hat{\mathbf{z}}$	(48*f*)	Zn III
$\mathbf{B}_{22}$	=	$\left(\frac{3}{4} + x_4\right)\,\mathbf{a}_1 + \left(\frac{3}{4} + x_4\right)\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} + \left(\frac{3}{4} + x_4\right)a\,\hat{\mathbf{z}}$	(48*f*)	Zn III
$\mathbf{B}_{23}$	=	$z_5\,\mathbf{a}_1 + z_5\,\mathbf{a}_2 + (2x_5 - z_5)\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + z_5 a\,\hat{\mathbf{z}}$	(96*g*)	Zn IV
$\mathbf{B}_{24}$	=	$z_5\,\mathbf{a}_1 + z_5\,\mathbf{a}_2 + \left(\frac{1}{2} - 2x_5 - z_5\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_5\right)a\,\hat{\mathbf{y}} + z_5 a\,\hat{\mathbf{z}}$	(96*g*)	Zn IV
$\mathbf{B}_{25}$	=	$(2x_5 - z_5)\,\mathbf{a}_1 + \left(\frac{1}{2} - 2x_5 - z_5\right)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_5\right)a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_5\right)a\,\hat{\mathbf{z}}$	(96*g*)	Zn IV
$\mathbf{B}_{26}$	=	$\left(\frac{1}{2} - 2x_5 - z_5\right)\,\mathbf{a}_1 + (2x_5 - z_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_5\right)a\,\hat{\mathbf{z}}$	(96*g*)	Zn IV

$\mathbf{B}_{27}$	=	$(2x_5 - z_5)\,\mathbf{a}_1 + z_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$z_5 a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}}$	(96*g*)	Zn IV
$\mathbf{B}_{28}$	=	$\left(\frac{1}{2} - 2x_5 - z_5\right)\mathbf{a}_1 + z_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$z_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_5\right)a\,\hat{\mathbf{z}}$	(96*g*)	Zn IV
$\mathbf{B}_{29}$	=	$z_5\,\mathbf{a}_1 + (2x_5 - z_5)\,\mathbf{a}_2 + \left(\frac{1}{2} - 2x_5 - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - z_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_5\right)a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}}$	(96*g*)	Zn IV
$\mathbf{B}_{30}$	=	$z_5\,\mathbf{a}_1 + \left(\frac{1}{2} - 2x_5 - z_5\right)\mathbf{a}_2 + (2x_5 - z_5)\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - z_5\right)a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_5\right)a\,\hat{\mathbf{z}}$	(96*g*)	Zn IV
$\mathbf{B}_{31}$	=	$z_5\,\mathbf{a}_1 + (2x_5 - z_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + z_5 a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}}$	(96*g*)	Zn IV
$\mathbf{B}_{32}$	=	$z_5\,\mathbf{a}_1 + \left(\frac{1}{2} - 2x_5 - z_5\right)\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_5\right)a\,\hat{\mathbf{x}} + z_5 a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_5\right)a\,\hat{\mathbf{z}}$	(96*g*)	Zn IV
$\mathbf{B}_{33}$	=	$\left(\frac{1}{2} - 2x_5 - z_5\right)\mathbf{a}_1 + z_5\,\mathbf{a}_2 + (2x_5 - z_5)\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - z_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_5\right)a\,\hat{\mathbf{z}}$	(96*g*)	Zn IV
$\mathbf{B}_{34}$	=	$(2x_5 - z_5)\,\mathbf{a}_1 + z_5\,\mathbf{a}_2 + \left(\frac{1}{2} - 2x_5 - z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - z_5\right)a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}}$	(96*g*)	Zn IV
$\mathbf{B}_{35}$	=	$-z_5\,\mathbf{a}_1 - z_5\,\mathbf{a}_2 + \left(\frac{1}{2} + 2x_5 + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_5\right)a\,\hat{\mathbf{y}} - z_5 a\,\hat{\mathbf{z}}$	(96*g*)	Zn IV
$\mathbf{B}_{36}$	=	$-z_5\,\mathbf{a}_1 - z_5\,\mathbf{a}_2 + (-2x_5 + z_5)\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} - z_5 a\,\hat{\mathbf{z}}$	(96*g*)	Zn IV
$\mathbf{B}_{37}$	=	$(-2x_5 + z_5)\,\mathbf{a}_1 + \left(\frac{1}{2} + 2x_5 + z_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_5\right)a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_5\right)a\,\hat{\mathbf{z}}$	(96*g*)	Zn IV
$\mathbf{B}_{38}$	=	$\left(\frac{1}{2} + 2x_5 + z_5\right)\mathbf{a}_1 + (-2x_5 + z_5)\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_5\right)a\,\hat{\mathbf{z}}$	(96*g*)	Zn IV
$\mathbf{B}_{39}$	=	$(-2x_5 + z_5)\,\mathbf{a}_1 - z_5\,\mathbf{a}_2 + \left(\frac{1}{2} + 2x_5 + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + z_5\right)a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}}$	(96*g*)	Zn IV
$\mathbf{B}_{40}$	=	$\left(\frac{1}{2} + 2x_5 + z_5\right)\mathbf{a}_1 - z_5\,\mathbf{a}_2 + (-2x_5 + z_5)\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + z_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + x_5\right)a\,\hat{\mathbf{z}}$	(96*g*)	Zn IV
$\mathbf{B}_{41}$	=	$-z_5\,\mathbf{a}_1 + (-2x_5 + z_5)\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} - z_5 a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}}$	(96*g*)	Zn IV
$\mathbf{B}_{42}$	=	$-z_5\,\mathbf{a}_1 + \left(\frac{1}{2} + 2x_5 + z_5\right)\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_5\right)a\,\hat{\mathbf{x}} - z_5 a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + x_5\right)a\,\hat{\mathbf{z}}$	(96*g*)	Zn IV
$\mathbf{B}_{43}$	=	$-z_5\,\mathbf{a}_1 + (-2x_5 + z_5)\,\mathbf{a}_2 + \left(\frac{1}{2} + 2x_5 + z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + z_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_5\right)a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}}$	(96*g*)	Zn IV
$\mathbf{B}_{44}$	=	$-z_5\,\mathbf{a}_1 + \left(\frac{1}{2} + 2x_5 + z_5\right)\mathbf{a}_2 + (-2x_5 + z_5)\,\mathbf{a}_3$	=	$\left(\frac{1}{4} + z_5\right)a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + x_5\right)a\,\hat{\mathbf{z}}$	(96*g*)	Zn IV
$\mathbf{B}_{45}$	=	$\left(\frac{1}{2} + 2x_5 + z_5\right)\mathbf{a}_1 - z_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-z_5 a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + x_5\right)a\,\hat{\mathbf{z}}$	(96*g*)	Zn IV
$\mathbf{B}_{46}$	=	$(-2x_5 + z_5)\,\mathbf{a}_1 - z_5\,\mathbf{a}_2 - z_5\,\mathbf{a}_3$	=	$-z_5 a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}}$	(96*g*)	Zn IV

References:

- S. Samson, *The Crystal Structure of the Intermetallic Compound $ZrZn_{22}$*, Acta Cryst. **14**, 1229–1236 (1961), doi:10.1107/S0365110X61003600.

Found in:

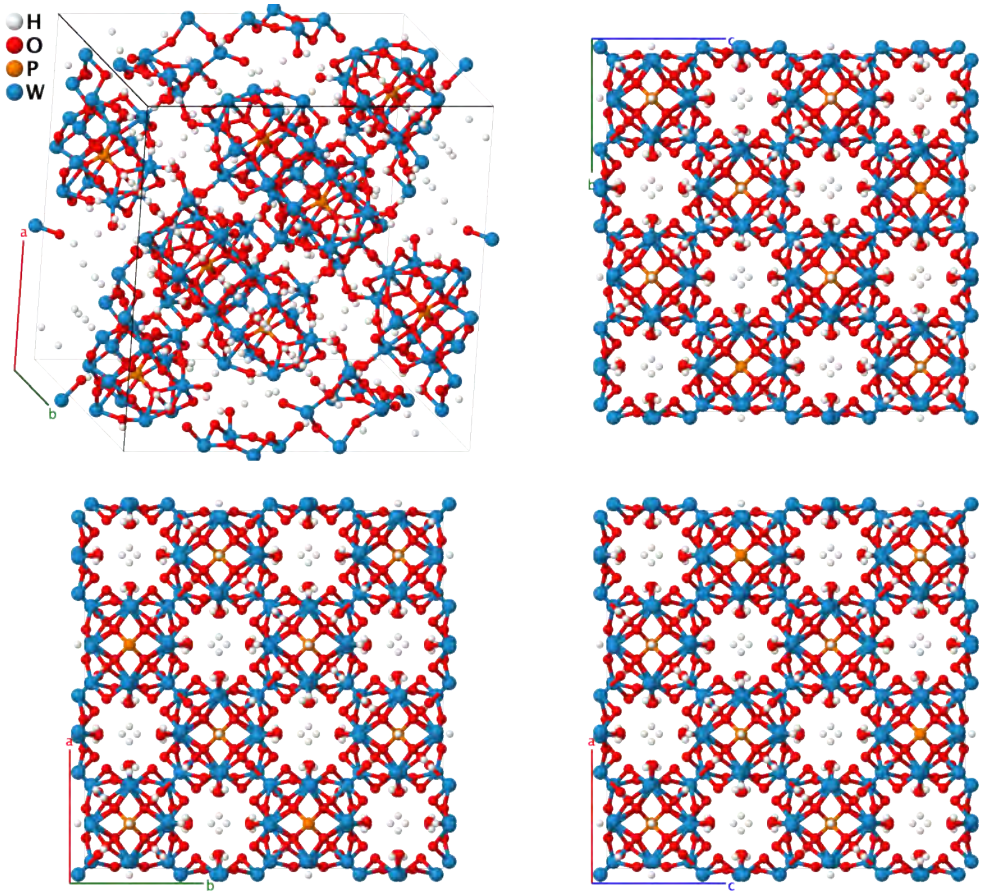
- T. B. Massalski, H. Okamoto, P. R. Subramanian, and L. Kacprzak, eds., *Binary Alloy Phase Diagrams*, vol. 3 (ASM International, Materials Park, Ohio, USA, 1990), 2^{nd} edn. Hf-Re to Zn-Zr.

Geometry files:

- CIF: pp. 1825
- POSCAR: pp. 1826

$H_3PW_{12}O_{40}\cdot 29H_2O$ ($H4_{21}$) Structure: A29B40CD12_cF656_227_ae2fg_e3g_b_g

http://aflow.org/prototype-encyclopedia/A29B40CD12_cF656_227_ae2fg_e3g_b_g



Prototype	:	$(H_2O)_{29}O_{40}PW_{12}$
AFLOW prototype label	:	A29B40CD12_cF656_227_ae2fg_e3g_b_g
***Strukturbericht* designation**	:	$H4_{21}$
Pearson symbol	:	cF656
Space group number	:	227
Space group symbol	:	$Fd\bar{3}m$
AFLOW prototype command	:	`aflow --proto=A29B40CD12_cF656_227_ae2fg_e3g_b_g --params=`$a, x_3, x_4, x_5, x_6, x_7, z_7, x_8, z_8, x_9, z_9, x_{10}, z_{10}, x_{11}, z_{11}$

Other compounds with this structure

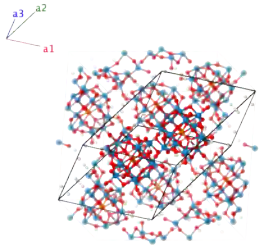
- $H_3PMo_{12}O_{40}\cdot 30H_2O$

- This compound is often colloquially called "PWA-29." On heating some water molecules will disassociate, leaving $H_3PW_{12}O_{40}\cdot 6H_2O$, $H_3PW_{12}O_{40}\cdot 5H_2O$ ($H4_{16}$), or $H_3PW_{12}O_{40}\cdot 3H_2O$.
- The three hydrogen atoms not formally associated with the water molecules are not located. Presumably they join with some water molecules to form H_3O^+ ions.

- Even the exact number and position of the water molecules is uncertain. (Clark, 1976), studying the related compound $H_3PMo_{12}O_{40}{\cdot}30H_2O$, states the the composition is approximately $30H_2O$, and that "only six of the water molecules occupy ordered sites."

Face-centered Cubic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{y}} + \tfrac{1}{2}\,a\,\hat{\mathbf{z}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{z}}\\ \mathbf{a}_3 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{y}}\end{aligned}$$



Basis vectors:

	Lattice Coordinates	Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	$\frac{1}{8}\,\mathbf{a}_1 + \frac{1}{8}\,\mathbf{a}_2 + \frac{1}{8}\,\mathbf{a}_3$	$\frac{1}{8}a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(8*a*)	H_2O I
$\mathbf{B}_2$	$\frac{7}{8}\,\mathbf{a}_1 + \frac{7}{8}\,\mathbf{a}_2 + \frac{7}{8}\,\mathbf{a}_3$	$\frac{7}{8}a\,\hat{\mathbf{x}} + \frac{7}{8}a\,\hat{\mathbf{y}} + \frac{7}{8}a\,\hat{\mathbf{z}}$	(8*a*)	H_2O I
$\mathbf{B}_3$	$\frac{3}{8}\,\mathbf{a}_1 + \frac{3}{8}\,\mathbf{a}_2 + \frac{3}{8}\,\mathbf{a}_3$	$\frac{3}{8}a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}}$	(8*b*)	P
$\mathbf{B}_4$	$\frac{5}{8}\,\mathbf{a}_1 + \frac{5}{8}\,\mathbf{a}_2 + \frac{5}{8}\,\mathbf{a}_3$	$\frac{5}{8}a\,\hat{\mathbf{x}} + \frac{5}{8}a\,\hat{\mathbf{y}} + \frac{5}{8}a\,\hat{\mathbf{z}}$	(8*b*)	P
$\mathbf{B}_5$	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	$x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(32*e*)	H_2O II
$\mathbf{B}_6$	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \left(\frac{1}{2} - 3x_3\right)\mathbf{a}_3$	$\left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(32*e*)	H_2O II
$\mathbf{B}_7$	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - 3x_3\right)\mathbf{a}_2 + x_3\,\mathbf{a}_3$	$\left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{z}}$	(32*e*)	H_2O II
$\mathbf{B}_8$	$\left(\frac{1}{2} - 3x_3\right)\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	$x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{z}}$	(32*e*)	H_2O II
$\mathbf{B}_9$	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\frac{1}{2} + 3x_3\right)\mathbf{a}_3$	$\left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(32*e*)	H_2O II
$\mathbf{B}_{10}$	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	$-x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(32*e*)	H_2O II
$\mathbf{B}_{11}$	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + 3x_3\right)\mathbf{a}_2 - x_3\,\mathbf{a}_3$	$\left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{z}}$	(32*e*)	H_2O II
$\mathbf{B}_{12}$	$\left(\frac{1}{2} + 3x_3\right)\mathbf{a}_1 - x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	$-x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{z}}$	(32*e*)	H_2O II
$\mathbf{B}_{13}$	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	$x_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	(32*e*)	O I
$\mathbf{B}_{14}$	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \left(\frac{1}{2} - 3x_4\right)\mathbf{a}_3$	$\left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	(32*e*)	O I
$\mathbf{B}_{15}$	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2} - 3x_4\right)\mathbf{a}_2 + x_4\,\mathbf{a}_3$	$\left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{z}}$	(32*e*)	O I
$\mathbf{B}_{16}$	$\left(\frac{1}{2} - 3x_4\right)\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	$x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{z}}$	(32*e*)	O I
$\mathbf{B}_{17}$	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \left(\frac{1}{2} + 3x_4\right)\mathbf{a}_3$	$\left(\frac{1}{4} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_4\right)a\,\hat{\mathbf{y}} - x_4a\,\hat{\mathbf{z}}$	(32*e*)	O I
$\mathbf{B}_{18}$	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	$-x_4a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} - x_4a\,\hat{\mathbf{z}}$	(32*e*)	O I
$\mathbf{B}_{19}$	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + 3x_4\right)\mathbf{a}_2 - x_4\,\mathbf{a}_3$	$\left(\frac{1}{4} + x_4\right)a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + x_4\right)a\,\hat{\mathbf{z}}$	(32*e*)	O I
$\mathbf{B}_{20}$	$\left(\frac{1}{2} + 3x_4\right)\mathbf{a}_1 - x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	$-x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + x_4\right)a\,\hat{\mathbf{z}}$	(32*e*)	O I
$\mathbf{B}_{21}$	$\left(\frac{1}{4} - x_5\right)\mathbf{a}_1 + x_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	$x_5a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(48*f*)	H_2O III
$\mathbf{B}_{22}$	$x_5\,\mathbf{a}_1 + \left(\frac{1}{4} - x_5\right)\mathbf{a}_2 + \left(\frac{1}{4} - x_5\right)\mathbf{a}_3$	$\left(\frac{1}{4} - x_5\right)a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(48*f*)	H_2O III
$\mathbf{B}_{23}$	$x_5\,\mathbf{a}_1 + \left(\frac{1}{4} - x_5\right)\mathbf{a}_2 + x_5\,\mathbf{a}_3$	$\frac{1}{8}a\,\hat{\mathbf{x}} + x_5a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(48*f*)	H_2O III
$\mathbf{B}_{24}$	$\left(\frac{1}{4} - x_5\right)\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \left(\frac{1}{4} - x_5\right)\mathbf{a}_3$	$\frac{1}{8}a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_5\right)a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(48*f*)	H_2O III

$$
\begin{array}{lcccll}
\mathbf{B}_{25} & = & x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \left(\frac{1}{4} - x_5\right)\mathbf{a}_3 & = & \frac{1}{8}a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}} & (48f)\ \mathrm{H_2O\ III} \\
\mathbf{B}_{26} & = & \left(\frac{1}{4} - x_5\right)\mathbf{a}_1 + \left(\frac{1}{4} - x_5\right)\mathbf{a}_2 + x_5\,\mathbf{a}_3 & = & \frac{1}{8}a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_5\right)a\,\hat{\mathbf{z}} & (48f)\ \mathrm{H_2O\ III} \\
\mathbf{B}_{27} & = & \left(\frac{3}{4} + x_5\right)\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \left(\frac{3}{4} + x_5\right)\mathbf{a}_3 & = & \frac{3}{8}a\,\hat{\mathbf{x}} + \left(\frac{3}{4} + x_5\right)a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}} & (48f)\ \mathrm{H_2O\ III} \\
\mathbf{B}_{28} & = & -x_5\,\mathbf{a}_1 + \left(\frac{3}{4} + x_5\right)\mathbf{a}_2 - x_5\,\mathbf{a}_3 & = & \frac{3}{8}a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}} & (48f)\ \mathrm{H_2O\ III} \\
\mathbf{B}_{29} & = & -x_5\,\mathbf{a}_1 + \left(\frac{3}{4} + x_5\right)\mathbf{a}_2 + \left(\frac{3}{4} + x_5\right)\mathbf{a}_3 & = & \left(\frac{3}{4} + x_5\right)a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}} & (48f)\ \mathrm{H_2O\ III} \\
\mathbf{B}_{30} & = & \left(\frac{3}{4} + x_5\right)\mathbf{a}_1 - x_5\,\mathbf{a}_2 - x_5\,\mathbf{a}_3 & = & -x_5 a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}} & (48f)\ \mathrm{H_2O\ III} \\
\mathbf{B}_{31} & = & -x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \left(\frac{3}{4} + x_5\right)\mathbf{a}_3 & = & \frac{3}{8}a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}} & (48f)\ \mathrm{H_2O\ III} \\
\mathbf{B}_{32} & = & \left(\frac{3}{4} + x_5\right)\mathbf{a}_1 + \left(\frac{3}{4} + x_5\right)\mathbf{a}_2 - x_5\,\mathbf{a}_3 & = & \frac{3}{8}a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} + \left(\frac{3}{4} + x_5\right)a\,\hat{\mathbf{z}} & (48f)\ \mathrm{H_2O\ III} \\
\mathbf{B}_{33} & = & \left(\frac{1}{4} - x_6\right)\mathbf{a}_1 + x_6\,\mathbf{a}_2 + x_6\,\mathbf{a}_3 & = & x_6 a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}} & (48f)\ \mathrm{H_2O\ IV} \\
\mathbf{B}_{34} & = & x_6\,\mathbf{a}_1 + \left(\frac{1}{4} - x_6\right)\mathbf{a}_2 + \left(\frac{1}{4} - x_6\right)\mathbf{a}_3 & = & \left(\frac{1}{4} - x_6\right)a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}} & (48f)\ \mathrm{H_2O\ IV} \\
\mathbf{B}_{35} & = & x_6\,\mathbf{a}_1 + \left(\frac{1}{4} - x_6\right)\mathbf{a}_2 + x_6\,\mathbf{a}_3 & = & \frac{1}{8}a\,\hat{\mathbf{x}} + x_6 a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}} & (48f)\ \mathrm{H_2O\ IV} \\
\mathbf{B}_{36} & = & \left(\frac{1}{4} - x_6\right)\mathbf{a}_1 + x_6\,\mathbf{a}_2 + \left(\frac{1}{4} - x_6\right)\mathbf{a}_3 & = & \frac{1}{8}a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_6\right)a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}} & (48f)\ \mathrm{H_2O\ IV} \\
\mathbf{B}_{37} & = & x_6\,\mathbf{a}_1 + x_6\,\mathbf{a}_2 + \left(\frac{1}{4} - x_6\right)\mathbf{a}_3 & = & \frac{1}{8}a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + x_6 a\,\hat{\mathbf{z}} & (48f)\ \mathrm{H_2O\ IV} \\
\mathbf{B}_{38} & = & \left(\frac{1}{4} - x_6\right)\mathbf{a}_1 + \left(\frac{1}{4} - x_6\right)\mathbf{a}_2 + x_6\,\mathbf{a}_3 & = & \frac{1}{8}a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_6\right)a\,\hat{\mathbf{z}} & (48f)\ \mathrm{H_2O\ IV} \\
\mathbf{B}_{39} & = & \left(\frac{3}{4} + x_6\right)\mathbf{a}_1 - x_6\,\mathbf{a}_2 + \left(\frac{3}{4} + x_6\right)\mathbf{a}_3 & = & \frac{3}{8}a\,\hat{\mathbf{x}} + \left(\frac{3}{4} + x_6\right)a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}} & (48f)\ \mathrm{H_2O\ IV} \\
\mathbf{B}_{40} & = & -x_6\,\mathbf{a}_1 + \left(\frac{3}{4} + x_6\right)\mathbf{a}_2 - x_6\,\mathbf{a}_3 & = & \frac{3}{8}a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}} & (48f)\ \mathrm{H_2O\ IV} \\
\mathbf{B}_{41} & = & -x_6\,\mathbf{a}_1 + \left(\frac{3}{4} + x_6\right)\mathbf{a}_2 + \left(\frac{3}{4} + x_6\right)\mathbf{a}_3 & = & \left(\frac{3}{4} + x_6\right)a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}} & (48f)\ \mathrm{H_2O\ IV} \\
\mathbf{B}_{42} & = & \left(\frac{3}{4} + x_6\right)\mathbf{a}_1 - x_6\,\mathbf{a}_2 - x_6\,\mathbf{a}_3 & = & -x_6 a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}} & (48f)\ \mathrm{H_2O\ IV} \\
\mathbf{B}_{43} & = & -x_6\,\mathbf{a}_1 - x_6\,\mathbf{a}_2 + \left(\frac{3}{4} + x_6\right)\mathbf{a}_3 & = & \frac{3}{8}a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} - x_6 a\,\hat{\mathbf{z}} & (48f)\ \mathrm{H_2O\ IV} \\
\mathbf{B}_{44} & = & \left(\frac{3}{4} + x_6\right)\mathbf{a}_1 + \left(\frac{3}{4} + x_6\right)\mathbf{a}_2 - x_6\,\mathbf{a}_3 & = & \frac{3}{8}a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} + \left(\frac{3}{4} + x_6\right)a\,\hat{\mathbf{z}} & (48f)\ \mathrm{H_2O\ IV} \\
\mathbf{B}_{45} & = & z_7\,\mathbf{a}_1 + z_7\,\mathbf{a}_2 + (2x_7 - z_7)\,\mathbf{a}_3 & = & x_7 a\,\hat{\mathbf{x}} + x_7 a\,\hat{\mathbf{y}} + z_7 a\,\hat{\mathbf{z}} & (96g)\ \mathrm{H_2O\ V} \\
\mathbf{B}_{46} & = & z_7\,\mathbf{a}_1 + z_7\,\mathbf{a}_2 + \left(\frac{1}{2} - 2x_7 - z_7\right)\mathbf{a}_3 & = & \left(\frac{1}{4} - x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_7\right)a\,\hat{\mathbf{y}} + z_7 a\,\hat{\mathbf{z}} & (96g)\ \mathrm{H_2O\ V} \\
\mathbf{B}_{47} & = & (2x_7 - z_7)\,\mathbf{a}_1 + \left(\frac{1}{2} - 2x_7 - z_7\right)\mathbf{a}_2 + z_7\,\mathbf{a}_3 & = & \left(\frac{1}{4} - x_7\right)a\,\hat{\mathbf{x}} + x_7 a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_7\right)a\,\hat{\mathbf{z}} & (96g)\ \mathrm{H_2O\ V} \\
\mathbf{B}_{48} & = & \left(\frac{1}{2} - 2x_7 - z_7\right)\mathbf{a}_1 + (2x_7 - z_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 & = & x_7 a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_7\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_7\right)a\,\hat{\mathbf{z}} & (96g)\ \mathrm{H_2O\ V} \\
\mathbf{B}_{49} & = & (2x_7 - z_7)\,\mathbf{a}_1 + z_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 & = & z_7 a\,\hat{\mathbf{x}} + x_7 a\,\hat{\mathbf{y}} + x_7 a\,\hat{\mathbf{z}} & (96g)\ \mathrm{H_2O\ V} \\
\mathbf{B}_{50} & = & \left(\frac{1}{2} - 2x_7 - z_7\right)\mathbf{a}_1 + z_7\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 & = & z_7 a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_7\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_7\right)a\,\hat{\mathbf{z}} & (96g)\ \mathrm{H_2O\ V} \\
\mathbf{B}_{51} & = & z_7\,\mathbf{a}_1 + (2x_7 - z_7)\,\mathbf{a}_2 + \left(\frac{1}{2} - 2x_7 - z_7\right)\mathbf{a}_3 & = & \left(\frac{1}{4} - z_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_7\right)a\,\hat{\mathbf{y}} + x_7 a\,\hat{\mathbf{z}} & (96g)\ \mathrm{H_2O\ V} \\
\mathbf{B}_{52} & = & z_7\,\mathbf{a}_1 + \left(\frac{1}{2} - 2x_7 - z_7\right)\mathbf{a}_2 + (2x_7 - z_7)\,\mathbf{a}_3 & = & \left(\frac{1}{4} - z_7\right)a\,\hat{\mathbf{x}} + x_7 a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_7\right)a\,\hat{\mathbf{z}} & (96g)\ \mathrm{H_2O\ V} \\
\mathbf{B}_{53} & = & z_7\,\mathbf{a}_1 + (2x_7 - z_7)\,\mathbf{a}_2 + z_7\,\mathbf{a}_3 & = & x_7 a\,\hat{\mathbf{x}} + z_7 a\,\hat{\mathbf{y}} + x_7 a\,\hat{\mathbf{z}} & (96g)\ \mathrm{H_2O\ V} \\
\mathbf{B}_{54} & = & z_7\,\mathbf{a}_1 + \left(\frac{1}{2} - 2x_7 - z_7\right)\mathbf{a}_2 + z_7\,\mathbf{a}_3 & = & \left(\frac{1}{4} - x_7\right)a\,\hat{\mathbf{x}} + z_7 a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_7\right)a\,\hat{\mathbf{z}} & (96g)\ \mathrm{H_2O\ V} \\
\mathbf{B}_{55} & = & \left(\frac{1}{2} - 2x_7 - z_7\right)\mathbf{a}_1 + z_7\,\mathbf{a}_2 + (2x_7 - z_7)\,\mathbf{a}_3 & = & x_7 a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - z_7\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_7\right)a\,\hat{\mathbf{z}} & (96g)\ \mathrm{H_2O\ V} \\
\mathbf{B}_{56} & = & (2x_7 - z_7)\,\mathbf{a}_1 + z_7\,\mathbf{a}_2 + \left(\frac{1}{2} - 2x_7 - z_7\right)\mathbf{a}_3 & = & \left(\frac{1}{4} - x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - z_7\right)a\,\hat{\mathbf{y}} + x_7 a\,\hat{\mathbf{z}} & (96g)\ \mathrm{H_2O\ V} \\
\mathbf{B}_{57} & = & -z_7\,\mathbf{a}_1 - z_7\,\mathbf{a}_2 + \left(\frac{1}{2} + 2x_7 + z_7\right)\mathbf{a}_3 & = & \left(\frac{1}{4} + x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_7\right)a\,\hat{\mathbf{y}} - z_7 a\,\hat{\mathbf{z}} & (96g)\ \mathrm{H_2O\ V}
\end{array}
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$\mathbf{B}_{58}$	=	$-z_7\,\mathbf{a}_1 - z_7\,\mathbf{a}_2 + (-2x_7 + z_7)\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} - x_7 a\,\hat{\mathbf{y}} - z_7 a\,\hat{\mathbf{z}}$	(96*g*)	H_2O V
$\mathbf{B}_{59}$	=	$(-2x_7 + z_7)\,\mathbf{a}_1 + \left(\frac{1}{2} + 2x_7 + z_7\right)\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_7\right)a\,\hat{\mathbf{x}} - x_7 a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_7\right)a\,\hat{\mathbf{z}}$	(96*g*)	H_2O V
$\mathbf{B}_{60}$	=	$\left(\frac{1}{2} + 2x_7 + z_7\right)\mathbf{a}_1 + (-2x_7 + z_7)\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_7\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_7\right)a\,\hat{\mathbf{z}}$	(96*g*)	H_2O V
$\mathbf{B}_{61}$	=	$(-2x_7 + z_7)\,\mathbf{a}_1 - z_7\,\mathbf{a}_2 + \left(\frac{1}{2} + 2x_7 + z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + z_7\right)a\,\hat{\mathbf{y}} - x_7 a\,\hat{\mathbf{z}}$	(96*g*)	H_2O V
$\mathbf{B}_{62}$	=	$\left(\frac{1}{2} + 2x_7 + z_7\right)\mathbf{a}_1 - z_7\,\mathbf{a}_2 + (-2x_7 + z_7)\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + z_7\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + x_7\right)a\,\hat{\mathbf{z}}$	(96*g*)	H_2O V
$\mathbf{B}_{63}$	=	$-z_7\,\mathbf{a}_1 + (-2x_7 + z_7)\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$-x_7 a\,\hat{\mathbf{x}} - z_7 a\,\hat{\mathbf{y}} - x_7 a\,\hat{\mathbf{z}}$	(96*g*)	H_2O V
$\mathbf{B}_{64}$	=	$-z_7\,\mathbf{a}_1 + \left(\frac{1}{2} + 2x_7 + z_7\right)\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_7\right)a\,\hat{\mathbf{x}} - z_7 a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + x_7\right)a\,\hat{\mathbf{z}}$	(96*g*)	H_2O V
$\mathbf{B}_{65}$	=	$-z_7\,\mathbf{a}_1 + (-2x_7 + z_7)\,\mathbf{a}_2 + \left(\frac{1}{2} + 2x_7 + z_7\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + z_7\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_7\right)a\,\hat{\mathbf{y}} - x_7 a\,\hat{\mathbf{z}}$	(96*g*)	H_2O V
$\mathbf{B}_{66}$	=	$-z_7\,\mathbf{a}_1 + \left(\frac{1}{2} + 2x_7 + z_7\right)\mathbf{a}_2 + (-2x_7 + z_7)\,\mathbf{a}_3$	=	$\left(\frac{1}{4} + z_7\right)a\,\hat{\mathbf{x}} - x_7 a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + x_7\right)a\,\hat{\mathbf{z}}$	(96*g*)	H_2O V
$\mathbf{B}_{67}$	=	$\left(\frac{1}{2} + 2x_7 + z_7\right)\mathbf{a}_1 - z_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$-z_7 a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_7\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + x_7\right)a\,\hat{\mathbf{z}}$	(96*g*)	H_2O V
$\mathbf{B}_{68}$	=	$(-2x_7 + z_7)\,\mathbf{a}_1 - z_7\,\mathbf{a}_2 - z_7\,\mathbf{a}_3$	=	$-z_7 a\,\hat{\mathbf{x}} - x_7 a\,\hat{\mathbf{y}} - x_7 a\,\hat{\mathbf{z}}$	(96*g*)	H_2O V
$\mathbf{B}_{69}$	=	$z_8\,\mathbf{a}_1 + z_8\,\mathbf{a}_2 + (2x_8 - z_8)\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + x_8 a\,\hat{\mathbf{y}} + z_8 a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{70}$	=	$z_8\,\mathbf{a}_1 + z_8\,\mathbf{a}_2 + \left(\frac{1}{2} - 2x_8 - z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_8\right)a\,\hat{\mathbf{y}} + z_8 a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{71}$	=	$(2x_8 - z_8)\,\mathbf{a}_1 + \left(\frac{1}{2} - 2x_8 - z_8\right)\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_8\right)a\,\hat{\mathbf{x}} + x_8 a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_8\right)a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{72}$	=	$\left(\frac{1}{2} - 2x_8 - z_8\right)\mathbf{a}_1 + (2x_8 - z_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_8\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_8\right)a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{73}$	=	$(2x_8 - z_8)\,\mathbf{a}_1 + z_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$z_8 a\,\hat{\mathbf{x}} + x_8 a\,\hat{\mathbf{y}} + x_8 a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{74}$	=	$\left(\frac{1}{2} - 2x_8 - z_8\right)\mathbf{a}_1 + z_8\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$z_8 a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_8\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_8\right)a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{75}$	=	$z_8\,\mathbf{a}_1 + (2x_8 - z_8)\,\mathbf{a}_2 + \left(\frac{1}{2} - 2x_8 - z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - z_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_8\right)a\,\hat{\mathbf{y}} + x_8 a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{76}$	=	$z_8\,\mathbf{a}_1 + \left(\frac{1}{2} - 2x_8 - z_8\right)\mathbf{a}_2 + (2x_8 - z_8)\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - z_8\right)a\,\hat{\mathbf{x}} + x_8 a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_8\right)a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{77}$	=	$z_8\,\mathbf{a}_1 + (2x_8 - z_8)\,\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + z_8 a\,\hat{\mathbf{y}} + x_8 a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{78}$	=	$z_8\,\mathbf{a}_1 + \left(\frac{1}{2} - 2x_8 - z_8\right)\mathbf{a}_2 + z_8\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_8\right)a\,\hat{\mathbf{x}} + z_8 a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_8\right)a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{79}$	=	$\left(\frac{1}{2} - 2x_8 - z_8\right)\mathbf{a}_1 + z_8\,\mathbf{a}_2 + (2x_8 - z_8)\,\mathbf{a}_3$	=	$x_8 a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - z_8\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_8\right)a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{80}$	=	$(2x_8 - z_8)\,\mathbf{a}_1 + z_8\,\mathbf{a}_2 + \left(\frac{1}{2} - 2x_8 - z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - z_8\right)a\,\hat{\mathbf{y}} + x_8 a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{81}$	=	$-z_8\,\mathbf{a}_1 - z_8\,\mathbf{a}_2 + \left(\frac{1}{2} + 2x_8 + z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_8\right)a\,\hat{\mathbf{y}} - z_8 a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{82}$	=	$-z_8\,\mathbf{a}_1 - z_8\,\mathbf{a}_2 + (-2x_8 + z_8)\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} - x_8 a\,\hat{\mathbf{y}} - z_8 a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{83}$	=	$(-2x_8 + z_8)\,\mathbf{a}_1 + \left(\frac{1}{2} + 2x_8 + z_8\right)\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_8\right)a\,\hat{\mathbf{x}} - x_8 a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_8\right)a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{84}$	=	$\left(\frac{1}{2} + 2x_8 + z_8\right)\mathbf{a}_1 + (-2x_8 + z_8)\,\mathbf{a}_2 - z_8\,\mathbf{a}_3$	=	$-x_8 a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_8\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + z_8\right)a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{85}$	=	$(-2x_8 + z_8)\,\mathbf{a}_1 - z_8\,\mathbf{a}_2 + \left(\frac{1}{2} + 2x_8 + z_8\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_8\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + z_8\right)a\,\hat{\mathbf{y}} - x_8 a\,\hat{\mathbf{z}}$	(96*g*)	O II

$$\mathbf{B}_{86} = \left(\tfrac{1}{2} + 2x_8 + z_8\right)\mathbf{a}_1 - z_8\,\mathbf{a}_2 + (-2x_8 + z_8)\,\mathbf{a}_3 = -x_8 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + z_8\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + x_8\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O II}$$

$$\mathbf{B}_{87} = -z_8\,\mathbf{a}_1 + (-2x_8 + z_8)\,\mathbf{a}_2 - z_8\,\mathbf{a}_3 = -x_8 a\,\hat{\mathbf{x}} - z_8 a\,\hat{\mathbf{y}} - x_8 a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O II}$$

$$\mathbf{B}_{88} = -z_8\,\mathbf{a}_1 + \left(\tfrac{1}{2} + 2x_8 + z_8\right)\mathbf{a}_2 - z_8\,\mathbf{a}_3 = \left(\tfrac{1}{4} + x_8\right)a\,\hat{\mathbf{x}} - z_8 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + x_8\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O II}$$

$$\mathbf{B}_{89} = -z_8\,\mathbf{a}_1 + (-2x_8 + z_8)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + 2x_8 + z_8\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + z_8\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_8\right)a\,\hat{\mathbf{y}} - x_8 a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O II}$$

$$\mathbf{B}_{90} = -z_8\,\mathbf{a}_1 + \left(\tfrac{1}{2} + 2x_8 + z_8\right)\mathbf{a}_2 + (-2x_8 + z_8)\,\mathbf{a}_3 = \left(\tfrac{1}{4} + z_8\right)a\,\hat{\mathbf{x}} - x_8 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + x_8\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O II}$$

$$\mathbf{B}_{91} = \left(\tfrac{1}{2} + 2x_8 + z_8\right)\mathbf{a}_1 - z_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3 = -z_8 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_8\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + x_8\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O II}$$

$$\mathbf{B}_{92} = (-2x_8 + z_8)\,\mathbf{a}_1 - z_8\,\mathbf{a}_2 - z_8\,\mathbf{a}_3 = -z_8 a\,\hat{\mathbf{x}} - x_8 a\,\hat{\mathbf{y}} - x_8 a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O II}$$

$$\mathbf{B}_{93} = z_9\,\mathbf{a}_1 + z_9\,\mathbf{a}_2 + (2x_9 - z_9)\,\mathbf{a}_3 = x_9 a\,\hat{\mathbf{x}} + x_9 a\,\hat{\mathbf{y}} + z_9 a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O III}$$

$$\mathbf{B}_{94} = z_9\,\mathbf{a}_1 + z_9\,\mathbf{a}_2 + \left(\tfrac{1}{2} - 2x_9 - z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - x_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_9\right)a\,\hat{\mathbf{y}} + z_9 a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O III}$$

$$\mathbf{B}_{95} = (2x_9 - z_9)\,\mathbf{a}_1 + \left(\tfrac{1}{2} - 2x_9 - z_9\right)\mathbf{a}_2 + z_9\,\mathbf{a}_3 = \left(\tfrac{1}{4} - x_9\right)a\,\hat{\mathbf{x}} + x_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - z_9\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O III}$$

$$\mathbf{B}_{96} = \left(\tfrac{1}{2} - 2x_9 - z_9\right)\mathbf{a}_1 + (2x_9 - z_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 = x_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - z_9\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O III}$$

$$\mathbf{B}_{97} = (2x_9 - z_9)\,\mathbf{a}_1 + z_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 = z_9 a\,\hat{\mathbf{x}} + x_9 a\,\hat{\mathbf{y}} + x_9 a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O III}$$

$$\mathbf{B}_{98} = \left(\tfrac{1}{2} - 2x_9 - z_9\right)\mathbf{a}_1 + z_9\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 = z_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - x_9\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O III}$$

$$\mathbf{B}_{99} = z_9\,\mathbf{a}_1 + (2x_9 - z_9)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - 2x_9 - z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - z_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_9\right)a\,\hat{\mathbf{y}} + x_9 a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O III}$$

$$\mathbf{B}_{100} = z_9\,\mathbf{a}_1 + \left(\tfrac{1}{2} - 2x_9 - z_9\right)\mathbf{a}_2 + (2x_9 - z_9)\,\mathbf{a}_3 = \left(\tfrac{1}{4} - z_9\right)a\,\hat{\mathbf{x}} + x_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - x_9\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O III}$$

$$\mathbf{B}_{101} = z_9\,\mathbf{a}_1 + (2x_9 - z_9)\,\mathbf{a}_2 + z_9\,\mathbf{a}_3 = x_9 a\,\hat{\mathbf{x}} + z_9 a\,\hat{\mathbf{y}} + x_9 a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O III}$$

$$\mathbf{B}_{102} = z_9\,\mathbf{a}_1 + \left(\tfrac{1}{2} - 2x_9 - z_9\right)\mathbf{a}_2 + z_9\,\mathbf{a}_3 = \left(\tfrac{1}{4} - x_9\right)a\,\hat{\mathbf{x}} + z_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - x_9\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O III}$$

$$\mathbf{B}_{103} = \left(\tfrac{1}{2} - 2x_9 - z_9\right)\mathbf{a}_1 + z_9\,\mathbf{a}_2 + (2x_9 - z_9)\,\mathbf{a}_3 = x_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - z_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - x_9\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O III}$$

$$\mathbf{B}_{104} = (2x_9 - z_9)\,\mathbf{a}_1 + z_9\,\mathbf{a}_2 + \left(\tfrac{1}{2} - 2x_9 - z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - x_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - z_9\right)a\,\hat{\mathbf{y}} + x_9 a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O III}$$

$$\mathbf{B}_{105} = -z_9\,\mathbf{a}_1 - z_9\,\mathbf{a}_2 + \left(\tfrac{1}{2} + 2x_9 + z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + x_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_9\right)a\,\hat{\mathbf{y}} - z_9 a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O III}$$

$$\mathbf{B}_{106} = -z_9\,\mathbf{a}_1 - z_9\,\mathbf{a}_2 + (-2x_9 + z_9)\,\mathbf{a}_3 = -x_9 a\,\hat{\mathbf{x}} - x_9 a\,\hat{\mathbf{y}} - z_9 a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O III}$$

$$\mathbf{B}_{107} = (-2x_9 + z_9)\,\mathbf{a}_1 + \left(\tfrac{1}{2} + 2x_9 + z_9\right)\mathbf{a}_2 - z_9\,\mathbf{a}_3 = \left(\tfrac{1}{4} + x_9\right)a\,\hat{\mathbf{x}} - x_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_9\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O III}$$

$$\mathbf{B}_{108} = \left(\tfrac{1}{2} + 2x_9 + z_9\right)\mathbf{a}_1 + (-2x_9 + z_9)\,\mathbf{a}_2 - z_9\,\mathbf{a}_3 = -x_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_9\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O III}$$

$$\mathbf{B}_{109} = (-2x_9 + z_9)\,\mathbf{a}_1 - z_9\,\mathbf{a}_2 + \left(\tfrac{1}{2} + 2x_9 + z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + x_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + z_9\right)a\,\hat{\mathbf{y}} - x_9 a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O III}$$

$$\mathbf{B}_{110} = \left(\tfrac{1}{2} + 2x_9 + z_9\right)\mathbf{a}_1 - z_9\,\mathbf{a}_2 + (-2x_9 + z_9)\,\mathbf{a}_3 = -x_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + z_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + x_9\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O III}$$

$$\mathbf{B}_{111} = -z_9\,\mathbf{a}_1 + (-2x_9 + z_9)\,\mathbf{a}_2 - z_9\,\mathbf{a}_3 = -x_9 a\,\hat{\mathbf{x}} - z_9 a\,\hat{\mathbf{y}} - x_9 a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O III}$$

$$\mathbf{B}_{112} = -z_9\,\mathbf{a}_1 + \left(\tfrac{1}{2} + 2x_9 + z_9\right)\mathbf{a}_2 - z_9\,\mathbf{a}_3 = \left(\tfrac{1}{4} + x_9\right)a\,\hat{\mathbf{x}} - z_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + x_9\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O III}$$

$$\mathbf{B}_{113} = -z_9\,\mathbf{a}_1 + (-2x_9 + z_9)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + 2x_9 + z_9\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + z_9\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_9\right)a\,\hat{\mathbf{y}} - x_9 a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O III}$$

$$\mathbf{B}_{114} = -z_9\,\mathbf{a}_1 + \left(\tfrac{1}{2} + 2x_9 + z_9\right)\mathbf{a}_2 + (-2x_9 + z_9)\,\mathbf{a}_3 = \left(\tfrac{1}{4} + z_9\right)a\,\hat{\mathbf{x}} - x_9 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + x_9\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O III}$$

$$\mathbf{B}_{115} = \left(\tfrac{1}{2} + 2x_9 + z_9\right)\mathbf{a}_1 - z_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3 = -z_9 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_9\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + x_9\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O III}$$

$$\mathbf{B}_{116} = (-2x_9 + z_9)\,\mathbf{a}_1 - z_9\,\mathbf{a}_2 - z_9\,\mathbf{a}_3 = -z_9 a\,\hat{\mathbf{x}} - x_9 a\,\hat{\mathbf{y}} - x_9 a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O III}$$

$$\mathbf{B}_{117} = z_{10}\,\mathbf{a}_1 + z_{10}\,\mathbf{a}_2 + (2x_{10} - z_{10})\,\mathbf{a}_3 = x_{10} a\,\hat{\mathbf{x}} + x_{10} a\,\hat{\mathbf{y}} + z_{10} a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O IV}$$

$$\mathbf{B}_{118} = z_{10}\,\mathbf{a}_1 + z_{10}\,\mathbf{a}_2 + \left(\tfrac{1}{2} - 2x_{10} - z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - x_{10}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_{10}\right)a\,\hat{\mathbf{y}} + z_{10} a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O IV}$$

$$\mathbf{B}_{119} = (2x_{10} - z_{10})\,\mathbf{a}_1 + \left(\tfrac{1}{2} - 2x_{10} - z_{10}\right)\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 = \left(\tfrac{1}{4} - x_{10}\right)a\,\hat{\mathbf{x}} + x_{10} a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - z_{10}\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O IV}$$

$$\mathbf{B}_{120} = \left(\tfrac{1}{2} - 2x_{10} - z_{10}\right)\mathbf{a}_1 + (2x_{10} - z_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 = x_{10} a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_{10}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - z_{10}\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O IV}$$

$$\mathbf{B}_{121} = (2x_{10} - z_{10})\,\mathbf{a}_1 + z_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 = z_{10} a\,\hat{\mathbf{x}} + x_{10} a\,\hat{\mathbf{y}} + x_{10} a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O IV}$$

$$\mathbf{B}_{122} = \left(\tfrac{1}{2} - 2x_{10} - z_{10}\right)\mathbf{a}_1 + z_{10}\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 = z_{10} a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_{10}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - x_{10}\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O IV}$$

$$\mathbf{B}_{123} = z_{10}\,\mathbf{a}_1 + (2x_{10} - z_{10})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - 2x_{10} - z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - z_{10}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_{10}\right)a\,\hat{\mathbf{y}} + x_{10} a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O IV}$$

$$\mathbf{B}_{124} = z_{10}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - 2x_{10} - z_{10}\right)\mathbf{a}_2 + (2x_{10} - z_{10})\,\mathbf{a}_3 = \left(\tfrac{1}{4} - z_{10}\right)a\,\hat{\mathbf{x}} + x_{10} a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - x_{10}\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O IV}$$

$$\mathbf{B}_{125} = z_{10}\,\mathbf{a}_1 + (2x_{10} - z_{10})\,\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 = x_{10} a\,\hat{\mathbf{x}} + z_{10} a\,\hat{\mathbf{y}} + x_{10} a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O IV}$$

$$\mathbf{B}_{126} = z_{10}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - 2x_{10} - z_{10}\right)\mathbf{a}_2 + z_{10}\,\mathbf{a}_3 = \left(\tfrac{1}{4} - x_{10}\right)a\,\hat{\mathbf{x}} + z_{10} a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - x_{10}\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O IV}$$

$$\mathbf{B}_{127} = \left(\tfrac{1}{2} - 2x_{10} - z_{10}\right)\mathbf{a}_1 + z_{10}\,\mathbf{a}_2 + (2x_{10} - z_{10})\,\mathbf{a}_3 = x_{10} a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - z_{10}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - x_{10}\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O IV}$$

$$\mathbf{B}_{128} = (2x_{10} - z_{10})\,\mathbf{a}_1 + z_{10}\,\mathbf{a}_2 + \left(\tfrac{1}{2} - 2x_{10} - z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - x_{10}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - z_{10}\right)a\,\hat{\mathbf{y}} + x_{10} a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O IV}$$

$$\mathbf{B}_{129} = -z_{10}\,\mathbf{a}_1 - z_{10}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + 2x_{10} + z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + x_{10}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_{10}\right)a\,\hat{\mathbf{y}} - z_{10} a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O IV}$$

$$\mathbf{B}_{130} = -z_{10}\,\mathbf{a}_1 - z_{10}\,\mathbf{a}_2 + (-2x_{10} + z_{10})\,\mathbf{a}_3 = -x_{10} a\,\hat{\mathbf{x}} - x_{10} a\,\hat{\mathbf{y}} - z_{10} a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O IV}$$

$$\mathbf{B}_{131} = (-2x_{10} + z_{10})\,\mathbf{a}_1 + \left(\tfrac{1}{2} + 2x_{10} + z_{10}\right)\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 = \left(\tfrac{1}{4} + x_{10}\right)a\,\hat{\mathbf{x}} - x_{10} a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_{10}\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O IV}$$

$$\mathbf{B}_{132} = \left(\tfrac{1}{2} + 2x_{10} + z_{10}\right)\mathbf{a}_1 + (-2x_{10} + z_{10})\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 = -x_{10} a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_{10}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_{10}\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O IV}$$

$$\mathbf{B}_{133} = (-2x_{10} + z_{10})\,\mathbf{a}_1 - z_{10}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + 2x_{10} + z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + x_{10}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + z_{10}\right)a\,\hat{\mathbf{y}} - x_{10} a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O IV}$$

$$\mathbf{B}_{134} = \left(\tfrac{1}{2} + 2x_{10} + z_{10}\right)\mathbf{a}_1 - z_{10}\,\mathbf{a}_2 + (-2x_{10} + z_{10})\,\mathbf{a}_3 = -x_{10} a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + z_{10}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + x_{10}\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O IV}$$

$$\mathbf{B}_{135} = -z_{10}\,\mathbf{a}_1 + (-2x_{10} + z_{10})\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 = -x_{10} a\,\hat{\mathbf{x}} - z_{10} a\,\hat{\mathbf{y}} - x_{10} a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O IV}$$

$$\mathbf{B}_{136} = -z_{10}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + 2x_{10} + z_{10}\right)\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 = \left(\tfrac{1}{4} + x_{10}\right)a\,\hat{\mathbf{x}} - z_{10} a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + x_{10}\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O IV}$$

$$\mathbf{B}_{137} = -z_{10}\,\mathbf{a}_1 + (-2x_{10} + z_{10})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + 2x_{10} + z_{10}\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + z_{10}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_{10}\right)a\,\hat{\mathbf{y}} - x_{10} a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O IV}$$

$$\mathbf{B}_{138} = -z_{10}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + 2x_{10} + z_{10}\right)\mathbf{a}_2 + (-2x_{10} + z_{10})\,\mathbf{a}_3 = \left(\tfrac{1}{4} + z_{10}\right)a\,\hat{\mathbf{x}} - x_{10} a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + x_{10}\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O IV}$$

$$\mathbf{B}_{139} = \left(\tfrac{1}{2} + 2x_{10} + z_{10}\right)\mathbf{a}_1 - z_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 = -z_{10} a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_{10}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + x_{10}\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O IV}$$

$$\mathbf{B}_{140} = (-2x_{10} + z_{10})\,\mathbf{a}_1 - z_{10}\,\mathbf{a}_2 - z_{10}\,\mathbf{a}_3 = -z_{10} a\,\hat{\mathbf{x}} - x_{10} a\,\hat{\mathbf{y}} - x_{10} a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O IV}$$

$$\mathbf{B}_{141} = z_{11}\,\mathbf{a}_1 + z_{11}\,\mathbf{a}_2 + (2x_{11} - z_{11})\,\mathbf{a}_3 = x_{11}a\,\hat{\mathbf{x}} + x_{11}a\,\hat{\mathbf{y}} + z_{11}a\,\hat{\mathbf{z}} \quad (96g) \quad \mathrm{W}$$

$$\mathbf{B}_{142} = z_{11}\,\mathbf{a}_1 + z_{11}\,\mathbf{a}_2 + \left(\tfrac{1}{2} - 2x_{11} - z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - x_{11}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_{11}\right)a\,\hat{\mathbf{y}} + z_{11}a\,\hat{\mathbf{z}} \quad (96g) \quad \mathrm{W}$$

$$\mathbf{B}_{143} = (2x_{11} - z_{11})\,\mathbf{a}_1 + \left(\tfrac{1}{2} - 2x_{11} - z_{11}\right)\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 = \left(\tfrac{1}{4} - x_{11}\right)a\,\hat{\mathbf{x}} + x_{11}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - z_{11}\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \mathrm{W}$$

$$\mathbf{B}_{144} = \left(\tfrac{1}{2} - 2x_{11} - z_{11}\right)\mathbf{a}_1 + (2x_{11} - z_{11})\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 = x_{11}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_{11}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - z_{11}\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \mathrm{W}$$

$$\mathbf{B}_{145} = (2x_{11} - z_{11})\,\mathbf{a}_1 + z_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 = z_{11}a\,\hat{\mathbf{x}} + x_{11}a\,\hat{\mathbf{y}} + x_{11}a\,\hat{\mathbf{z}} \quad (96g) \quad \mathrm{W}$$

$$\mathbf{B}_{146} = \left(\tfrac{1}{2} - 2x_{11} - z_{11}\right)\mathbf{a}_1 + z_{11}\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 = z_{11}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_{11}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - x_{11}\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \mathrm{W}$$

$$\mathbf{B}_{147} = z_{11}\,\mathbf{a}_1 + (2x_{11} - z_{11})\,\mathbf{a}_2 + \left(\tfrac{1}{2} - 2x_{11} - z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - z_{11}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_{11}\right)a\,\hat{\mathbf{y}} + x_{11}a\,\hat{\mathbf{z}} \quad (96g) \quad \mathrm{W}$$

$$\mathbf{B}_{148} = z_{11}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - 2x_{11} - z_{11}\right)\mathbf{a}_2 + (2x_{11} - z_{11})\,\mathbf{a}_3 = \left(\tfrac{1}{4} - z_{11}\right)a\,\hat{\mathbf{x}} + x_{11}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - x_{11}\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \mathrm{W}$$

$$\mathbf{B}_{149} = z_{11}\,\mathbf{a}_1 + (2x_{11} - z_{11})\,\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 = x_{11}a\,\hat{\mathbf{x}} + z_{11}a\,\hat{\mathbf{y}} + x_{11}a\,\hat{\mathbf{z}} \quad (96g) \quad \mathrm{W}$$

$$\mathbf{B}_{150} = z_{11}\,\mathbf{a}_1 + \left(\tfrac{1}{2} - 2x_{11} - z_{11}\right)\mathbf{a}_2 + z_{11}\,\mathbf{a}_3 = \left(\tfrac{1}{4} - x_{11}\right)a\,\hat{\mathbf{x}} + z_{11}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - x_{11}\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \mathrm{W}$$

$$\mathbf{B}_{151} = \left(\tfrac{1}{2} - 2x_{11} - z_{11}\right)\mathbf{a}_1 + z_{11}\,\mathbf{a}_2 + (2x_{11} - z_{11})\,\mathbf{a}_3 = x_{11}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - z_{11}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - x_{11}\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \mathrm{W}$$

$$\mathbf{B}_{152} = (2x_{11} - z_{11})\,\mathbf{a}_1 + z_{11}\,\mathbf{a}_2 + \left(\tfrac{1}{2} - 2x_{11} - z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - x_{11}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - z_{11}\right)a\,\hat{\mathbf{y}} + x_{11}a\,\hat{\mathbf{z}} \quad (96g) \quad \mathrm{W}$$

$$\mathbf{B}_{153} = -z_{11}\,\mathbf{a}_1 - z_{11}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + 2x_{11} + z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + x_{11}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_{11}\right)a\,\hat{\mathbf{y}} - z_{11}a\,\hat{\mathbf{z}} \quad (96g) \quad \mathrm{W}$$

$$\mathbf{B}_{154} = -z_{11}\,\mathbf{a}_1 - z_{11}\,\mathbf{a}_2 + (-2x_{11} + z_{11})\,\mathbf{a}_3 = -x_{11}a\,\hat{\mathbf{x}} - x_{11}a\,\hat{\mathbf{y}} - z_{11}a\,\hat{\mathbf{z}} \quad (96g) \quad \mathrm{W}$$

$$\mathbf{B}_{155} = (-2x_{11} + z_{11})\,\mathbf{a}_1 + \left(\tfrac{1}{2} + 2x_{11} + z_{11}\right)\mathbf{a}_2 - z_{11}\,\mathbf{a}_3 = \left(\tfrac{1}{4} + x_{11}\right)a\,\hat{\mathbf{x}} - x_{11}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_{11}\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \mathrm{W}$$

$$\mathbf{B}_{156} = \left(\tfrac{1}{2} + 2x_{11} + z_{11}\right)\mathbf{a}_1 + (-2x_{11} + z_{11})\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3 = -x_{11}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_{11}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_{11}\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \mathrm{W}$$

$$\mathbf{B}_{157} = (-2x_{11} + z_{11})\,\mathbf{a}_1 - z_{11}\,\mathbf{a}_2 + \left(\tfrac{1}{2} + 2x_{11} + z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + x_{11}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + z_{11}\right)a\,\hat{\mathbf{y}} - x_{11}a\,\hat{\mathbf{z}} \quad (96g) \quad \mathrm{W}$$

$$\mathbf{B}_{158} = \left(\tfrac{1}{2} + 2x_{11} + z_{11}\right)\mathbf{a}_1 - z_{11}\,\mathbf{a}_2 + (-2x_{11} + z_{11})\,\mathbf{a}_3 = -x_{11}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + z_{11}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + x_{11}\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \mathrm{W}$$

$$\mathbf{B}_{159} = -z_{11}\,\mathbf{a}_1 + (-2x_{11} + z_{11})\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3 = -x_{11}a\,\hat{\mathbf{x}} - z_{11}a\,\hat{\mathbf{y}} - x_{11}a\,\hat{\mathbf{z}} \quad (96g) \quad \mathrm{W}$$

$$\mathbf{B}_{160} = -z_{11}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + 2x_{11} + z_{11}\right)\mathbf{a}_2 - z_{11}\,\mathbf{a}_3 = \left(\tfrac{1}{4} + x_{11}\right)a\,\hat{\mathbf{x}} - z_{11}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + x_{11}\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \mathrm{W}$$

$$\mathbf{B}_{161} = -z_{11}\,\mathbf{a}_1 + (-2x_{11} + z_{11})\,\mathbf{a}_2 + \left(\tfrac{1}{2} + 2x_{11} + z_{11}\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + z_{11}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_{11}\right)a\,\hat{\mathbf{y}} - x_{11}a\,\hat{\mathbf{z}} \quad (96g) \quad \mathrm{W}$$

$$\mathbf{B}_{162} = -z_{11}\,\mathbf{a}_1 + \left(\tfrac{1}{2} + 2x_{11} + z_{11}\right)\mathbf{a}_2 + (-2x_{11} + z_{11})\,\mathbf{a}_3 = \left(\tfrac{1}{4} + z_{11}\right)a\,\hat{\mathbf{x}} - x_{11}a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + x_{11}\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \mathrm{W}$$

$$\mathbf{B}_{163} = \left(\tfrac{1}{2} + 2x_{11} + z_{11}\right)\mathbf{a}_1 - z_{11}\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3 = -z_{11}a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_{11}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + x_{11}\right)a\,\hat{\mathbf{z}} \quad (96g) \quad \mathrm{W}$$

$$\mathbf{B}_{164} = (-2x_{11} + z_{11})\,\mathbf{a}_1 - z_{11}\,\mathbf{a}_2 - z_{11}\,\mathbf{a}_3 = -z_{11}a\,\hat{\mathbf{x}} - x_{11}a\,\hat{\mathbf{y}} - x_{11}a\,\hat{\mathbf{z}} \quad (96g) \quad \mathrm{W}$$

References:

- A. J. Bradley and J. W. Illingworth, *The Crystal Structure of $H_3PW_{12}O_{40}{\cdot}29H_2O$*, Proc. Roy. Soc. Lond. A **157**, 113–131 (1936), doi:10.1098/rspa.1936.0183.

- C. J. Clark and D. Hall, *Dodecamolybdophosphoric acid* circa *30-hydrate*, Acta Crystallogr. Sect. B Struct. Sci. **32**, 1545–1547 (1976), doi:10.1107/S0567740876005748.

Found in:

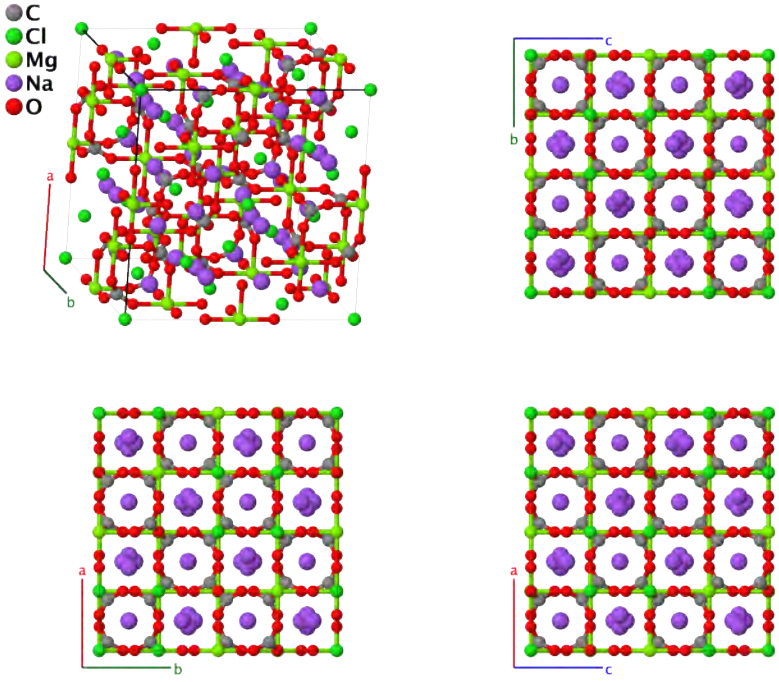
- C. Gottfried, ed., *Strukturbericht Band IV 1936* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1938).

Geometry files:

- CIF: pp. 1826
- POSCAR: pp. 1827

$G7_3$ [Northupite, $Na_3MgCl(CO_3)_2$] (*obsolete*) Structure: A2BCD3E6_cF208_227_e_c_d_f_g

http://aflow.org/prototype-encyclopedia/A2BCD3E6_cF208_227_e_c_d_f_g

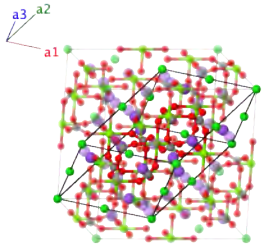


Prototype	:	$C_2ClMgNa_3O_6$
AFLOW prototype label	:	A2BCD3E6_cF208_227_e_c_d_f_g
***Strukturbericht* designation**	:	$G7_3$
Pearson symbol	:	cF208
Space group number	:	227
Space group symbol	:	$Fd\bar{3}m$
AFLOW prototype command	:	`aflow --proto=A2BCD3E6_cF208_227_e_c_d_f_g --params=`a, x_3, x_4, x_5, z_5

- This is the original structure determined by (Shiba, 1931) and given the designation $G7_3$ in (Hermann, 1937). (Negro, 1975) showed that the correct structure was actually related to cubic pyrochlore, however the two structures are very similar, and a displacement of the oxygen atoms by less than 1 Å brings the two structures into agreement.

Face-centered Cubic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(16*c*)	Cl
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}}$	(16*c*)	Cl
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(16*c*)	Cl
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$\frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(16*c*)	Cl
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(16*d*)	Mg
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(16*d*)	Mg
$\mathbf{B}_7$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(16*d*)	Mg
$\mathbf{B}_8$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(16*d*)	Mg
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(32*e*)	C
$\mathbf{B}_{10}$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \left(\frac{1}{2} - 3x_3\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	(32*e*)	C
$\mathbf{B}_{11}$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - 3x_3\right)\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{z}}$	(32*e*)	C
$\mathbf{B}_{12}$	=	$\left(\frac{1}{2} - 3x_3\right)\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{z}}$	(32*e*)	C
$\mathbf{B}_{13}$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\frac{1}{2} + 3x_3\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(32*e*)	C
$\mathbf{B}_{14}$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	(32*e*)	C
$\mathbf{B}_{15}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + 3x_3\right)\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{z}}$	(32*e*)	C
$\mathbf{B}_{16}$	=	$\left(\frac{1}{2} + 3x_3\right)\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{z}}$	(32*e*)	C
$\mathbf{B}_{17}$	=	$\left(\frac{1}{4} - x_4\right)\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Na
$\mathbf{B}_{18}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{4} - x_4\right)\,\mathbf{a}_2 + \left(\frac{1}{4} - x_4\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Na
$\mathbf{B}_{19}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{4} - x_4\right)\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Na
$\mathbf{B}_{20}$	=	$\left(\frac{1}{4} - x_4\right)\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \left(\frac{1}{4} - x_4\right)\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Na
$\mathbf{B}_{21}$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \left(\frac{1}{4} - x_4\right)\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	(48*f*)	Na
$\mathbf{B}_{22}$	=	$\left(\frac{1}{4} - x_4\right)\,\mathbf{a}_1 + \left(\frac{1}{4} - x_4\right)\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{z}}$	(48*f*)	Na
$\mathbf{B}_{23}$	=	$\left(\frac{3}{4} + x_4\right)\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \left(\frac{3}{4} + x_4\right)\,\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} + \left(\frac{3}{4} + x_4\right)a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Na
$\mathbf{B}_{24}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{3}{4} + x_4\right)\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Na
$\mathbf{B}_{25}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{3}{4} + x_4\right)\,\mathbf{a}_2 + \left(\frac{3}{4} + x_4\right)\,\mathbf{a}_3$	=	$\left(\frac{3}{4} + x_4\right)a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Na
$\mathbf{B}_{26}$	=	$\left(\frac{3}{4} + x_4\right)\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Na
$\mathbf{B}_{27}$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \left(\frac{3}{4} + x_4\right)\,\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} - x_4a\,\hat{\mathbf{z}}$	(48*f*)	Na
$\mathbf{B}_{28}$	=	$\left(\frac{3}{4} + x_4\right)\,\mathbf{a}_1 + \left(\frac{3}{4} + x_4\right)\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} + \left(\frac{3}{4} + x_4\right)a\,\hat{\mathbf{z}}$	(48*f*)	Na
$\mathbf{B}_{29}$	=	$z_5\,\mathbf{a}_1 + z_5\,\mathbf{a}_2 + (2x_5 - z_5)\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}} + x_5a\,\hat{\mathbf{y}} + z_5a\,\hat{\mathbf{z}}$	(96*g*)	O
$\mathbf{B}_{30}$	=	$z_5\,\mathbf{a}_1 + z_5\,\mathbf{a}_2 + \left(\frac{1}{2} - 2x_5 - z_5\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_5\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_5\right)a\,\hat{\mathbf{y}} + z_5a\,\hat{\mathbf{z}}$	(96*g*)	O
$\mathbf{B}_{31}$	=	$(2x_5 - z_5)\,\mathbf{a}_1 + \left(\frac{1}{2} - 2x_5 - z_5\right)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_5\right)a\,\hat{\mathbf{x}} + x_5a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_5\right)a\,\hat{\mathbf{z}}$	(96*g*)	O
$\mathbf{B}_{32}$	=	$\left(\frac{1}{2} - 2x_5 - z_5\right)\,\mathbf{a}_1 + (2x_5 - z_5)\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_5\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - z_5\right)a\,\hat{\mathbf{z}}$	(96*g*)	O
$\mathbf{B}_{33}$	=	$(2x_5 - z_5)\,\mathbf{a}_1 + z_5\,\mathbf{a}_2 + z_5\,\mathbf{a}_3$	=	$z_5a\,\hat{\mathbf{x}} + x_5a\,\hat{\mathbf{y}} + x_5a\,\hat{\mathbf{z}}$	(96*g*)	O

$\mathbf{B}_{34}$	=	$\left(\frac{1}{2}-2x_5-z_5\right)\mathbf{a}_1+z_5\,\mathbf{a}_2+z_5\,\mathbf{a}_3$	=	$z_5a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-x_5\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-x_5\right)a\,\hat{\mathbf{z}}$	(96*g*)	O
$\mathbf{B}_{35}$	=	$z_5\,\mathbf{a}_1+(2x_5-z_5)\,\mathbf{a}_2+\left(\frac{1}{2}-2x_5-z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}-z_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-x_5\right)a\,\hat{\mathbf{y}}+x_5a\,\hat{\mathbf{z}}$	(96*g*)	O
$\mathbf{B}_{36}$	=	$z_5\,\mathbf{a}_1+\left(\frac{1}{2}-2x_5-z_5\right)\mathbf{a}_2+(2x_5-z_5)\,\mathbf{a}_3$	=	$\left(\frac{1}{4}-z_5\right)a\,\hat{\mathbf{x}}+x_5a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-x_5\right)a\,\hat{\mathbf{z}}$	(96*g*)	O
$\mathbf{B}_{37}$	=	$z_5\,\mathbf{a}_1+(2x_5-z_5)\,\mathbf{a}_2+z_5\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}}+z_5a\,\hat{\mathbf{y}}+x_5a\,\hat{\mathbf{z}}$	(96*g*)	O
$\mathbf{B}_{38}$	=	$z_5\,\mathbf{a}_1+\left(\frac{1}{2}-2x_5-z_5\right)\mathbf{a}_2+z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{4}-x_5\right)a\,\hat{\mathbf{x}}+z_5a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-x_5\right)a\,\hat{\mathbf{z}}$	(96*g*)	O
$\mathbf{B}_{39}$	=	$\left(\frac{1}{2}-2x_5-z_5\right)\mathbf{a}_1+z_5\,\mathbf{a}_2+(2x_5-z_5)\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-z_5\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-x_5\right)a\,\hat{\mathbf{z}}$	(96*g*)	O
$\mathbf{B}_{40}$	=	$(2x_5-z_5)\,\mathbf{a}_1+z_5\,\mathbf{a}_2+\left(\frac{1}{2}-2x_5-z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}-x_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-z_5\right)a\,\hat{\mathbf{y}}+x_5a\,\hat{\mathbf{z}}$	(96*g*)	O
$\mathbf{B}_{41}$	=	$-z_5\,\mathbf{a}_1-z_5\,\mathbf{a}_2+\left(\frac{1}{2}+2x_5+z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}+x_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+x_5\right)a\,\hat{\mathbf{y}}-z_5a\,\hat{\mathbf{z}}$	(96*g*)	O
$\mathbf{B}_{42}$	=	$-z_5\,\mathbf{a}_1-z_5\,\mathbf{a}_2+(-2x_5+z_5)\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}}-x_5a\,\hat{\mathbf{y}}-z_5a\,\hat{\mathbf{z}}$	(96*g*)	O
$\mathbf{B}_{43}$	=	$(-2x_5+z_5)\,\mathbf{a}_1+\left(\frac{1}{2}+2x_5+z_5\right)\mathbf{a}_2-z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{4}+x_5\right)a\,\hat{\mathbf{x}}-x_5a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+z_5\right)a\,\hat{\mathbf{z}}$	(96*g*)	O
$\mathbf{B}_{44}$	=	$\left(\frac{1}{2}+2x_5+z_5\right)\mathbf{a}_1+(-2x_5+z_5)\,\mathbf{a}_2-z_5\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+x_5\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+z_5\right)a\,\hat{\mathbf{z}}$	(96*g*)	O
$\mathbf{B}_{45}$	=	$(-2x_5+z_5)\,\mathbf{a}_1-z_5\,\mathbf{a}_2+\left(\frac{1}{2}+2x_5+z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}+x_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+z_5\right)a\,\hat{\mathbf{y}}-x_5a\,\hat{\mathbf{z}}$	(96*g*)	O
$\mathbf{B}_{46}$	=	$\left(\frac{1}{2}+2x_5+z_5\right)\mathbf{a}_1-z_5\,\mathbf{a}_2+(-2x_5+z_5)\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+z_5\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+x_5\right)a\,\hat{\mathbf{z}}$	(96*g*)	O
$\mathbf{B}_{47}$	=	$-z_5\,\mathbf{a}_1+(-2x_5+z_5)\,\mathbf{a}_2-z_5\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}}-z_5a\,\hat{\mathbf{y}}-x_5a\,\hat{\mathbf{z}}$	(96*g*)	O
$\mathbf{B}_{48}$	=	$-z_5\,\mathbf{a}_1+\left(\frac{1}{2}+2x_5+z_5\right)\mathbf{a}_2-z_5\,\mathbf{a}_3$	=	$\left(\frac{1}{4}+x_5\right)a\,\hat{\mathbf{x}}-z_5a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+x_5\right)a\,\hat{\mathbf{z}}$	(96*g*)	O
$\mathbf{B}_{49}$	=	$-z_5\,\mathbf{a}_1+(-2x_5+z_5)\,\mathbf{a}_2+\left(\frac{1}{2}+2x_5+z_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}+z_5\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+x_5\right)a\,\hat{\mathbf{y}}-x_5a\,\hat{\mathbf{z}}$	(96*g*)	O
$\mathbf{B}_{50}$	=	$-z_5\,\mathbf{a}_1+\left(\frac{1}{2}+2x_5+z_5\right)\mathbf{a}_2+(-2x_5+z_5)\,\mathbf{a}_3$	=	$\left(\frac{1}{4}+z_5\right)a\,\hat{\mathbf{x}}-x_5a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+x_5\right)a\,\hat{\mathbf{z}}$	(96*g*)	O
$\mathbf{B}_{51}$	=	$\left(\frac{1}{2}+2x_5+z_5\right)\mathbf{a}_1-z_5\,\mathbf{a}_2-z_5\,\mathbf{a}_3$	=	$-z_5a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+x_5\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+x_5\right)a\,\hat{\mathbf{z}}$	(96*g*)	O
$\mathbf{B}_{52}$	=	$(-2x_5+z_5)\,\mathbf{a}_1-z_5\,\mathbf{a}_2-z_5\,\mathbf{a}_3$	=	$-z_5a\,\hat{\mathbf{x}}-x_5a\,\hat{\mathbf{y}}-x_5a\,\hat{\mathbf{z}}$	(96*g*)	O

References:

- H. Shiba and T. Watanabé, *Les structures des cristaux de northupite, de northupite bromée et de tychite*, C. R. Acad. Sci. C **193**, 1421–1423 (1931). http://rruff.info/uploads/CRHSAS193_1421.pdf.
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Found in:

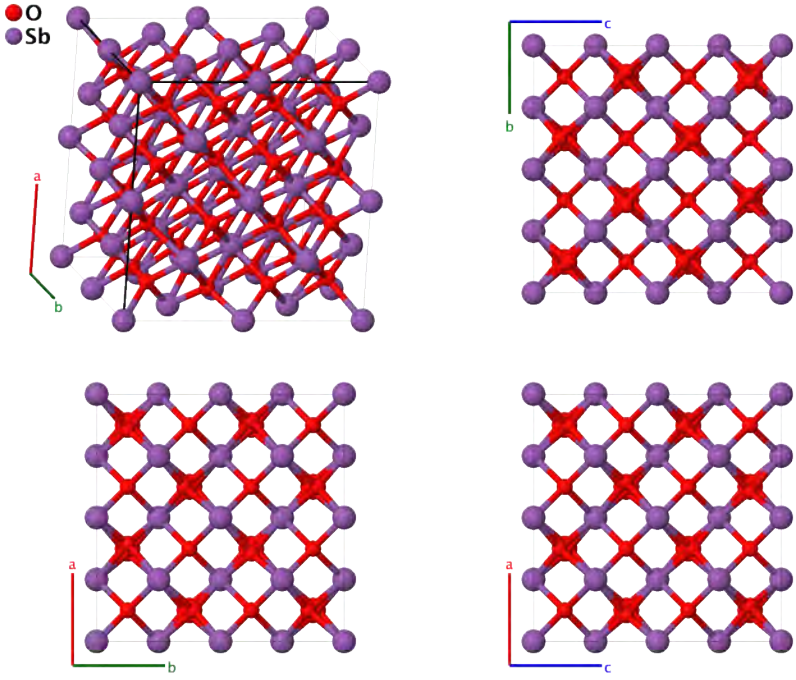
- R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

Geometry files:

- CIF: pp. 1828
- POSCAR: pp. 1829

$D6_2$ (Sb_2O_4) (*obsolete*) Structure: A2B_cF96_227_abf_cd

http://aflow.org/prototype-encyclopedia/A2B_cF96_227_abf_cd

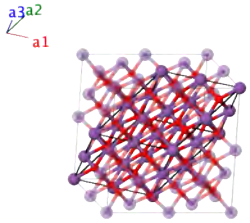


Prototype	:	O_2Sb
AFLOW prototype label	:	A2B_cF96_227_abf_cd
***Strukturbericht* designation**	:	$D6_2$
Pearson symbol	:	cF96
Space group number	:	227
Space group symbol	:	$Fd\bar{3}m$
AFLOW prototype command	:	`aflow --proto=A2B_cF96_227_abf_cd --params=`a, x_5

- Shortly after (Gottfried, 1937) gave this compound the *Strukturbericht* designation $D6_2$, (Dihiström, 1937) showed that they were actually determining the structure of Sb_3O_6OH, making this structure obsolete. Indeed, (Herrman, 1943) formally withdraws this from the *Strukturbericht* list, saying "The type and description [in (Gottfried, 1937)] should be deleted, as the x-rays were not based on the supposed substance." We present it for its historical interest.
- Modern experiments have determined that SbO_2 appears as cervantite (α-Sb_2O_4) or clinocervantite (β-Sb_2O_4).

Face-centered Cubic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{y}} + \frac{1}{2}\,a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{1}{2}\,a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{1}{2}\,a\,\hat{\mathbf{y}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{8}\,\mathbf{a}_1 + \frac{1}{8}\,\mathbf{a}_2 + \frac{1}{8}\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(8*a*)	O I
$\mathbf{B}_2$	=	$\frac{7}{8}\,\mathbf{a}_1 + \frac{7}{8}\,\mathbf{a}_2 + \frac{7}{8}\,\mathbf{a}_3$	=	$\frac{7}{8}a\,\hat{\mathbf{x}} + \frac{7}{8}a\,\hat{\mathbf{y}} + \frac{7}{8}a\,\hat{\mathbf{z}}$	(8*a*)	O I
$\mathbf{B}_3$	=	$\frac{3}{8}\,\mathbf{a}_1 + \frac{3}{8}\,\mathbf{a}_2 + \frac{3}{8}\,\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}}$	(8*b*)	O II
$\mathbf{B}_4$	=	$\frac{5}{8}\,\mathbf{a}_1 + \frac{5}{8}\,\mathbf{a}_2 + \frac{5}{8}\,\mathbf{a}_3$	=	$\frac{5}{8}a\,\hat{\mathbf{x}} + \frac{5}{8}a\,\hat{\mathbf{y}} + \frac{5}{8}a\,\hat{\mathbf{z}}$	(8*b*)	O II
$\mathbf{B}_5$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(16*c*)	Sb I
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}}$	(16*c*)	Sb I
$\mathbf{B}_7$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(16*c*)	Sb I
$\mathbf{B}_8$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$\frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(16*c*)	Sb I
$\mathbf{B}_9$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(16*d*)	Sb II
$\mathbf{B}_{10}$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(16*d*)	Sb II
$\mathbf{B}_{11}$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(16*d*)	Sb II
$\mathbf{B}_{12}$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(16*d*)	Sb II
$\mathbf{B}_{13}$	=	$\left(\frac{1}{4} - x_5\right)\mathbf{a}_1 + x_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$x_5 a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(48*f*)	O III
$\mathbf{B}_{14}$	=	$x_5\,\mathbf{a}_1 + \left(\frac{1}{4} - x_5\right)\mathbf{a}_2 + \left(\frac{1}{4} - x_5\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_5\right)a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(48*f*)	O III
$\mathbf{B}_{15}$	=	$x_5\,\mathbf{a}_1 + \left(\frac{1}{4} - x_5\right)\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + x_5 a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(48*f*)	O III
$\mathbf{B}_{16}$	=	$\left(\frac{1}{4} - x_5\right)\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \left(\frac{1}{4} - x_5\right)\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_5\right)a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	(48*f*)	O III
$\mathbf{B}_{17}$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \left(\frac{1}{4} - x_5\right)\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + x_5 a\,\hat{\mathbf{z}}$	(48*f*)	O III
$\mathbf{B}_{18}$	=	$\left(\frac{1}{4} - x_5\right)\mathbf{a}_1 + \left(\frac{1}{4} - x_5\right)\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_5\right)a\,\hat{\mathbf{z}}$	(48*f*)	O III
$\mathbf{B}_{19}$	=	$\left(\frac{3}{4} + x_5\right)\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \left(\frac{3}{4} + x_5\right)\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} + \left(\frac{3}{4} + x_5\right)a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}}$	(48*f*)	O III
$\mathbf{B}_{20}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{3}{4} + x_5\right)\mathbf{a}_2 - x_5\,\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} - x_5 a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}}$	(48*f*)	O III
$\mathbf{B}_{21}$	=	$-x_5\,\mathbf{a}_1 + \left(\frac{3}{4} + x_5\right)\mathbf{a}_2 + \left(\frac{3}{4} + x_5\right)\mathbf{a}_3$	=	$\left(\frac{3}{4} + x_5\right)a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}}$	(48*f*)	O III
$\mathbf{B}_{22}$	=	$\left(\frac{3}{4} + x_5\right)\mathbf{a}_1 - x_5\,\mathbf{a}_2 - x_5\,\mathbf{a}_3$	=	$-x_5 a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}}$	(48*f*)	O III
$\mathbf{B}_{23}$	=	$-x_5\,\mathbf{a}_1 - x_5\,\mathbf{a}_2 + \left(\frac{3}{4} + x_5\right)\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} - x_5 a\,\hat{\mathbf{z}}$	(48*f*)	O III
$\mathbf{B}_{24}$	=	$\left(\frac{3}{4} + x_5\right)\mathbf{a}_1 + \left(\frac{3}{4} + x_5\right)\mathbf{a}_2 - x_5\,\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} + \left(\frac{3}{4} + x_5\right)a\,\hat{\mathbf{z}}$	(48*f*)	O III

References:

- G. Natta and M. Baccaredda, *Tetrossido di antimonio e antimoniati*, Zeitschrift für Kristallographie - Crystalline Materials

85, 271–296 (1933), doi:10.1524/zkri.1933.85.1.271.
- K. Dihlström and A. Westgren, *Über den Bau des sogenannten Antimontetroxyds und der damit isomorphen Verbindung* $BiTa_2O_6F$, Z. Anorg. Allg. Chem. **235**, 153–160 (1937), doi:10.1002/zaac.19372350121.
- K. Herrmann, ed., *Strukturbericht Band VII 1939* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1943).

Found in:

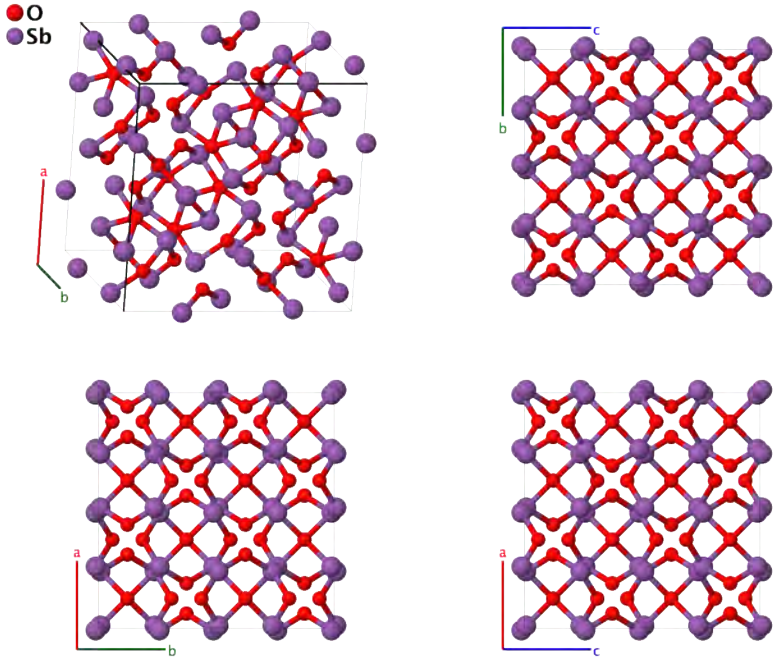
- C. Gottfried and F. Schossberger, eds., *Strukturbericht Band III 1933-1935* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Geometry files:

- CIF: pp. 1830
- POSCAR: pp. 1831

Senarmontite (Sb_2O_3, $D6_1$) Structure: A3B2_cF80_227_f_e

http://aflow.org/prototype-encyclopedia/A3B2_cF80_227_f_e



Prototype	:	O_3Sb_2
AFLOW prototype label	:	A3B2_cF80_227_f_e
***Strukturbericht* designation**	:	$D6_1$
Pearson symbol	:	cF80
Space group number	:	227
Space group symbol	:	$Fd\bar{3}m$
AFLOW prototype command	:	`aflow --proto=A3B2_cF80_227_f_e --params=`a, x_1, x_2

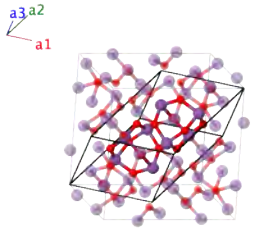
Other compounds with this structure

- As_2O_3 (arsenolite)

- (Ewald, 1931) designated this as *Strukturbericht* $D6_1$, however (Parthé, 1993) and (Villars, 1991) label this as *Strukturbericht* $D5_4$, and Parthé uses As_2O_3 as the prototype. While this structure obviously fits better with the $D5$ series (A_2B_3) than $D6$ (A_2B_4), the $D5_4$ structure was (inadvertently?) omitted from (Hermann, 1937), which jumps from $D5_3$ to $D5_5$. We will follow this historical record (Ewald, 1931) here.
- This is the cubic form of Sb_2O_3. For the orthorhombic form see the valentinite ($D5_{11}$) structure.
- (Svensson, 1975) gave the atomic coordinates in setting 1 of space group $Fd\bar{3}m$ #227. We used FINDSYM to shift the coordinates to the standard setting 2.

Face-centered Cubic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}\,a\,\hat{\mathbf{y}} + \frac{1}{2}\,a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{1}{2}\,a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}\,a\,\hat{\mathbf{x}} + \frac{1}{2}\,a\,\hat{\mathbf{y}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$x_1\,\mathbf{a}_1 + x_1\,\mathbf{a}_2 + x_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + x_1 a\,\hat{\mathbf{y}} + x_1 a\,\hat{\mathbf{z}}$	$(32e)$	Sb
$\mathbf{B}_2$	=	$x_1\,\mathbf{a}_1 + x_1\,\mathbf{a}_2 + \left(\frac{1}{2} - 3x_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_1\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_1\right)a\,\hat{\mathbf{y}} + x_1 a\,\hat{\mathbf{z}}$	$(32e)$	Sb
$\mathbf{B}_3$	=	$x_1\,\mathbf{a}_1 + \left(\frac{1}{2} - 3x_1\right)\mathbf{a}_2 + x_1\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_1\right)a\,\hat{\mathbf{x}} + x_1 a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_1\right)a\,\hat{\mathbf{z}}$	$(32e)$	Sb
$\mathbf{B}_4$	=	$\left(\frac{1}{2} - 3x_1\right)\mathbf{a}_1 + x_1\,\mathbf{a}_2 + x_1\,\mathbf{a}_3$	=	$x_1 a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_1\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_1\right)a\,\hat{\mathbf{z}}$	$(32e)$	Sb
$\mathbf{B}_5$	=	$-x_1\,\mathbf{a}_1 - x_1\,\mathbf{a}_2 + \left(\frac{1}{2} + 3x_1\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_1\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_1\right)a\,\hat{\mathbf{y}} - x_1 a\,\hat{\mathbf{z}}$	$(32e)$	Sb
$\mathbf{B}_6$	=	$-x_1\,\mathbf{a}_1 - x_1\,\mathbf{a}_2 - x_1\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} - x_1 a\,\hat{\mathbf{y}} - x_1 a\,\hat{\mathbf{z}}$	$(32e)$	Sb
$\mathbf{B}_7$	=	$-x_1\,\mathbf{a}_1 + \left(\frac{1}{2} + 3x_1\right)\mathbf{a}_2 - x_1\,\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_1\right)a\,\hat{\mathbf{x}} - x_1 a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + x_1\right)a\,\hat{\mathbf{z}}$	$(32e)$	Sb
$\mathbf{B}_8$	=	$\left(\frac{1}{2} + 3x_1\right)\mathbf{a}_1 - x_1\,\mathbf{a}_2 - x_1\,\mathbf{a}_3$	=	$-x_1 a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_1\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + x_1\right)a\,\hat{\mathbf{z}}$	$(32e)$	Sb
$\mathbf{B}_9$	=	$\left(\frac{1}{4} - x_2\right)\mathbf{a}_1 + x_2\,\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$x_2 a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	$(48f)$	O
$\mathbf{B}_{10}$	=	$x_2\,\mathbf{a}_1 + \left(\frac{1}{4} - x_2\right)\mathbf{a}_2 + \left(\frac{1}{4} - x_2\right)\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_2\right)a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	$(48f)$	O
$\mathbf{B}_{11}$	=	$x_2\,\mathbf{a}_1 + \left(\frac{1}{4} - x_2\right)\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + x_2 a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	$(48f)$	O
$\mathbf{B}_{12}$	=	$\left(\frac{1}{4} - x_2\right)\mathbf{a}_1 + x_2\,\mathbf{a}_2 + \left(\frac{1}{4} - x_2\right)\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_2\right)a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	$(48f)$	O
$\mathbf{B}_{13}$	=	$x_2\,\mathbf{a}_1 + x_2\,\mathbf{a}_2 + \left(\frac{1}{4} - x_2\right)\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + x_2 a\,\hat{\mathbf{z}}$	$(48f)$	O
$\mathbf{B}_{14}$	=	$\left(\frac{1}{4} - x_2\right)\mathbf{a}_1 + \left(\frac{1}{4} - x_2\right)\mathbf{a}_2 + x_2\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_2\right)a\,\hat{\mathbf{z}}$	$(48f)$	O
$\mathbf{B}_{15}$	=	$\left(\frac{3}{4} + x_2\right)\mathbf{a}_1 - x_2\,\mathbf{a}_2 + \left(\frac{3}{4} + x_2\right)\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} + \left(\frac{3}{4} + x_2\right)a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}}$	$(48f)$	O
$\mathbf{B}_{16}$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{3}{4} + x_2\right)\mathbf{a}_2 - x_2\,\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} - x_2 a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}}$	$(48f)$	O
$\mathbf{B}_{17}$	=	$-x_2\,\mathbf{a}_1 + \left(\frac{3}{4} + x_2\right)\mathbf{a}_2 + \left(\frac{3}{4} + x_2\right)\mathbf{a}_3$	=	$\left(\frac{3}{4} + x_2\right)a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}}$	$(48f)$	O
$\mathbf{B}_{18}$	=	$\left(\frac{3}{4} + x_2\right)\mathbf{a}_1 - x_2\,\mathbf{a}_2 - x_2\,\mathbf{a}_3$	=	$-x_2 a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} + \frac{3}{8}a\,\hat{\mathbf{z}}$	$(48f)$	O
$\mathbf{B}_{19}$	=	$-x_2\,\mathbf{a}_1 - x_2\,\mathbf{a}_2 + \left(\frac{3}{4} + x_2\right)\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} - x_2 a\,\hat{\mathbf{z}}$	$(48f)$	O
$\mathbf{B}_{20}$	=	$\left(\frac{3}{4} + x_2\right)\mathbf{a}_1 + \left(\frac{3}{4} + x_2\right)\mathbf{a}_2 - x_2\,\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}} + \frac{3}{8}a\,\hat{\mathbf{y}} + \left(\frac{3}{4} + x_2\right)a\,\hat{\mathbf{z}}$	$(48f)$	O

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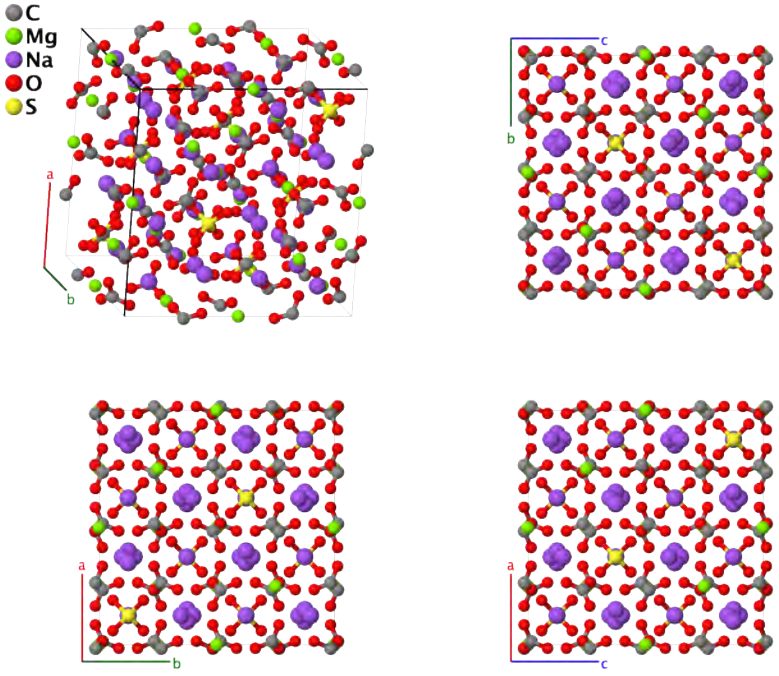
Geometry files:
- CIF: pp. 1831
- POSCAR: pp. 1832

$H5_6$ [Tychite, $Na_6Mg_2SO_4(CO_3)_4$)] (*obsolete*) Structure: A4B2C6D16E_cF232_227_e_d_f_eg_a

http://aflow.org/prototype-encyclopedia/A4B2C6D16E_cF232_227_e_d_f_eg_a

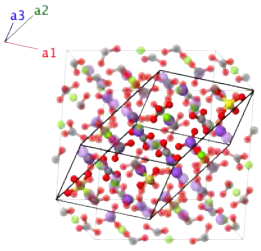


Prototype	:	$C_4Mg_2Na_6O_{16}S$
AFLOW prototype label	:	A4B2C6D16E_cF232_227_e_d_f_eg_a
***Strukturbericht* designation**	:	$H5_6$
Pearson symbol	:	cF232
Space group number	:	227
Space group symbol	:	$Fd\bar{3}m$
AFLOW prototype command	:	`aflow --proto=A4B2C6D16E_cF232_227_e_d_f_eg_a --params=`$a, x_3, x_4, x_5, x_6, z_6$

- This is the original structure determined by (Shiba, 1931) and given the designation $H5_6$ in (Hermann, 1937). (Schmidt, 2006) showed that the true structure is in space group $Fd\bar{3}$ #203, however the two structures are very similar, and a displacement of the oxygen atoms by less than 1 Å brings the two structures into agreement.
- (Hermann, 1937) gives the chemical formula as $Na_6Mg_2SO_4(CO_3)_2$, but the given Wyckoff positions are in agreement with the correct formula, $Na_6Mg_2SO_4(CO_3)_4$.

Face-centered Cubic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= \tfrac{1}{2}\,a\,\hat{\mathbf{y}} + \tfrac{1}{2}\,a\,\hat{\mathbf{z}}\\ \mathbf{a}_2 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{z}}\\ \mathbf{a}_3 &= \tfrac{1}{2}\,a\,\hat{\mathbf{x}} + \tfrac{1}{2}\,a\,\hat{\mathbf{y}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$\frac{1}{8}\,\mathbf{a}_1 + \frac{1}{8}\,\mathbf{a}_2 + \frac{1}{8}\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	$(8a)$	S
$\mathbf{B}_2$	=	$\frac{7}{8}\,\mathbf{a}_1 + \frac{7}{8}\,\mathbf{a}_2 + \frac{7}{8}\,\mathbf{a}_3$	=	$\frac{7}{8}a\,\hat{\mathbf{x}} + \frac{7}{8}a\,\hat{\mathbf{y}} + \frac{7}{8}a\,\hat{\mathbf{z}}$	$(8a)$	S
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	$(16d)$	Mg
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	$(16d)$	Mg
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	$(16d)$	Mg
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	$(16d)$	Mg
$\mathbf{B}_7$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	$(32e)$	C
$\mathbf{B}_8$	=	$x_3\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + \left(\frac{1}{2} - 3x_3\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}}$	$(32e)$	C
$\mathbf{B}_9$	=	$x_3\,\mathbf{a}_1 + \left(\frac{1}{2} - 3x_3\right)\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{z}}$	$(32e)$	C
$\mathbf{B}_{10}$	=	$\left(\frac{1}{2} - 3x_3\right)\,\mathbf{a}_1 + x_3\,\mathbf{a}_2 + x_3\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_3\right)a\,\hat{\mathbf{z}}$	$(32e)$	C
$\mathbf{B}_{11}$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 + \left(\frac{1}{2} + 3x_3\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	$(32e)$	C
$\mathbf{B}_{12}$	=	$-x_3\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}}$	$(32e)$	C
$\mathbf{B}_{13}$	=	$-x_3\,\mathbf{a}_1 + \left(\frac{1}{2} + 3x_3\right)\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{z}}$	$(32e)$	C
$\mathbf{B}_{14}$	=	$\left(\frac{1}{2} + 3x_3\right)\,\mathbf{a}_1 - x_3\,\mathbf{a}_2 - x_3\,\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + x_3\right)a\,\hat{\mathbf{z}}$	$(32e)$	C
$\mathbf{B}_{15}$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	$(32e)$	O I
$\mathbf{B}_{16}$	=	$x_4\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + \left(\frac{1}{2} - 3x_4\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{y}} + x_4a\,\hat{\mathbf{z}}$	$(32e)$	O I
$\mathbf{B}_{17}$	=	$x_4\,\mathbf{a}_1 + \left(\frac{1}{2} - 3x_4\right)\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{x}} + x_4a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{z}}$	$(32e)$	O I
$\mathbf{B}_{18}$	=	$\left(\frac{1}{2} - 3x_4\right)\,\mathbf{a}_1 + x_4\,\mathbf{a}_2 + x_4\,\mathbf{a}_3$	=	$x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} - x_4\right)a\,\hat{\mathbf{z}}$	$(32e)$	O I
$\mathbf{B}_{19}$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 + \left(\frac{1}{2} + 3x_4\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_4\right)a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_4\right)a\,\hat{\mathbf{y}} - x_4a\,\hat{\mathbf{z}}$	$(32e)$	O I
$\mathbf{B}_{20}$	=	$-x_4\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} - x_4a\,\hat{\mathbf{z}}$	$(32e)$	O I
$\mathbf{B}_{21}$	=	$-x_4\,\mathbf{a}_1 + \left(\frac{1}{2} + 3x_4\right)\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$\left(\frac{1}{4} + x_4\right)a\,\hat{\mathbf{x}} - x_4a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + x_4\right)a\,\hat{\mathbf{z}}$	$(32e)$	O I
$\mathbf{B}_{22}$	=	$\left(\frac{1}{2} + 3x_4\right)\,\mathbf{a}_1 - x_4\,\mathbf{a}_2 - x_4\,\mathbf{a}_3$	=	$-x_4a\,\hat{\mathbf{x}} + \left(\frac{1}{4} + x_4\right)a\,\hat{\mathbf{y}} + \left(\frac{1}{4} + x_4\right)a\,\hat{\mathbf{z}}$	$(32e)$	O I
$\mathbf{B}_{23}$	=	$\left(\frac{1}{4} - x_5\right)\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$x_5a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	$(48f)$	Na
$\mathbf{B}_{24}$	=	$x_5\,\mathbf{a}_1 + \left(\frac{1}{4} - x_5\right)\,\mathbf{a}_2 + \left(\frac{1}{4} - x_5\right)\,\mathbf{a}_3$	=	$\left(\frac{1}{4} - x_5\right)a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	$(48f)$	Na
$\mathbf{B}_{25}$	=	$x_5\,\mathbf{a}_1 + \left(\frac{1}{4} - x_5\right)\,\mathbf{a}_2 + x_5\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + x_5a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	$(48f)$	Na
$\mathbf{B}_{26}$	=	$\left(\frac{1}{4} - x_5\right)\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \left(\frac{1}{4} - x_5\right)\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \left(\frac{1}{4} - x_5\right)a\,\hat{\mathbf{y}} + \frac{1}{8}a\,\hat{\mathbf{z}}$	$(48f)$	Na
$\mathbf{B}_{27}$	=	$x_5\,\mathbf{a}_1 + x_5\,\mathbf{a}_2 + \left(\frac{1}{4} - x_5\right)\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}} + \frac{1}{8}a\,\hat{\mathbf{y}} + x_5a\,\hat{\mathbf{z}}$	$(48f)$	Na

$\mathbf{B}_{28}$	=	$\left(\frac{1}{4}-x_5\right)\mathbf{a}_1+\left(\frac{1}{4}-x_5\right)\mathbf{a}_2+x_5\,\mathbf{a}_3$	=	$\frac{1}{8}a\,\hat{\mathbf{x}}+\frac{1}{8}a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-x_5\right)a\,\hat{\mathbf{z}}$	(48*f*)	Na
$\mathbf{B}_{29}$	=	$\left(\frac{3}{4}+x_5\right)\mathbf{a}_1-x_5\,\mathbf{a}_2+\left(\frac{3}{4}+x_5\right)\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}}+\left(\frac{3}{4}+x_5\right)a\,\hat{\mathbf{y}}+\frac{3}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Na
$\mathbf{B}_{30}$	=	$-x_5\,\mathbf{a}_1+\left(\frac{3}{4}+x_5\right)\mathbf{a}_2-x_5\,\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}}-x_5a\,\hat{\mathbf{y}}+\frac{3}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Na
$\mathbf{B}_{31}$	=	$-x_5\,\mathbf{a}_1+\left(\frac{3}{4}+x_5\right)\mathbf{a}_2+\left(\frac{3}{4}+x_5\right)\mathbf{a}_3$	=	$\left(\frac{3}{4}+x_5\right)a\,\hat{\mathbf{x}}+\frac{3}{8}a\,\hat{\mathbf{y}}+\frac{3}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Na
$\mathbf{B}_{32}$	=	$\left(\frac{3}{4}+x_5\right)\mathbf{a}_1-x_5\,\mathbf{a}_2-x_5\,\mathbf{a}_3$	=	$-x_5a\,\hat{\mathbf{x}}+\frac{3}{8}a\,\hat{\mathbf{y}}+\frac{3}{8}a\,\hat{\mathbf{z}}$	(48*f*)	Na
$\mathbf{B}_{33}$	=	$-x_5\,\mathbf{a}_1-x_5\,\mathbf{a}_2+\left(\frac{3}{4}+x_5\right)\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}}+\frac{3}{8}a\,\hat{\mathbf{y}}-x_5a\,\hat{\mathbf{z}}$	(48*f*)	Na
$\mathbf{B}_{34}$	=	$\left(\frac{3}{4}+x_5\right)\mathbf{a}_1+\left(\frac{3}{4}+x_5\right)\mathbf{a}_2-x_5\,\mathbf{a}_3$	=	$\frac{3}{8}a\,\hat{\mathbf{x}}+\frac{3}{8}a\,\hat{\mathbf{y}}+\left(\frac{3}{4}+x_5\right)a\,\hat{\mathbf{z}}$	(48*f*)	Na
$\mathbf{B}_{35}$	=	$z_6\,\mathbf{a}_1+z_6\,\mathbf{a}_2+(2x_6-z_6)\,\mathbf{a}_3$	=	$x_6a\,\hat{\mathbf{x}}+x_6a\,\hat{\mathbf{y}}+z_6a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{36}$	=	$z_6\,\mathbf{a}_1+z_6\,\mathbf{a}_2+\left(\frac{1}{2}-2x_6-z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}-x_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-x_6\right)a\,\hat{\mathbf{y}}+z_6a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{37}$	=	$(2x_6-z_6)\,\mathbf{a}_1+$ $\left(\frac{1}{2}-2x_6-z_6\right)\mathbf{a}_2+z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{4}-x_6\right)a\,\hat{\mathbf{x}}+x_6a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-z_6\right)a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{38}$	=	$\left(\frac{1}{2}-2x_6-z_6\right)\mathbf{a}_1+$ $(2x_6-z_6)\,\mathbf{a}_2+z_6\,\mathbf{a}_3$	=	$x_6a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-x_6\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-z_6\right)a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{39}$	=	$(2x_6-z_6)\,\mathbf{a}_1+z_6\,\mathbf{a}_2+z_6\,\mathbf{a}_3$	=	$z_6a\,\hat{\mathbf{x}}+x_6a\,\hat{\mathbf{y}}+x_6a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{40}$	=	$\left(\frac{1}{2}-2x_6-z_6\right)\mathbf{a}_1+z_6\,\mathbf{a}_2+z_6\,\mathbf{a}_3$	=	$z_6a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-x_6\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-x_6\right)a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{41}$	=	$z_6\,\mathbf{a}_1+(2x_6-z_6)\,\mathbf{a}_2+$ $\left(\frac{1}{2}-2x_6-z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}-z_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-x_6\right)a\,\hat{\mathbf{y}}+x_6a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{42}$	=	$z_6\,\mathbf{a}_1+\left(\frac{1}{2}-2x_6-z_6\right)\mathbf{a}_2+$ $(2x_6-z_6)\,\mathbf{a}_3$	=	$\left(\frac{1}{4}-z_6\right)a\,\hat{\mathbf{x}}+x_6a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-x_6\right)a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{43}$	=	$z_6\,\mathbf{a}_1+(2x_6-z_6)\,\mathbf{a}_2+z_6\,\mathbf{a}_3$	=	$x_6a\,\hat{\mathbf{x}}+z_6a\,\hat{\mathbf{y}}+x_6a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{44}$	=	$z_6\,\mathbf{a}_1+\left(\frac{1}{2}-2x_6-z_6\right)\mathbf{a}_2+z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{4}-x_6\right)a\,\hat{\mathbf{x}}+z_6a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-x_6\right)a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{45}$	=	$\left(\frac{1}{2}-2x_6-z_6\right)\mathbf{a}_1+z_6\,\mathbf{a}_2+$ $(2x_6-z_6)\,\mathbf{a}_3$	=	$x_6a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-z_6\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-x_6\right)a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{46}$	=	$(2x_6-z_6)\,\mathbf{a}_1+z_6\,\mathbf{a}_2+$ $\left(\frac{1}{2}-2x_6-z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}-x_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-z_6\right)a\,\hat{\mathbf{y}}+x_6a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{47}$	=	$-z_6\,\mathbf{a}_1-z_6\,\mathbf{a}_2+\left(\frac{1}{2}+2x_6+z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}+x_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+x_6\right)a\,\hat{\mathbf{y}}-z_6a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{48}$	=	$-z_6\,\mathbf{a}_1-z_6\,\mathbf{a}_2+(-2x_6+z_6)\,\mathbf{a}_3$	=	$-x_6a\,\hat{\mathbf{x}}-x_6a\,\hat{\mathbf{y}}-z_6a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{49}$	=	$(-2x_6+z_6)\,\mathbf{a}_1+$ $\left(\frac{1}{2}+2x_6+z_6\right)\mathbf{a}_2-z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{4}+x_6\right)a\,\hat{\mathbf{x}}-x_6a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+z_6\right)a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{50}$	=	$\left(\frac{1}{2}+2x_6+z_6\right)\mathbf{a}_1+$ $(-2x_6+z_6)\,\mathbf{a}_2-z_6\,\mathbf{a}_3$	=	$-x_6a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+x_6\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+z_6\right)a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{51}$	=	$(-2x_6+z_6)\,\mathbf{a}_1-z_6\,\mathbf{a}_2+$ $\left(\frac{1}{2}+2x_6+z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}+x_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+z_6\right)a\,\hat{\mathbf{y}}-x_6a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{52}$	=	$\left(\frac{1}{2}+2x_6+z_6\right)\mathbf{a}_1-z_6\,\mathbf{a}_2+$ $(-2x_6+z_6)\,\mathbf{a}_3$	=	$-x_6a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+z_6\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+x_6\right)a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{53}$	=	$-z_6\,\mathbf{a}_1+(-2x_6+z_6)\,\mathbf{a}_2-z_6\,\mathbf{a}_3$	=	$-x_6a\,\hat{\mathbf{x}}-z_6a\,\hat{\mathbf{y}}-x_6a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{54}$	=	$-z_6\,\mathbf{a}_1+\left(\frac{1}{2}+2x_6+z_6\right)\mathbf{a}_2-z_6\,\mathbf{a}_3$	=	$\left(\frac{1}{4}+x_6\right)a\,\hat{\mathbf{x}}-z_6a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+x_6\right)a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{55}$	=	$-z_6\,\mathbf{a}_1+(-2x_6+z_6)\,\mathbf{a}_2+$ $\left(\frac{1}{2}+2x_6+z_6\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}+z_6\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+x_6\right)a\,\hat{\mathbf{y}}-x_6a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{56}$	=	$-z_6\,\mathbf{a}_1+\left(\frac{1}{2}+2x_6+z_6\right)\mathbf{a}_2+$ $(-2x_6+z_6)\,\mathbf{a}_3$	=	$\left(\frac{1}{4}+z_6\right)a\,\hat{\mathbf{x}}-x_6a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+x_6\right)a\,\hat{\mathbf{z}}$	(96*g*)	O II
$\mathbf{B}_{57}$	=	$\left(\frac{1}{2}+2x_6+z_6\right)\mathbf{a}_1-z_6\,\mathbf{a}_2-z_6\,\mathbf{a}_3$	=	$-z_6a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+x_6\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+x_6\right)a\,\hat{\mathbf{z}}$	(96*g*)	O II

$$\mathbf{B}_{58} = (-2x_6 + z_6)\,\mathbf{a}_1 - z_6\,\mathbf{a}_2 - z_6\,\mathbf{a}_3 = -z_6 a\,\hat{\mathbf{x}} - x_6 a\,\hat{\mathbf{y}} - x_6 a\,\hat{\mathbf{z}} \quad (96g) \quad \text{O II}$$

References:

- H. Shiba and T. Watanabé, *Les structures des cristaux de northupite, de northupite bromée et de tychite*, C. R. Acad. Sci. C **193**, 1421–1423 (1931). http://rruff.info/uploads/CRHSAS193_1421.pdf.
- C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

Found in:

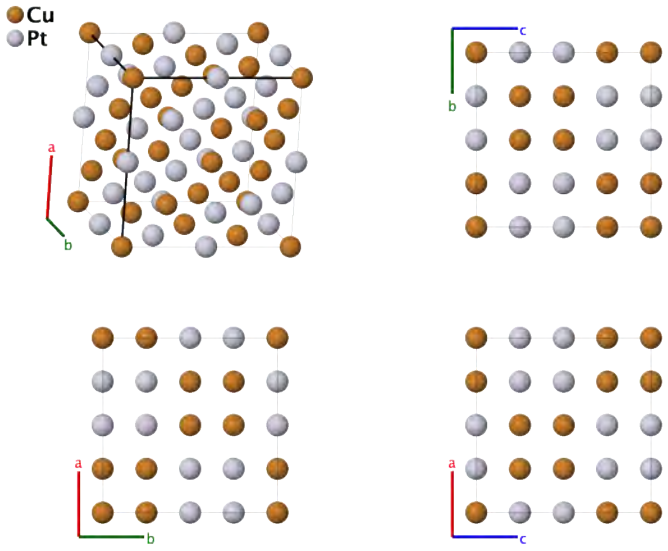
- G. R. Schmidt, J. Reynard, H. Yang, and R. T. Downs, *Tychite, $Na_6Mg_2(SO_4)(CO_3)_4$: Structure analysis and Raman spectroscopic data*, Acta Crystallogr. E **62**, i207–i209 (2006), doi:10.1107/S160053680603491X.

Geometry files:

- CIF: pp. 1833
- POSCAR: pp. 1834

Cubic CuPt ($L1_3$ (I), $D4$) Structure: AB_cF32_227_c_d

http://aflow.org/prototype-encyclopedia/AB_cF32_227_c_d

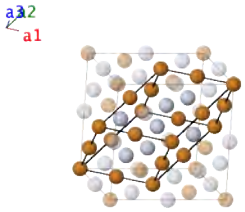


Prototype	:	CuPt
AFLOW prototype label	:	AB_cF32_227_c_d
***Strukturbericht* designation**	:	$L1_3$ (I)
Pearson symbol	:	cF32
Space group number	:	227
Space group symbol	:	$Fd\bar{3}m$
AFLOW prototype command	:	`aflow --proto=AB_cF32_227_c_d --params=`a

- (Johansson, 1929) described two possible structures for CuPt. (Ewald, 1929) and later (Villars, 2007) used the description to determine the space group and atomic positions. This page describes the cubic structure, which (Ewald, 1929) labeled *Strukturbericht* $L1_1$. The other structure is rhombohedral, and was listed as $L1_1$. (Villars, 2007) prefers the later structure, listing the current one as "superceded."
- (Barrett, 1980) noted that even slight additions of platinum above stoichiometry will cause a change in the crystal structure.
- This structure is equivalent to the $D4$ structure of (Lu, 1991).
- This structure should not be confused with the $CuPt_3$ structure, which has also been given the $L1_3$ label, and which we will often refer to as $L1_3$ (II).

Face-centered Cubic primitive vectors:

$$\mathbf{a}_1 = \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(16*c*)	Cu
$\mathbf{B}_2$	=	$\frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}}$	(16*c*)	Cu
$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(16*c*)	Cu
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1$	=	$\frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(16*c*)	Cu
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(16*d*)	Pt
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(16*d*)	Pt
$\mathbf{B}_7$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(16*d*)	Pt
$\mathbf{B}_8$	=	$\frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(16*d*)	Pt

References:

- C. H. Johansson and J. O. Linde, *Gitterstruktur und elektrisches Leitvermögen der Mischkristallreihen Au-Cu, Pd-Cu und Pt-Cu*, Ann. Phys. **387**, 449–478 (1927), doi:10.1002/andp.19273870402.
- P. Villars, K. Cenzual, J. Daams, R. Gladyshevskii, O. Shcherban, V. Dubenskyy, N. Melnichenko-Koblyuk, O. Pavlyuk, I. Savysyuk, S. Stoyko, and L. Sysa, *Structure Types. Part 5: Space Groups (173) $P6_3$ - (166) R-3m · CuPt: Datasheet from Landolt-Börnstein - Group III Condensed Matter · Volume 43A5: "Structure Types. Part 5: Space Groups (173) $P6_3$ - (166) R-3m" in SpringerMaterials*, doi:10.1007/978-3-540-46933-9_359. Copyright 2007 Springer-Verlag", Part of SpringerMaterials.
- C. Barrett and T. B. Massalski, *Structure of Metals – Crystallographic Methods, Principles, and Data* (Pergammon Press, Oxford, New York, 1980).
- Z. W. Lu, S.-H. Wei, A. Zunger, S. Frota-Pessoa, and L. G. Ferreira, *First-principles statistical mechanics of structural stability of intermetallic compounds*, Phys. Rev. B **44**, 512–544 (1991), doi:10.1103/PhysRevB.44.512.

Found in:

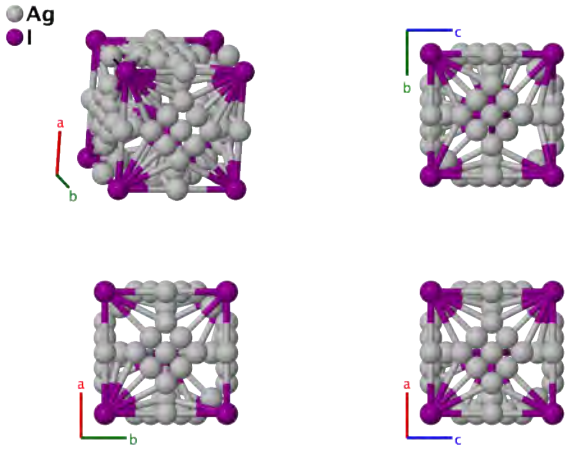
- P. P. Ewald and C. Hermann, eds., *Strukturbericht 1913-1928* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1931).

Geometry files:

- CIF: pp. 1834
- POSCAR: pp. 1835

α-AgI ($B23$) Structure: A21B_cI44_229_bdh_a

http://aflow.org/prototype-encyclopedia/A21B_cI44_229_bdh_a

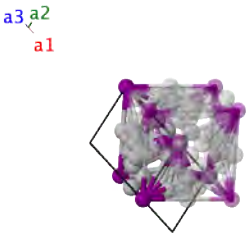


Prototype	:	AgI
AFLOW prototype label	:	A21B_cI44_229_bdh_a
***Strukturbericht* designation**	:	$B23$
Pearson symbol	:	cI44
Space group number	:	229
Space group symbol	:	$Im\bar{3}m$
AFLOW prototype command	:	`aflow --proto=A21B_cI44_229_bdh_a --params=`a, y_4

- Under ambient conditions, silver iodide exists as a mixture of β-AgI, which has the wurtzite ($B4$) structure, and γ-AgI, which has the zincblende ($B3$) structure (Hull, 2004). Above 420 K, AgI transforms to this superionic α phase. The iodine atom sits at the ($2a$) site of the bcc lattice of space group #229, while the silver atom is randomly distributed on one of the ($6b$), ($12d$), and ($24h$) Wyckoff sites in each unit cell. On average, then, each of the 21 Ag sites listed above is occupied only 4.762% of the time in any given primitive cell. This easy transport between sites drives the superionic behavior of α-AgI.

Body-centered Cubic primitive vectors:

$$\mathbf{a}_1 = -\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_2 = \tfrac{1}{2}a\,\hat{\mathbf{x}} - \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}a\,\hat{\mathbf{z}}$$
$$\mathbf{a}_3 = \tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} - \tfrac{1}{2}a\,\hat{\mathbf{z}}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	($2a$)	I
$\mathbf{B}_2$	=	$\tfrac{1}{2}\,\mathbf{a}_2 + \tfrac{1}{2}\,\mathbf{a}_3$	=	$\tfrac{1}{2}a\,\hat{\mathbf{x}}$	($6b$)	Ag I

$\mathbf{B}_3$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}}$	(6*b*)	Ag I
$\mathbf{B}_4$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2$	=	$\frac{1}{2}a\,\hat{\mathbf{z}}$	(6*b*)	Ag I
$\mathbf{B}_5$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(12*d*)	Ag II
$\mathbf{B}_6$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{x}} + \frac{1}{2}a\,\hat{\mathbf{y}}$	(12*d*)	Ag II
$\mathbf{B}_7$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{3}{4}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{y}}$	(12*d*)	Ag II
$\mathbf{B}_8$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{4}\,\mathbf{a}_3$	=	$\frac{1}{4}a\,\hat{\mathbf{y}} + \frac{1}{2}a\,\hat{\mathbf{z}}$	(12*d*)	Ag II
$\mathbf{B}_9$	=	$\frac{3}{4}\,\mathbf{a}_1 + \frac{1}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{y}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(12*d*)	Ag II
$\mathbf{B}_{10}$	=	$\frac{1}{4}\,\mathbf{a}_1 + \frac{3}{4}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{4}a\,\hat{\mathbf{z}}$	(12*d*)	Ag II
$\mathbf{B}_{11}$	=	$2y_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + y_4\,\mathbf{a}_3$	=	$y_4a\,\hat{\mathbf{y}} + y_4a\,\hat{\mathbf{z}}$	(24*h*)	Ag III
$\mathbf{B}_{12}$	=	$y_4\,\mathbf{a}_2 - y_4\,\mathbf{a}_3$	=	$-y_4a\,\hat{\mathbf{y}} + y_4a\,\hat{\mathbf{z}}$	(24*h*)	Ag III
$\mathbf{B}_{13}$	=	$-y_4\,\mathbf{a}_2 + y_4\,\mathbf{a}_3$	=	$y_4a\,\hat{\mathbf{y}} - y_4a\,\hat{\mathbf{z}}$	(24*h*)	Ag III
$\mathbf{B}_{14}$	=	$-2y_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - y_4\,\mathbf{a}_3$	=	$-y_4a\,\hat{\mathbf{y}} - y_4a\,\hat{\mathbf{z}}$	(24*h*)	Ag III
$\mathbf{B}_{15}$	=	$y_4\,\mathbf{a}_1 + 2y_4\,\mathbf{a}_2 + y_4\,\mathbf{a}_3$	=	$y_4a\,\hat{\mathbf{x}} + y_4a\,\hat{\mathbf{z}}$	(24*h*)	Ag III
$\mathbf{B}_{16}$	=	$-y_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_3$	=	$y_4a\,\hat{\mathbf{x}} - y_4a\,\hat{\mathbf{z}}$	(24*h*)	Ag III
$\mathbf{B}_{17}$	=	$y_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_3$	=	$-y_4a\,\hat{\mathbf{x}} + y_4a\,\hat{\mathbf{z}}$	(24*h*)	Ag III
$\mathbf{B}_{18}$	=	$-y_4\,\mathbf{a}_1 - 2y_4\,\mathbf{a}_2 - y_4\,\mathbf{a}_3$	=	$-y_4a\,\hat{\mathbf{x}} - y_4a\,\hat{\mathbf{z}}$	(24*h*)	Ag III
$\mathbf{B}_{19}$	=	$y_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2 + 2y_4\,\mathbf{a}_3$	=	$y_4a\,\hat{\mathbf{x}} + y_4a\,\hat{\mathbf{y}}$	(24*h*)	Ag III
$\mathbf{B}_{20}$	=	$y_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2$	=	$-y_4a\,\hat{\mathbf{x}} + y_4a\,\hat{\mathbf{y}}$	(24*h*)	Ag III
$\mathbf{B}_{21}$	=	$-y_4\,\mathbf{a}_1 + y_4\,\mathbf{a}_2$	=	$y_4a\,\hat{\mathbf{x}} - y_4a\,\hat{\mathbf{y}}$	(24*h*)	Ag III
$\mathbf{B}_{22}$	=	$-y_4\,\mathbf{a}_1 - y_4\,\mathbf{a}_2 - 2y_4\,\mathbf{a}_3$	=	$-y_4a\,\hat{\mathbf{x}} - y_4a\,\hat{\mathbf{y}}$	(24*h*)	Ag III

References:

- L. W. Strock, *Kristallstruktur des Hochtemperatur-Jodsilbers α-AgJ*, Z. Physik. Chem. B **25**, 441–459 (1934), doi:10.1515/zpch-1934-2535.
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Found in:

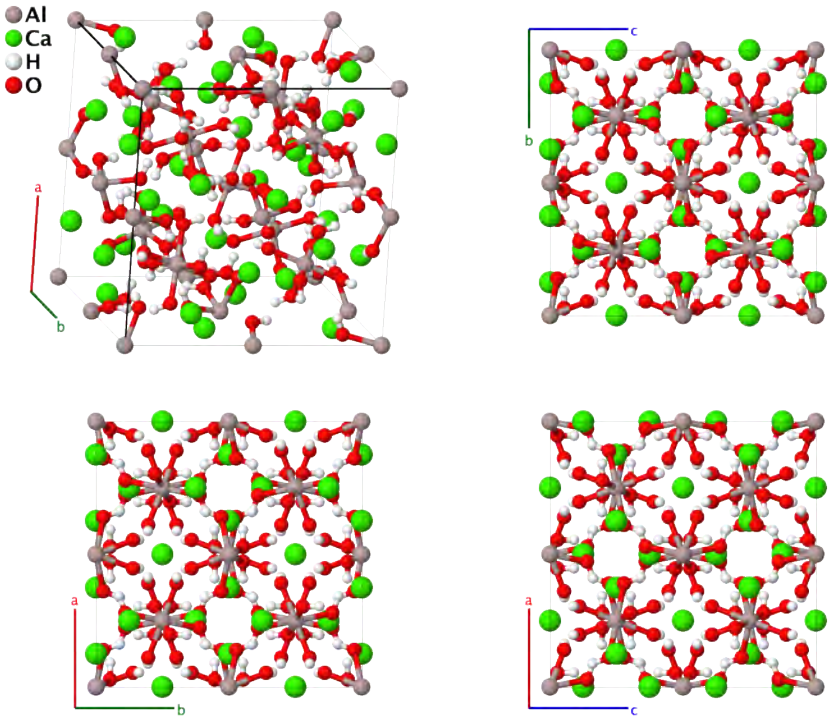
- S. Hoshino, *Crystal Structure and Phase Transition of Some Metallic Halides IV On the Anomalous Structure of α-AgI*, J. Phys. Soc. Jpn. **12**, 315–326 (1957), doi:10.1143/JPSJ.12.315.

Geometry files:

- CIF: pp. 1835
- POSCAR: pp. 1836

$Ca_3Al_2(OH)_{12}$ ($J2_3$) Structure: A2B3C12D12_cI232_230_a_c_h_h

http://aflow.org/prototype-encyclopedia/A2B3C12D12_cI232_230_a_c_h_h

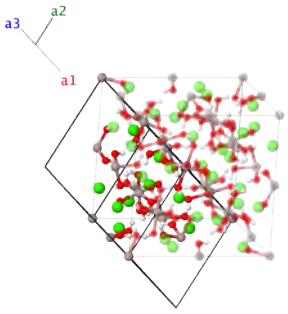


Prototype	:	$Al_2Ca_3H_{12}O_{12}$
AFLOW prototype label	:	A2B3C12D12_cI232_230_a_c_h_h
***Strukturbericht* designation**	:	$J2_3$
Pearson symbol	:	cI232
Space group number	:	230
Space group symbol	:	$Ia\bar{3}d$
AFLOW prototype command	:	`aflow --proto=A2B3C12D12_cI232_230_a_c_h_h --params=`$a, x_3, y_3, z_3, x_4, y_4, z_4$

- The original determination of this structure by (Brandenberger, 1933) did not locate the hydrogen atoms, and according to (Gottfried, 1937) used the coordinates of garnet, $S1_4$. (Bartl, 1986) was able to locate the hydrogen atoms, and as they do not change the space group we include them in the $J2_3$ structure.

Body-centered Cubic primitive vectors:

$$\begin{aligned}\mathbf{a}_1 &= -\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}a\,\hat{\mathbf{z}}\\ \mathbf{a}_2 &= \tfrac{1}{2}a\,\hat{\mathbf{x}} - \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{2}a\,\hat{\mathbf{z}}\\ \mathbf{a}_3 &= \tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} - \tfrac{1}{2}a\,\hat{\mathbf{z}}\end{aligned}$$



Basis vectors:

		Lattice Coordinates		Cartesian Coordinates	Wyckoff Position	Atom Type
$\mathbf{B}_1$	=	$0\,\mathbf{a}_1 + 0\,\mathbf{a}_2 + 0\,\mathbf{a}_3$	=	$0\,\hat{\mathbf{x}} + 0\,\hat{\mathbf{y}} + 0\,\hat{\mathbf{z}}$	(16*a*)	Al
$\mathbf{B}_2$	=	$\tfrac{1}{2}\,\mathbf{a}_1 + \tfrac{1}{2}\,\mathbf{a}_3$	=	$\tfrac{1}{2}a\,\hat{\mathbf{y}}$	(16*a*)	Al
$\mathbf{B}_3$	=	$\tfrac{1}{2}\,\mathbf{a}_2 + \tfrac{1}{2}\,\mathbf{a}_3$	=	$\tfrac{1}{2}a\,\hat{\mathbf{x}}$	(16*a*)	Al
$\mathbf{B}_4$	=	$\tfrac{1}{2}\,\mathbf{a}_1 + \tfrac{1}{2}\,\mathbf{a}_2$	=	$\tfrac{1}{2}a\,\hat{\mathbf{z}}$	(16*a*)	Al
$\mathbf{B}_5$	=	$\tfrac{1}{2}\,\mathbf{a}_1$	=	$\tfrac{3}{4}a\,\hat{\mathbf{x}} + \tfrac{1}{4}a\,\hat{\mathbf{y}} + \tfrac{1}{4}a\,\hat{\mathbf{z}}$	(16*a*)	Al
$\mathbf{B}_6$	=	$\tfrac{1}{2}\,\mathbf{a}_1 + \tfrac{1}{2}\,\mathbf{a}_2 + \tfrac{1}{2}\,\mathbf{a}_3$	=	$\tfrac{1}{4}a\,\hat{\mathbf{x}} + \tfrac{1}{4}a\,\hat{\mathbf{y}} + \tfrac{1}{4}a\,\hat{\mathbf{z}}$	(16*a*)	Al
$\mathbf{B}_7$	=	$\tfrac{1}{2}\,\mathbf{a}_3$	=	$\tfrac{1}{4}a\,\hat{\mathbf{x}} + \tfrac{1}{4}a\,\hat{\mathbf{y}} + \tfrac{3}{4}a\,\hat{\mathbf{z}}$	(16*a*)	Al
$\mathbf{B}_8$	=	$\tfrac{1}{2}\,\mathbf{a}_2$	=	$\tfrac{1}{4}a\,\hat{\mathbf{x}} + \tfrac{3}{4}a\,\hat{\mathbf{y}} + \tfrac{1}{4}a\,\hat{\mathbf{z}}$	(16*a*)	Al
$\mathbf{B}_9$	=	$\tfrac{1}{4}\,\mathbf{a}_1 + \tfrac{3}{8}\,\mathbf{a}_2 + \tfrac{1}{8}\,\mathbf{a}_3$	=	$\tfrac{1}{8}a\,\hat{\mathbf{x}} + \tfrac{1}{4}a\,\hat{\mathbf{z}}$	(24*c*)	Ca
$\mathbf{B}_{10}$	=	$\tfrac{3}{4}\,\mathbf{a}_1 + \tfrac{1}{8}\,\mathbf{a}_2 + \tfrac{3}{8}\,\mathbf{a}_3$	=	$-\tfrac{1}{8}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{4}a\,\hat{\mathbf{z}}$	(24*c*)	Ca
$\mathbf{B}_{11}$	=	$\tfrac{1}{8}\,\mathbf{a}_1 + \tfrac{1}{4}\,\mathbf{a}_2 + \tfrac{3}{8}\,\mathbf{a}_3$	=	$\tfrac{1}{4}a\,\hat{\mathbf{x}} + \tfrac{1}{8}a\,\hat{\mathbf{y}}$	(24*c*)	Ca
$\mathbf{B}_{12}$	=	$\tfrac{3}{8}\,\mathbf{a}_1 + \tfrac{3}{4}\,\mathbf{a}_2 + \tfrac{1}{8}\,\mathbf{a}_3$	=	$\tfrac{1}{4}a\,\hat{\mathbf{x}} - \tfrac{1}{8}a\,\hat{\mathbf{y}} + \tfrac{1}{2}a\,\hat{\mathbf{z}}$	(24*c*)	Ca
$\mathbf{B}_{13}$	=	$\tfrac{3}{8}\,\mathbf{a}_1 + \tfrac{1}{8}\,\mathbf{a}_2 + \tfrac{1}{4}\,\mathbf{a}_3$	=	$\tfrac{1}{4}a\,\hat{\mathbf{y}} + \tfrac{1}{8}a\,\hat{\mathbf{z}}$	(24*c*)	Ca
$\mathbf{B}_{14}$	=	$\tfrac{1}{8}\,\mathbf{a}_1 + \tfrac{3}{8}\,\mathbf{a}_2 + \tfrac{3}{4}\,\mathbf{a}_3$	=	$\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{4}a\,\hat{\mathbf{y}} - \tfrac{1}{8}a\,\hat{\mathbf{z}}$	(24*c*)	Ca
$\mathbf{B}_{15}$	=	$\tfrac{3}{4}\,\mathbf{a}_1 + \tfrac{5}{8}\,\mathbf{a}_2 + \tfrac{7}{8}\,\mathbf{a}_3$	=	$\tfrac{3}{8}a\,\hat{\mathbf{x}} + \tfrac{1}{2}a\,\hat{\mathbf{y}} + \tfrac{1}{4}a\,\hat{\mathbf{z}}$	(24*c*)	Ca
$\mathbf{B}_{16}$	=	$\tfrac{1}{4}\,\mathbf{a}_1 + \tfrac{7}{8}\,\mathbf{a}_2 + \tfrac{5}{8}\,\mathbf{a}_3$	=	$\tfrac{5}{8}a\,\hat{\mathbf{x}} + \tfrac{1}{4}a\,\hat{\mathbf{z}}$	(24*c*)	Ca
$\mathbf{B}_{17}$	=	$\tfrac{7}{8}\,\mathbf{a}_1 + \tfrac{3}{4}\,\mathbf{a}_2 + \tfrac{5}{8}\,\mathbf{a}_3$	=	$\tfrac{1}{4}a\,\hat{\mathbf{x}} + \tfrac{3}{8}a\,\hat{\mathbf{y}} + \tfrac{1}{2}a\,\hat{\mathbf{z}}$	(24*c*)	Ca
$\mathbf{B}_{18}$	=	$\tfrac{5}{8}\,\mathbf{a}_1 + \tfrac{1}{4}\,\mathbf{a}_2 + \tfrac{7}{8}\,\mathbf{a}_3$	=	$\tfrac{1}{4}a\,\hat{\mathbf{x}} + \tfrac{5}{8}a\,\hat{\mathbf{y}}$	(24*c*)	Ca
$\mathbf{B}_{19}$	=	$\tfrac{5}{8}\,\mathbf{a}_1 + \tfrac{7}{8}\,\mathbf{a}_2 + \tfrac{3}{4}\,\mathbf{a}_3$	=	$\tfrac{1}{2}a\,\hat{\mathbf{x}} + \tfrac{1}{4}a\,\hat{\mathbf{y}} + \tfrac{3}{8}a\,\hat{\mathbf{z}}$	(24*c*)	Ca
$\mathbf{B}_{20}$	=	$\tfrac{7}{8}\,\mathbf{a}_1 + \tfrac{5}{8}\,\mathbf{a}_2 + \tfrac{1}{4}\,\mathbf{a}_3$	=	$\tfrac{1}{4}a\,\hat{\mathbf{y}} + \tfrac{5}{8}a\,\hat{\mathbf{z}}$	(24*c*)	Ca
$\mathbf{B}_{21}$	=	$(y_3 + z_3)\,\mathbf{a}_1 + (x_3 + z_3)\,\mathbf{a}_2 + (x_3 + y_3)\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} + y_3a\,\hat{\mathbf{y}} + z_3a\,\hat{\mathbf{z}}$	(96*h*)	H
$\mathbf{B}_{22}$	=	$\left(\tfrac{1}{2} - y_3 + z_3\right)\mathbf{a}_1 + (-x_3 + z_3)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_3 - y_3\right)\mathbf{a}_3$	=	$-x_3a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_3\right)a\,\hat{\mathbf{y}} + z_3a\,\hat{\mathbf{z}}$	(96*h*)	H
$\mathbf{B}_{23}$	=	$(y_3 - z_3)\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_3 - z_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_3 + y_3\right)\mathbf{a}_3$	=	$\left(\tfrac{1}{2} - x_3\right)a\,\hat{\mathbf{x}} + y_3a\,\hat{\mathbf{y}} - z_3a\,\hat{\mathbf{z}}$	(96*h*)	H
$\mathbf{B}_{24}$	=	$\left(\tfrac{1}{2} - y_3 - z_3\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_3 - z_3\right)\mathbf{a}_2 + (x_3 - y_3)\,\mathbf{a}_3$	=	$x_3a\,\hat{\mathbf{x}} - y_3a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_3\right)a\,\hat{\mathbf{z}}$	(96*h*)	H

$$\mathbf{B}_{25} = (x_3+y_3)\,\mathbf{a}_1 + (y_3+z_3)\,\mathbf{a}_2 + (x_3+z_3)\,\mathbf{a}_3 = z_3a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} + y_3a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{26} = \left(\tfrac{1}{2}-x_3-y_3\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_3+z_3\right)\mathbf{a}_2 + (-x_3+z_3)\,\mathbf{a}_3 = z_3a\,\hat{\mathbf{x}} - x_3a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-y_3\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{27} = \left(\tfrac{1}{2}-x_3+y_3\right)\mathbf{a}_1 + (y_3-z_3)\,\mathbf{a}_2 + \left(\tfrac{1}{2}-x_3-z_3\right)\mathbf{a}_3 = -z_3a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-x_3\right)a\,\hat{\mathbf{y}} + y_3a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{28} = (x_3-y_3)\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_3-z_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2}+x_3-z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-z_3\right)a\,\hat{\mathbf{x}} + x_3a\,\hat{\mathbf{y}} - y_3a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{29} = (x_3+z_3)\,\mathbf{a}_1 + (x_3+y_3)\,\mathbf{a}_2 + (y_3+z_3)\,\mathbf{a}_3 = y_3a\,\hat{\mathbf{x}} + z_3a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{30} = (-x_3+z_3)\,\mathbf{a}_1 + \left(\tfrac{1}{2}-x_3-y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-y_3+z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{2}-y_3\right)a\,\hat{\mathbf{x}} + z_3a\,\hat{\mathbf{y}} - x_3a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{31} = \left(\tfrac{1}{2}-x_3-z_3\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-x_3+y_3\right)\mathbf{a}_2 + (y_3-z_3)\,\mathbf{a}_3 = y_3a\,\hat{\mathbf{x}} - z_3a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2}-x_3\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{32} = \left(\tfrac{1}{2}+x_3-z_3\right)\mathbf{a}_1 + (x_3-y_3)\,\mathbf{a}_2 + \left(\tfrac{1}{2}-y_3-z_3\right)\mathbf{a}_3 = -y_3a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2}-z_3\right)a\,\hat{\mathbf{y}} + x_3a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{33} = \left(\tfrac{1}{2}+x_3-z_3\right)\mathbf{a}_1 + (y_3-z_3)\,\mathbf{a}_2 + (x_3+y_3)\,\mathbf{a}_3 = \left(\tfrac{3}{4}+y_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4}+x_3\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4}-z_3\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{34} = \left(\tfrac{1}{2}-x_3-z_3\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-y_3-z_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-x_3-y_3\right)\mathbf{a}_3 = \left(\tfrac{1}{4}-y_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4}-x_3\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4}-z_3\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{35} = (-x_3+z_3)\,\mathbf{a}_1 + (y_3+z_3)\,\mathbf{a}_2 + \left(\tfrac{1}{2}-x_3+y_3\right)\mathbf{a}_3 = \left(\tfrac{1}{4}+y_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4}-x_3\right)a\,\hat{\mathbf{y}} + \left(\tfrac{3}{4}+z_3\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{36} = (x_3+z_3)\,\mathbf{a}_1 + \left(\tfrac{1}{2}-y_3+z_3\right)\mathbf{a}_2 + (x_3-y_3)\,\mathbf{a}_3 = \left(\tfrac{1}{4}-y_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{3}{4}+x_3\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4}+z_3\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{37} = \left(\tfrac{1}{2}-y_3+z_3\right)\mathbf{a}_1 + (x_3-y_3)\,\mathbf{a}_2 + (x_3+z_3)\,\mathbf{a}_3 = \left(\tfrac{3}{4}+x_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4}+z_3\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4}-y_3\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{38} = (y_3+z_3)\,\mathbf{a}_1 + \left(\tfrac{1}{2}-x_3+y_3\right)\mathbf{a}_2 + (-x_3+z_3)\,\mathbf{a}_3 = \left(\tfrac{1}{4}-x_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{3}{4}+z_3\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4}+y_3\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{39} = \left(\tfrac{1}{2}-y_3-z_3\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-x_3-y_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-x_3-z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{4}-x_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4}-z_3\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4}-y_3\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{40} = (y_3-z_3)\,\mathbf{a}_1 + (x_3+y_3)\,\mathbf{a}_2 + \left(\tfrac{1}{2}+x_3-z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{4}+x_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4}-z_3\right)a\,\hat{\mathbf{y}} + \left(\tfrac{3}{4}+y_3\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{41} = \left(\tfrac{1}{2}-x_3+y_3\right)\mathbf{a}_1 + (-x_3+z_3)\,\mathbf{a}_2 + (y_3+z_3)\,\mathbf{a}_3 = \left(\tfrac{3}{4}+z_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4}+y_3\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4}-x_3\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{42} = (x_3-y_3)\,\mathbf{a}_1 + (x_3+z_3)\,\mathbf{a}_2 + \left(\tfrac{1}{2}-y_3+z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{4}+z_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4}-y_3\right)a\,\hat{\mathbf{y}} + \left(\tfrac{3}{4}+x_3\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{43} = (x_3+y_3)\,\mathbf{a}_1 + \left(\tfrac{1}{2}+x_3-z_3\right)\mathbf{a}_2 + (y_3-z_3)\,\mathbf{a}_3 = \left(\tfrac{1}{4}-z_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{3}{4}+y_3\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4}+x_3\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{44} = \left(\tfrac{1}{2}-x_3-y_3\right)\mathbf{a}_1 + \left(\tfrac{1}{2}-x_3-z_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2}-y_3-z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{4}-z_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4}-y_3\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4}-x_3\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{45} = (-y_3-z_3)\,\mathbf{a}_1 + (-x_3-z_3)\,\mathbf{a}_2 + (-x_3-y_3)\,\mathbf{a}_3 = -x_3a\,\hat{\mathbf{x}} - y_3a\,\hat{\mathbf{y}} - z_3a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{46} = \left(\tfrac{1}{2}+y_3-z_3\right)\mathbf{a}_1+(x_3-z_3)\,\mathbf{a}_2+\left(\tfrac{1}{2}+x_3+y_3\right)\mathbf{a}_3 = x_3a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+y_3\right)a\,\hat{\mathbf{y}}-z_3a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{47} = (-y_3+z_3)\,\mathbf{a}_1+\left(\tfrac{1}{2}+x_3+z_3\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_3-y_3\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+x_3\right)a\,\hat{\mathbf{x}}-y_3a\,\hat{\mathbf{y}}+z_3a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{48} = \left(\tfrac{1}{2}+y_3+z_3\right)\mathbf{a}_1+\left(\tfrac{1}{2}-x_3+z_3\right)\mathbf{a}_2+(-x_3+y_3)\,\mathbf{a}_3 = -x_3a\,\hat{\mathbf{x}}+y_3a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+z_3\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{49} = (-x_3-y_3)\,\mathbf{a}_1+(-y_3-z_3)\,\mathbf{a}_2+(-x_3-z_3)\,\mathbf{a}_3 = -z_3a\,\hat{\mathbf{x}}-x_3a\,\hat{\mathbf{y}}-y_3a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{50} = \left(\tfrac{1}{2}+x_3+y_3\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_3-z_3\right)\mathbf{a}_2+(x_3-z_3)\,\mathbf{a}_3 = -z_3a\,\hat{\mathbf{x}}+x_3a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+y_3\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{51} = \left(\tfrac{1}{2}+x_3-y_3\right)\mathbf{a}_1+(-y_3+z_3)\,\mathbf{a}_2+\left(\tfrac{1}{2}+x_3+z_3\right)\mathbf{a}_3 = z_3a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+x_3\right)a\,\hat{\mathbf{y}}-y_3a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{52} = (-x_3+y_3)\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_3+z_3\right)\mathbf{a}_2+\left(\tfrac{1}{2}-x_3+z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+z_3\right)a\,\hat{\mathbf{x}}-x_3a\,\hat{\mathbf{y}}+y_3a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{53} = (-x_3-z_3)\,\mathbf{a}_1+(-x_3-y_3)\,\mathbf{a}_2+(-y_3-z_3)\,\mathbf{a}_3 = -y_3a\,\hat{\mathbf{x}}-z_3a\,\hat{\mathbf{y}}-x_3a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{54} = (x_3-z_3)\,\mathbf{a}_1+\left(\tfrac{1}{2}+x_3+y_3\right)\mathbf{a}_2+\left(\tfrac{1}{2}+y_3-z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{2}+y_3\right)a\,\hat{\mathbf{x}}-z_3a\,\hat{\mathbf{y}}+x_3a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{55} = \left(\tfrac{1}{2}+x_3+z_3\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_3-y_3\right)\mathbf{a}_2+(-y_3+z_3)\,\mathbf{a}_3 = -y_3a\,\hat{\mathbf{x}}+z_3a\,\hat{\mathbf{y}}+\left(\tfrac{1}{2}+x_3\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{56} = \left(\tfrac{1}{2}-x_3+z_3\right)\mathbf{a}_1+(-x_3+y_3)\,\mathbf{a}_2+\left(\tfrac{1}{2}+y_3+z_3\right)\mathbf{a}_3 = y_3a\,\hat{\mathbf{x}}+\left(\tfrac{1}{2}+z_3\right)a\,\hat{\mathbf{y}}-x_3a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{57} = \left(\tfrac{1}{2}-x_3+z_3\right)\mathbf{a}_1+(-y_3+z_3)\,\mathbf{a}_2+(-x_3-y_3)\,\mathbf{a}_3 = -\left(y_3+\tfrac{1}{4}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{4}-x_3\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{4}+z_3\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{58} = \left(\tfrac{1}{2}+x_3+z_3\right)\mathbf{a}_1+\left(\tfrac{1}{2}+y_3+z_3\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_3+y_3\right)\mathbf{a}_3 = \left(\tfrac{1}{4}+y_3\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{4}+x_3\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{4}+z_3\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{59} = (x_3-z_3)\,\mathbf{a}_1+(-y_3-z_3)\,\mathbf{a}_2+\left(\tfrac{1}{2}+x_3-y_3\right)\mathbf{a}_3 = \left(\tfrac{1}{4}-y_3\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{4}+x_3\right)a\,\hat{\mathbf{y}}-\left(z_3+\tfrac{1}{4}\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{60} = (-x_3-z_3)\,\mathbf{a}_1+\left(\tfrac{1}{2}+y_3-z_3\right)\mathbf{a}_2+(-x_3+y_3)\,\mathbf{a}_3 = \left(\tfrac{1}{4}+y_3\right)a\,\hat{\mathbf{x}}-\left(x_3+\tfrac{1}{4}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{4}-z_3\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{61} = \left(\tfrac{1}{2}+y_3-z_3\right)\mathbf{a}_1+(-x_3+y_3)\,\mathbf{a}_2+(-x_3-z_3)\,\mathbf{a}_3 = -\left(x_3+\tfrac{1}{4}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{4}-z_3\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{4}+y_3\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{62} = (-y_3-z_3)\,\mathbf{a}_1+\left(\tfrac{1}{2}+x_3-y_3\right)\mathbf{a}_2+(x_3-z_3)\,\mathbf{a}_3 = \left(\tfrac{1}{4}+x_3\right)a\,\hat{\mathbf{x}}-\left(z_3+\tfrac{1}{4}\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{4}-y_3\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{63} = \left(\tfrac{1}{2}+y_3+z_3\right)\mathbf{a}_1+\left(\tfrac{1}{2}+x_3+y_3\right)\mathbf{a}_2+\left(\tfrac{1}{2}+x_3+z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{4}+x_3\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{4}+z_3\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{4}+y_3\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{64} = (-y_3+z_3)\,\mathbf{a}_1+(-x_3-y_3)\,\mathbf{a}_2+\left(\tfrac{1}{2}-x_3+z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{4}-x_3\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{4}+z_3\right)a\,\hat{\mathbf{y}}-\left(y_3+\tfrac{1}{4}\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{65} = \left(\tfrac{1}{2}+x_3-y_3\right)\mathbf{a}_1+(x_3-z_3)\,\mathbf{a}_2+(-y_3-z_3)\,\mathbf{a}_3 = -\left(z_3+\tfrac{1}{4}\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{4}-y_3\right)a\,\hat{\mathbf{y}}+\left(\tfrac{1}{4}+x_3\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{66} = (-x_3+y_3)\,\mathbf{a}_1+(-x_3-z_3)\,\mathbf{a}_2+\left(\tfrac{1}{2}+y_3-z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{4}-z_3\right)a\,\hat{\mathbf{x}}+\left(\tfrac{1}{4}+y_3\right)a\,\hat{\mathbf{y}}-\left(x_3+\tfrac{1}{4}\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{67} = (-x_3 - y_3)\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_3 + z_3\right)\mathbf{a}_2 + (-y_3 + z_3)\,\mathbf{a}_3 = \left(\tfrac{1}{4} + z_3\right)a\,\hat{\mathbf{x}} - \left(y_3 + \tfrac{1}{4}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - x_3\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{68} = \left(\tfrac{1}{2} + x_3 + y_3\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_3 + z_3\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + y_3 + z_3\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + z_3\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + y_3\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + x_3\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{H}$$

$$\mathbf{B}_{69} = (y_4 + z_4)\,\mathbf{a}_1 + (x_4 + z_4)\,\mathbf{a}_2 + (x_4 + y_4)\,\mathbf{a}_3 = x_4 a\,\hat{\mathbf{x}} + y_4 a\,\hat{\mathbf{y}} + z_4 a\,\hat{\mathbf{z}} \quad (96h) \quad \text{O}$$

$$\mathbf{B}_{70} = \left(\tfrac{1}{2} - y_4 + z_4\right)\mathbf{a}_1 + (-x_4 + z_4)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_4 - y_4\right)\mathbf{a}_3 = -x_4 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - y_4\right)a\,\hat{\mathbf{y}} + z_4 a\,\hat{\mathbf{z}} \quad (96h) \quad \text{O}$$

$$\mathbf{B}_{71} = (y_4 - z_4)\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_4 - z_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_4 + y_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - x_4\right)a\,\hat{\mathbf{x}} + y_4 a\,\hat{\mathbf{y}} - z_4 a\,\hat{\mathbf{z}} \quad (96h) \quad \text{O}$$

$$\mathbf{B}_{72} = \left(\tfrac{1}{2} - y_4 - z_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_4 - z_4\right)\mathbf{a}_2 + (x_4 - y_4)\,\mathbf{a}_3 = x_4 a\,\hat{\mathbf{x}} - y_4 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - z_4\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{O}$$

$$\mathbf{B}_{73} = (x_4 + y_4)\,\mathbf{a}_1 + (y_4 + z_4)\,\mathbf{a}_2 + (x_4 + z_4)\,\mathbf{a}_3 = z_4 a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}} + y_4 a\,\hat{\mathbf{z}} \quad (96h) \quad \text{O}$$

$$\mathbf{B}_{74} = \left(\tfrac{1}{2} - x_4 - y_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_4 + z_4\right)\mathbf{a}_2 + (-x_4 + z_4)\,\mathbf{a}_3 = z_4 a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - y_4\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{O}$$

$$\mathbf{B}_{75} = \left(\tfrac{1}{2} - x_4 + y_4\right)\mathbf{a}_1 + (y_4 - z_4)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_4 - z_4\right)\mathbf{a}_3 = -z_4 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - x_4\right)a\,\hat{\mathbf{y}} + y_4 a\,\hat{\mathbf{z}} \quad (96h) \quad \text{O}$$

$$\mathbf{B}_{76} = (x_4 - y_4)\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_4 - z_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_4 - z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - z_4\right)a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}} - y_4 a\,\hat{\mathbf{z}} \quad (96h) \quad \text{O}$$

$$\mathbf{B}_{77} = (x_4 + z_4)\,\mathbf{a}_1 + (x_4 + y_4)\,\mathbf{a}_2 + (y_4 + z_4)\,\mathbf{a}_3 = y_4 a\,\hat{\mathbf{x}} + z_4 a\,\hat{\mathbf{y}} + x_4 a\,\hat{\mathbf{z}} \quad (96h) \quad \text{O}$$

$$\mathbf{B}_{78} = (-x_4 + z_4)\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_4 - y_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - y_4 + z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2} - y_4\right)a\,\hat{\mathbf{x}} + z_4 a\,\hat{\mathbf{y}} - x_4 a\,\hat{\mathbf{z}} \quad (96h) \quad \text{O}$$

$$\mathbf{B}_{79} = \left(\tfrac{1}{2} - x_4 - z_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_4 + y_4\right)\mathbf{a}_2 + (y_4 - z_4)\,\mathbf{a}_3 = y_4 a\,\hat{\mathbf{x}} - z_4 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} - x_4\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{O}$$

$$\mathbf{B}_{80} = \left(\tfrac{1}{2} + x_4 - z_4\right)\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - y_4 - z_4\right)\mathbf{a}_3 = -y_4 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} - z_4\right)a\,\hat{\mathbf{y}} + x_4 a\,\hat{\mathbf{z}} \quad (96h) \quad \text{O}$$

$$\mathbf{B}_{81} = \left(\tfrac{1}{2} + x_4 - z_4\right)\mathbf{a}_1 + (y_4 - z_4)\,\mathbf{a}_2 + (x_4 + y_4)\,\mathbf{a}_3 = \left(\tfrac{3}{4} + y_4\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - z_4\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{O}$$

$$\mathbf{B}_{82} = \left(\tfrac{1}{2} - x_4 - z_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - y_4 - z_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_4 - y_4\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - y_4\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - z_4\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{O}$$

$$\mathbf{B}_{83} = (-x_4 + z_4)\,\mathbf{a}_1 + (y_4 + z_4)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - x_4 + y_4\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + y_4\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{3}{4} + z_4\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{O}$$

$$\mathbf{B}_{84} = (x_4 + z_4)\,\mathbf{a}_1 + \left(\tfrac{1}{2} - y_4 + z_4\right)\mathbf{a}_2 + (x_4 - y_4)\,\mathbf{a}_3 = \left(\tfrac{1}{4} - y_4\right)a\,\hat{\mathbf{x}} + \left(\tfrac{3}{4} + x_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_4\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{O}$$

$$\mathbf{B}_{85} = \left(\tfrac{1}{2} - y_4 + z_4\right)\mathbf{a}_1 + (x_4 - y_4)\,\mathbf{a}_2 + (x_4 + z_4)\,\mathbf{a}_3 = \left(\tfrac{3}{4} + x_4\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + z_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - y_4\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{O}$$

$$\mathbf{B}_{86} = (y_4 + z_4)\,\mathbf{a}_1 + \left(\tfrac{1}{2} - x_4 + y_4\right)\mathbf{a}_2 + (-x_4 + z_4)\,\mathbf{a}_3 = \left(\tfrac{1}{4} - x_4\right)a\,\hat{\mathbf{x}} + \left(\tfrac{3}{4} + z_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + y_4\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{O}$$

$$\mathbf{B}_{87} = \left(\tfrac{1}{2} - y_4 - z_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_4 - y_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_4 - z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - x_4\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - z_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - y_4\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \text{O}$$

$$\mathbf{B}_{88} = (y_4 - z_4)\,\mathbf{a}_1 + (x_4 + y_4)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_4 - z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + x_4\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - z_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{3}{4} + y_4\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \mathrm{O}$$

$$\mathbf{B}_{89} = \left(\tfrac{1}{2} - x_4 + y_4\right)\mathbf{a}_1 + (-x_4 + z_4)\,\mathbf{a}_2 + (y_4 + z_4)\,\mathbf{a}_3 = \left(\tfrac{3}{4} + z_4\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + y_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - x_4\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \mathrm{O}$$

$$\mathbf{B}_{90} = (x_4 - y_4)\,\mathbf{a}_1 + (x_4 + z_4)\,\mathbf{a}_2 + \left(\tfrac{1}{2} - y_4 + z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + z_4\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - y_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{3}{4} + x_4\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \mathrm{O}$$

$$\mathbf{B}_{91} = (x_4 + y_4)\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_4 - z_4\right)\mathbf{a}_2 + (y_4 - z_4)\,\mathbf{a}_3 = \left(\tfrac{1}{4} - z_4\right)a\,\hat{\mathbf{x}} + \left(\tfrac{3}{4} + y_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + x_4\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \mathrm{O}$$

$$\mathbf{B}_{92} = \left(\tfrac{1}{2} - x_4 - y_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_4 - z_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - y_4 - z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - z_4\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - y_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - x_4\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \mathrm{O}$$

$$\mathbf{B}_{93} = (-y_4 - z_4)\,\mathbf{a}_1 + (-x_4 - z_4)\,\mathbf{a}_2 + (-x_4 - y_4)\,\mathbf{a}_3 = -x_4 a\,\hat{\mathbf{x}} - y_4 a\,\hat{\mathbf{y}} - z_4 a\,\hat{\mathbf{z}} \quad (96h) \quad \mathrm{O}$$

$$\mathbf{B}_{94} = \left(\tfrac{1}{2} + y_4 - z_4\right)\mathbf{a}_1 + (x_4 - z_4)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_4 + y_4\right)\mathbf{a}_3 = x_4 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + y_4\right)a\,\hat{\mathbf{y}} - z_4 a\,\hat{\mathbf{z}} \quad (96h) \quad \mathrm{O}$$

$$\mathbf{B}_{95} = (-y_4 + z_4)\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_4 + z_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_4 - y_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + x_4\right)a\,\hat{\mathbf{x}} - y_4 a\,\hat{\mathbf{y}} + z_4 a\,\hat{\mathbf{z}} \quad (96h) \quad \mathrm{O}$$

$$\mathbf{B}_{96} = \left(\tfrac{1}{2} + y_4 + z_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2} - x_4 + z_4\right)\mathbf{a}_2 + (-x_4 + y_4)\,\mathbf{a}_3 = -x_4 a\,\hat{\mathbf{x}} + y_4 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + z_4\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \mathrm{O}$$

$$\mathbf{B}_{97} = (-x_4 - y_4)\,\mathbf{a}_1 + (-y_4 - z_4)\,\mathbf{a}_2 + (-x_4 - z_4)\,\mathbf{a}_3 = -z_4 a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}} - y_4 a\,\hat{\mathbf{z}} \quad (96h) \quad \mathrm{O}$$

$$\mathbf{B}_{98} = \left(\tfrac{1}{2} + x_4 + y_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_4 - z_4\right)\mathbf{a}_2 + (x_4 - z_4)\,\mathbf{a}_3 = -z_4 a\,\hat{\mathbf{x}} + x_4 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + y_4\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \mathrm{O}$$

$$\mathbf{B}_{99} = \left(\tfrac{1}{2} + x_4 - y_4\right)\mathbf{a}_1 + (-y_4 + z_4)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_4 + z_4\right)\mathbf{a}_3 = z_4 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + x_4\right)a\,\hat{\mathbf{y}} - y_4 a\,\hat{\mathbf{z}} \quad (96h) \quad \mathrm{O}$$

$$\mathbf{B}_{100} = (-x_4 + y_4)\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_4 + z_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} - x_4 + z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + z_4\right)a\,\hat{\mathbf{x}} - x_4 a\,\hat{\mathbf{y}} + y_4 a\,\hat{\mathbf{z}} \quad (96h) \quad \mathrm{O}$$

$$\mathbf{B}_{101} = (-x_4 - z_4)\,\mathbf{a}_1 + (-x_4 - y_4)\,\mathbf{a}_2 + (-y_4 - z_4)\,\mathbf{a}_3 = -y_4 a\,\hat{\mathbf{x}} - z_4 a\,\hat{\mathbf{y}} - x_4 a\,\hat{\mathbf{z}} \quad (96h) \quad \mathrm{O}$$

$$\mathbf{B}_{102} = (x_4 - z_4)\,\mathbf{a}_1 + \left(\tfrac{1}{2} + x_4 + y_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + y_4 - z_4\right)\mathbf{a}_3 = \left(\tfrac{1}{2} + y_4\right)a\,\hat{\mathbf{x}} - z_4 a\,\hat{\mathbf{y}} + x_4 a\,\hat{\mathbf{z}} \quad (96h) \quad \mathrm{O}$$

$$\mathbf{B}_{103} = \left(\tfrac{1}{2} + x_4 + z_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + x_4 - y_4\right)\mathbf{a}_2 + (-y_4 + z_4)\,\mathbf{a}_3 = -y_4 a\,\hat{\mathbf{x}} + z_4 a\,\hat{\mathbf{y}} + \left(\tfrac{1}{2} + x_4\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \mathrm{O}$$

$$\mathbf{B}_{104} = \left(\tfrac{1}{2} - x_4 + z_4\right)\mathbf{a}_1 + (-x_4 + y_4)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + y_4 + z_4\right)\mathbf{a}_3 = y_4 a\,\hat{\mathbf{x}} + \left(\tfrac{1}{2} + z_4\right)a\,\hat{\mathbf{y}} - x_4 a\,\hat{\mathbf{z}} \quad (96h) \quad \mathrm{O}$$

$$\mathbf{B}_{105} = \left(\tfrac{1}{2} - x_4 + z_4\right)\mathbf{a}_1 + (-y_4 + z_4)\,\mathbf{a}_2 + (-x_4 - y_4)\,\mathbf{a}_3 = -\left(y_4 + \tfrac{1}{4}\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} - x_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_4\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \mathrm{O}$$

$$\mathbf{B}_{106} = \left(\tfrac{1}{2} + x_4 + z_4\right)\mathbf{a}_1 + \left(\tfrac{1}{2} + y_4 + z_4\right)\mathbf{a}_2 + \left(\tfrac{1}{2} + x_4 + y_4\right)\mathbf{a}_3 = \left(\tfrac{1}{4} + y_4\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_4\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} + z_4\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \mathrm{O}$$

$$\mathbf{B}_{107} = (x_4 - z_4)\,\mathbf{a}_1 + (-y_4 - z_4)\,\mathbf{a}_2 + \left(\tfrac{1}{2} + x_4 - y_4\right)\mathbf{a}_3 = \left(\tfrac{1}{4} - y_4\right)a\,\hat{\mathbf{x}} + \left(\tfrac{1}{4} + x_4\right)a\,\hat{\mathbf{y}} - \left(z_4 + \tfrac{1}{4}\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \mathrm{O}$$

$$\mathbf{B}_{108} = (-x_4 - z_4)\,\mathbf{a}_1 + \left(\tfrac{1}{2} + y_4 - z_4\right)\mathbf{a}_2 + (-x_4 + y_4)\,\mathbf{a}_3 = \left(\tfrac{1}{4} + y_4\right)a\,\hat{\mathbf{x}} - \left(x_4 + \tfrac{1}{4}\right)a\,\hat{\mathbf{y}} + \left(\tfrac{1}{4} - z_4\right)a\,\hat{\mathbf{z}} \quad (96h) \quad \mathrm{O}$$

$\mathbf{B}_{109}$	=	$\left(\frac{1}{2}+y_4-z_4\right)\mathbf{a}_1+(-x_4+y_4)\,\mathbf{a}_2+(-x_4-z_4)\,\mathbf{a}_3$	=	$-\left(x_4+\frac{1}{4}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-z_4\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+y_4\right)a\,\hat{\mathbf{z}}$	(96*h*)	O
$\mathbf{B}_{110}$	=	$(-y_4-z_4)\,\mathbf{a}_1+\left(\frac{1}{2}+x_4-y_4\right)\mathbf{a}_2+(x_4-z_4)\,\mathbf{a}_3$	=	$\left(\frac{1}{4}+x_4\right)a\,\hat{\mathbf{x}}-\left(z_4+\frac{1}{4}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-y_4\right)a\,\hat{\mathbf{z}}$	(96*h*)	O
$\mathbf{B}_{111}$	=	$\left(\frac{1}{2}+y_4+z_4\right)\mathbf{a}_1+\left(\frac{1}{2}+x_4+y_4\right)\mathbf{a}_2+\left(\frac{1}{2}+x_4+z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}+x_4\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+z_4\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+y_4\right)a\,\hat{\mathbf{z}}$	(96*h*)	O
$\mathbf{B}_{112}$	=	$(-y_4+z_4)\,\mathbf{a}_1+(-x_4-y_4)\,\mathbf{a}_2+\left(\frac{1}{2}-x_4+z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}-x_4\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+z_4\right)a\,\hat{\mathbf{y}}-\left(y_4+\frac{1}{4}\right)a\,\hat{\mathbf{z}}$	(96*h*)	O
$\mathbf{B}_{113}$	=	$\left(\frac{1}{2}+x_4-y_4\right)\mathbf{a}_1+(x_4-z_4)\,\mathbf{a}_2+(-y_4-z_4)\,\mathbf{a}_3$	=	$-\left(z_4+\frac{1}{4}\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}-y_4\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+x_4\right)a\,\hat{\mathbf{z}}$	(96*h*)	O
$\mathbf{B}_{114}$	=	$(-x_4+y_4)\,\mathbf{a}_1+(-x_4-z_4)\,\mathbf{a}_2+\left(\frac{1}{2}+y_4-z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}-z_4\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+y_4\right)a\,\hat{\mathbf{y}}-\left(x_4+\frac{1}{4}\right)a\,\hat{\mathbf{z}}$	(96*h*)	O
$\mathbf{B}_{115}$	=	$(-x_4-y_4)\,\mathbf{a}_1+\left(\frac{1}{2}-x_4+z_4\right)\mathbf{a}_2+(-y_4+z_4)\,\mathbf{a}_3$	=	$\left(\frac{1}{4}+z_4\right)a\,\hat{\mathbf{x}}-\left(y_4+\frac{1}{4}\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}-x_4\right)a\,\hat{\mathbf{z}}$	(96*h*)	O
$\mathbf{B}_{116}$	=	$\left(\frac{1}{2}+x_4+y_4\right)\mathbf{a}_1+\left(\frac{1}{2}+x_4+z_4\right)\mathbf{a}_2+\left(\frac{1}{2}+y_4+z_4\right)\mathbf{a}_3$	=	$\left(\frac{1}{4}+z_4\right)a\,\hat{\mathbf{x}}+\left(\frac{1}{4}+y_4\right)a\,\hat{\mathbf{y}}+\left(\frac{1}{4}+x_4\right)a\,\hat{\mathbf{z}}$	(96*h*)	O

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- H. Bartl, *Tricalciumaluminathexahydrat, $Ca_3[Al(OH)_6]_2$, Bindungslängen und -valenzen aus Röntgeneinkristallmessungen*, Fresen. Z. Anal. Chem. **324**, 124–126 (1986), doi:10.1007/BF00473351.
- C. Gottfried and F. Schossberger, eds., *Strukturbericht Band III 1933-1935* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).
- E. Brandenberger, *Kristallstrukturelle Untersuchungen an Calciumaluminathydraten*, Schweiz. Mineral. Petrog. Mitt. **13** (1933).

Geometry files:

- CIF: pp. 1836
- POSCAR: pp. 1837

CIF and POSCAR Files

$NaC_5H_{11}O_8S$: A5B11CD8E_aP26_1_5a_11a_a_8a_a - CIF

```
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_chemical_formula_sum 'C5 H11 Na O8 S'

loop_
_publ_author_name
 'A. H. Haines'
 'D. L. Hughes'
_journal_name_full_name
;
 Acta Crystallographica Section E: Crystallographic Communications
;
_journal_volume 72
_journal_year 2016
_journal_page_first 628
_journal_page_last 631
_publ_Section_title
;
 Crystal structure of sodium (1S)-D-lyxit-1-yl sulfonate
;

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      ↪ 0.552,0.42,0.642,0.781,0.158,-0.018,0.345,0.159,0.144,0.035,
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C2 C 1 a 0.29660 0.23360 0.74080 1.00000
C3 C 1 a 0.05620 0.30940 0.86872 1.00000
C4 C 1 a 0.16940 0.41620 -0.01970 1.00000
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H1 H 1 a 0.15800 -0.02200 0.67900 1.00000
H2 H 1 a 0.45000 0.11700 0.79400 1.00000
H3 H 1 a -0.09600 0.43300 0.81600 1.00000
H4 H 1 a 0.24600 0.54800 -0.08100 1.00000
H5 H 1 a 0.76400 0.63200 0.03700 1.00000
H6 H 1 a 0.86700 0.40900 0.17100 1.00000
H7 H 1 a 0.89300 0.24800 0.50300 1.00000
H8 H 1 a 0.55200 0.42000 0.64200 1.00000
H9 H 1 a 0.78100 0.15800 -0.01800 1.00000
H10 H 1 a 0.34500 0.15900 0.14400 1.00000
H11 H 1 a 0.03500 0.79900 0.11300 1.00000
Na1 Na 1 a 0.14247 0.62758 0.42059 1.00000
O1 O 1 a -0.04310 0.29760 0.55444 1.00000
O2 O 1 a 0.55970 0.24470 0.39414 1.00000
O3 O 1 a 0.70800 0.83690 0.55456 1.00000
O4 O 1 a 0.33050 -0.05200 0.37321 1.00000
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O7 O 1 a 0.40560 0.23550 0.05675 1.00000
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S1 S 1 a 0.47392 0.02789 0.47548 1.00000
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$NaC_5H_{11}O_8S$: A5B11CD8E_aP26_1_5a_11a_a_8a_a - POSCAR

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      ↪ y9,z9,x10,y10,z10,x11,y11,z11,x12,y12,z12,x13,y13,z13,x14,y14,
      ↪ z14,x15,y15,z15,x16,y16,z16,x17,y17,z17,x18,y18,z18,x19,y19,z19
      ↪ ,x20,y20,z20,x21,y21,z21,x22,y22,z22,x23,y23,z23,x24,y24,z24,
      ↪ x25,y25,z25,x26,y26,z26 --params=4.8558,1.20466246551,
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      ↪ ,0.433,0.816,0.246,0.548,-0.081,0.764,0.632,0.037,0.867,0.409,
      ↪ 0.171,0.893,0.248,0.503,0.552,0.42,0.642,0.781,0.158,-0.018,
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      ↪ Crystallogr. E 72, 628-631 (2016)
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  -0.03760000000000    0.10350000000000   -0.05381000000000     O     (1a)
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$Ni(NO_3)_2(H_2O)_6$: A12B2CD12_aP54_2_12i_2i_i_12i - CIF

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 'M. T. Camellini'
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_journal_volume 27
_journal_year 1971
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;
 The crystal structures of nitrates of divalent hexaaquocations. III.
      ↪ Hexaaquonickel nitrate
;

# Found in Reaction of [Ni(H$_{2}$O)$_{6}$](NO$_{3}$)$_{2}$ with gaseous
      ↪ NH$_{3}$; crystal growth {\em via} in-situ solvation, 2015

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      ↪ 0.4207,0.873,0.6648,0.3707,0.6081,0.19,0.09,0.2325,0.0606,
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H3 H 2 i -0.02330 0.12830 0.35500 1.00000
H4 H 2 i 0.20200 0.13500 0.44200 1.00000
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H9 H 2 i 0.40000 0.35500 0.39170 1.00000
H10 H 2 i 0.47040 0.34510 0.18170 1.00000
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N2 N 2 i 0.37070 0.60810 0.19000 1.00000
Ni1 Ni 2 i 0.09000 0.23250 0.06060 1.00000
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O4 O 2 i 0.53860 0.62830 0.28040 1.00000
O5 O 2 i 0.25870 0.60350 0.31580 1.00000
O6 O 2 i 0.31060 0.59040 -0.04240 1.00000
O7 O 2 i 0.01350 0.36550 0.26940 1.00000
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O9 O 2 i 0.17000 0.10380 0.85330 1.00000
O10 O 2 i 0.07730 0.31490 0.77600 1.00000
O11 O 2 i 0.36140 0.34960 0.21790 1.00000
O12 O 2 i 0.81660 0.12620 0.90310 1.00000
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$Ni(NO_3)_2(H_2O)_6$: A12B2CD12_aP54_2_12i_2i_i_12i - POSCAR

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      ↪ y9,z9,x10,y10,z10,x11,y11,z11,x12,y12,z12,x13,y13,z13,x14,y14,
      ↪ z14,x15,y15,z15,x16,y16,z16,x17,y17,z17,x18,y18,z18,x19,y19,z19
      ↪ ,x20,y20,z20,x21,y21,z21,x22,y22,z22,x23,y23,z23,x24,y24,z24,
      ↪ x25,y25,z25,x26,y26,z26,x27,y27,z27 --params=7.694,
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      ↪ 0.442,0.3033,0.1153,-0.095,0.07,0.04,0.7433,-0.0353,0.2655,
      ↪ 0.6483,0.1356,0.3861,0.7517,0.4,0.355,0.3917,0.4704,0.3451,
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      ↪ 0.6648,0.3707,0.6081,0.19,0.09,0.2325,0.0606,0.4696,0.8648,
      ↪ 0.8845,0.5323,0.8766,0.5422,0.2611,0.8757,0.5844,0.5386,0.6283,
      ↪ 0.2804,0.2587,0.6035,0.3158,0.3106,0.5904,-0.0424,0.0135,0.3655
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      ↪ 0.776,0.3614,0.3496,0.2179,0.8166,0.1262,0.9031 & P-1 C_{i}^{1}
      ↪ #2 (i^27) & aP54 & None & H12N2NiO12 & H12N2NiO12 & F. Bigoli
      ↪ et al., Acta Crystallogr. Sect. B Struct. Sci. 27, 1427-1434 (
      ↪ 1971)
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$Co_2B_2O_5$: A2B2C5_aP18_2_2i_2i_5i - CIF

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_journal_year 1950
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;
 The Crystal Structure of Cobaltpyroborate
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B1 B 2 i 0.67000 0.67000 0.34000 1.00000
B2 B 2 i 0.36000 0.88000 0.17000 1.00000
Co1 Co 2 i 0.74300 0.21300 0.36000 1.00000
Co2 Co 2 i 0.24500 0.37400 0.10000 1.00000
O1 O 2 i 0.24400 0.70900 0.05400 1.00000
O2 O 2 i 0.20800 0.09200 0.19000 1.00000
```

```
O3  O  2  i  0.73500  0.47600  0.24900  1.00000
O4  O  2  i  0.56200  0.84200  0.28800  1.00000
O5  O  2  i  0.77000  0.69800  0.48900  1.00000
```

$Co_2B_2O_5$: A2B2C5_aP18_2_2i_2i_5i - POSCAR

```
A2B2C5_aP18_2_2i_2i_5i & a,b/a,c/a,alpha,beta,gamma,x1,y1,z1,x2,y2,z2,x3
    ↪ ,y3,z3,x4,y4,z4,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9 --
    ↪ params=3.16,1.87974683544,2.83227848101,103.9,91.0,92.0,0.67,
    ↪ 0.67,0.34,0.36,0.88,0.17,0.743,0.213,0.36,0.245,0.374,0.1,0.244
    ↪ ,0.709,0.054,0.208,0.092,0.19,0.735,0.476,0.249,0.562,0.842,
    ↪ 0.288,0.77,0.698,0.489 & P-1 C_{i}^{1} #2 (i^9) & aP18 & None &
    ↪  B2Co2O5 & B2Co2O5 & S. V. Berger, Acta Chem. Scand. 4,
    ↪ 1054-1065 (1950)
   1.00000000000000
   3.16000000000000     0.00000000000000     0.00000000000000
  -0.20730301041286     5.93638151249343     0.00000000000000
  -0.15619903761369    -2.15680611598055     8.68483098504038
     B     Co     O
     4      4     10
Direct
   0.67000000000000     0.67000000000000     0.34000000000000    B    (2i)
  -0.67000000000000    -0.67000000000000    -0.34000000000000    B    (2i)
   0.36000000000000     0.88000000000000     0.17000000000000    B    (2i)
  -0.36000000000000    -0.88000000000000    -0.17000000000000    B    (2i)
   0.74300000000000     0.21300000000000     0.36000000000000    Co   (2i)
  -0.74300000000000    -0.21300000000000    -0.36000000000000    Co   (2i)
   0.24500000000000     0.37400000000000     0.10000000000000    Co   (2i)
  -0.24500000000000    -0.37400000000000    -0.10000000000000    Co   (2i)
   0.24400000000000     0.70900000000000     0.05400000000000    O    (2i)
  -0.24400000000000    -0.70900000000000    -0.05400000000000    O    (2i)
   0.20800000000000     0.09200000000000     0.19000000000000    O    (2i)
  -0.20800000000000    -0.09200000000000    -0.19000000000000    O    (2i)
   0.73500000000000     0.47600000000000     0.24900000000000    O    (2i)
  -0.73500000000000    -0.47600000000000    -0.24900000000000    O    (2i)
   0.56200000000000     0.84200000000000     0.28800000000000    O    (2i)
  -0.56200000000000    -0.84200000000000    -0.28800000000000    O    (2i)
   0.77000000000000     0.69800000000000     0.48900000000000    O    (2i)
  -0.77000000000000    -0.69800000000000    -0.48900000000000    O    (2i)
```

Kyanite (Al_2SiO_5, $S0_1$): A2B5C_aP32_2_4i_10i_2i - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Kyanite'
_chemical_formula_sum 'Al2 O5 Si'

loop_
_publ_author_name
 'H. Yang'
 'R. T. Downs'
 'L. W. Finger'
 'R. M. Hazen'
 'C. T. Prewitt'
_journal_name_full_name
;
 American Mineralogist
;
_journal_volume 82
_journal_year 1997
_journal_page_first 467
_journal_page_last 474
_publ_Section_title
;
 Compressibility and crystal structure of kyanite, Al$_{2}$SiO$_{5}$, at
    ↪ high pressure
;

_aflow_title 'Kyanite (Al$_{2}$SiO$_{5}$, $S0_{1}$) Structure'
_aflow_proto 'A2B5C_aP32_2_4i_10i_2i'
_aflow_params 'a,b/a,c/a,\alpha,\beta,\gamma,x_{1},y_{1},z_{1},x_{2},y_{
    ↪ 2},z_{2},x_{3},y_{3},z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{5},
    ↪ x_{6},y_{6},z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{
    ↪ 9},z_{9},x_{10},y_{10},z_{10},x_{11},y_{11},z_{11},x_{12},y_{12
    ↪ },z_{12},x_{13},y_{13},z_{13},x_{14},y_{14},z_{14},x_{15},y_{15
    ↪ },z_{15},x_{16},y_{16},z_{16}'
_aflow_params_values '7.12,1.10223314607,0.782837078652,89.974,101.177,
    ↪ 106.0,0.32533,0.70412,0.45812,0.2974,0.69882,0.9504,0.0998,
    ↪ 0.38615,0.64043,0.11205,0.9175,0.16469,0.10933,0.14685,0.12866,
    ↪ 0.12287,0.68535,0.18113,0.27507,0.45443,0.95474,0.28353,0.9357,
    ↪ 0.93567,0.10836,0.1521,0.66671,0.12192,0.63063,0.63939,0.28226,
    ↪ 0.44512,0.42868,0.29156,0.94684,0.46574,0.50074,0.27519,0.24405
    ↪ ,0.50154,0.23099,0.75595,0.29625,0.06488,0.70657,0.29102,
    ↪ 0.33168,0.18937'
_aflow_Strukturbericht '$S0_{1}$'
_aflow_Pearson 'aP32'

_symmetry_space_group_name_H-M "P -1"
_symmetry_Int_Tables_number 2

_cell_length_a    7.12000
_cell_length_b    7.84790
_cell_length_c    5.57380
_cell_angle_alpha 89.97400
_cell_angle_beta  101.17700
_cell_angle_gamma 106.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Al1  Al  2  i  0.32533  0.70412  0.45812  1.00000
Al2  Al  2  i  0.29740  0.69882  0.95040  1.00000
Al3  Al  2  i  0.09980  0.38615  0.64043  1.00000
Al4  Al  2  i  0.11205  0.91750  0.16469  1.00000
O1  O  2  i  0.10933  0.14685  0.12866  1.00000
O2  O  2  i  0.12287  0.68535  0.18113  1.00000
O3  O  2  i  0.27507  0.45443  0.95474  1.00000
O4  O  2  i  0.28353  0.93570  0.93567  1.00000
O5  O  2  i  0.10836  0.15210  0.66671  1.00000
O6  O  2  i  0.12192  0.63063  0.63939  1.00000
O7  O  2  i  0.28226  0.44512  0.42868  1.00000
O8  O  2  i  0.29156  0.94684  0.46574  1.00000
O9  O  2  i  0.50074  0.27519  0.24405  1.00000
O10  O  2  i  0.50154  0.23099  0.75595  1.00000
Si1  Si  2  i  0.29625  0.06488  0.70657  1.00000
Si2  Si  2  i  0.29102  0.33168  0.18937  1.00000
```

Kyanite (Al_2SiO_5, $S0_1$): A2B5C_aP32_2_4i_10i_2i - POSCAR

```
A2B5C_aP32_2_4i_10i_2i & a,b/a,c/a,alpha,beta,gamma,x1,y1,z1,x2,y2,z2,x3
    ↪ ,y3,z3,x4,y4,z4,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,
    ↪ x10,y10,z10,x11,y11,z11,x12,y12,z12,x13,y13,z13,x14,y14,z14,x15
    ↪ ,y15,z15,x16,y16,z16 --params=7.12,1.10223314607,0.782837078652
    ↪ ,89.974,101.177,106.0,0.32533,0.70412,0.45812,0.2974,0.69882,
    ↪ 0.9504,0.0998,0.38615,0.64043,0.11205,0.9175,0.16469,0.10933,
    ↪ 0.14685,0.12866,0.12287,0.68535,0.18113,0.27507,0.45443,0.95474
    ↪ ,0.28353,0.9357,0.93567,0.10836,0.1521,0.66671,0.12192,0.63063,
    ↪ 0.63939,0.28226,0.44512,0.42868,0.29156,0.94684,0.46574,0.50074
    ↪ ,0.27519,0.24405,0.50154,0.23099,0.75595,0.29625,0.06488,
    ↪ 0.70657,0.29102,0.33168,0.18937 & P-1 C_{i}^{1} #2 (i^16) &
    ↪ aP32 & $S0_{1}$ & Al2O5Si & Kyanite & H. Yang et al., Am.
    ↪ Mineral. 82, 467-474 (1997)
   1.00000000000000
   7.12000000000000     0.00000000000000     0.00000000000000
  -2.16317440471623     7.54388566355433     0.00000000000000
  -1.08042848514409    -0.30717664286521     5.45944715503020
     Al     O     Si
      8    20      4
Direct
   0.32533000000000     0.70412000000000     0.45812000000000    Al   (2i)
  -0.32533000000000    -0.70412000000000    -0.45812000000000    Al   (2i)
   0.29740000000000     0.69882000000000     0.95040000000000    Al   (2i)
  -0.29740000000000    -0.69882000000000    -0.95040000000000    Al   (2i)
   0.09980000000000     0.38615000000000     0.64043000000000    Al   (2i)
  -0.09980000000000    -0.38615000000000    -0.64043000000000    Al   (2i)
   0.11205000000000     0.91750000000000     0.16469000000000    Al   (2i)
  -0.11205000000000    -0.91750000000000    -0.16469000000000    Al   (2i)
   0.10933000000000     0.14685000000000     0.12866000000000    O    (2i)
  -0.10933000000000    -0.14685000000000    -0.12866000000000    O    (2i)
   0.12287000000000     0.68535000000000     0.18113000000000    O    (2i)
  -0.12287000000000    -0.68535000000000    -0.18113000000000    O    (2i)
   0.27507000000000     0.45443000000000     0.95474000000000    O    (2i)
  -0.27507000000000    -0.45443000000000    -0.95474000000000    O    (2i)
   0.28353000000000     0.93570000000000     0.93567000000000    O    (2i)
  -0.28353000000000    -0.93570000000000    -0.93567000000000    O    (2i)
   0.10836000000000     0.15210000000000     0.66671000000000    O    (2i)
  -0.10836000000000    -0.15210000000000    -0.66671000000000    O    (2i)
   0.12192000000000     0.63063000000000     0.63939000000000    O    (2i)
  -0.12192000000000    -0.63063000000000    -0.63939000000000    O    (2i)
   0.28226000000000     0.44512000000000     0.42868000000000    O    (2i)
  -0.28226000000000    -0.44512000000000    -0.42868000000000    O    (2i)
   0.29156000000000     0.94684000000000     0.46574000000000    O    (2i)
  -0.29156000000000    -0.94684000000000    -0.46574000000000    O    (2i)
   0.50074000000000     0.27519000000000     0.24405000000000    O    (2i)
  -0.50074000000000    -0.27519000000000    -0.24405000000000    O    (2i)
   0.50154000000000     0.23099000000000     0.75595000000000    O    (2i)
  -0.50154000000000    -0.23099000000000    -0.75595000000000    O    (2i)
   0.29625000000000     0.06488000000000     0.70657000000000    Si   (2i)
  -0.29625000000000    -0.06488000000000    -0.70657000000000    Si   (2i)
   0.29102000000000     0.33168000000000     0.18937000000000    Si   (2i)
  -0.29102000000000    -0.33168000000000    -0.18937000000000    Si   (2i)
```

α-$Ho_2Si_2O_7$: A2B7C2_aP44_2_4i_14i_4i - CIF

```
# CIF file
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_audit_creation_method FINDSYM

_chemical_name_mineral 'Ho2O7Si2'
_chemical_formula_sum 'Ho2 O7 Si2'

loop_
_publ_author_name
 'J. Felsche'
_journal_name_full_name
;
 Naturwissenschaften
;
_journal_volume 59
_journal_year 1972
_journal_page_first 35
_journal_page_last 36
_publ_Section_title
;
 A new silicate structure containing linear [Si$_{3}$O$_{10}$] groups
;
```

```
# Found in Revision of the crystallographic data of polymorphic Y${2}
      ↪ $Si$_{2}$O$_{7}$ and Y${2}$SiO${5}$ compounds, 2004

_aflow_title '$\alpha$-Ho$_{2}$Si$_{2}$O$_{7}$ Structure'
_aflow_proto 'A2B7C2_aP44_2_4i_14i_4i'
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      ↪ 2},z_{2},x_{3},y_{3},z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{5},
      ↪ x_{6},y_{6},z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{
      ↪ 9},z_{9},x_{10},y_{10},z_{10},x_{11},y_{11},z_{11},x_{12},y_{12
      ↪ },z_{12},x_{13},y_{13},z_{13},x_{14},y_{14},z_{14},x_{15},y_{15
      ↪ },z_{15},x_{16},y_{16},z_{16},x_{17},y_{17},z_{17},x_{18},y_{18
      ↪ },z_{18},x_{19},y_{19},z_{19},x_{20},y_{20},z_{20},x_{21},y_{21
      ↪ },z_{21},x_{22},y_{22},z_{22}'
_aflow_params_values '6.612,1.00862068966,1.82773744707,85.81,89.38,
      ↪ 88.57,-0.0521,0.331,0.11666,0.8845,0.0908,0.35915,0.3705,0.7756
      ↪ ,0.36947,0.6657,0.828,0.1059,0.6408,0.4893,0.1258,0.6226,0.1401
      ↪ ,0.2096,0.2968,0.2956,0.0948,0.4002,0.4255,0.304,0.5879,0.1717,
      ↪ 0.4454,0.224,0.095,0.3785,0.2937,0.422,0.5099,0.2862,0.2179,
      ↪ 0.687,-0.0374,0.2281,0.5714,0.0752,0.5789,0.6864,0.2414,-0.0859
      ↪ ,-0.001,0.3397,0.7841,0.1903,-0.0041,0.6708,0.0797,0.0023,
      ↪ 0.0137,0.1857,0.1539,0.8505,0.1168,0.4862,0.3353,0.1761,0.3781,
      ↪ 0.2726,0.4051,0.1457,0.3719,0.6179'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'aP44'

_symmetry_space_group_name_H-M "P -1"
_symmetry_Int_Tables_number 2

_cell_length_a 6.61200
_cell_length_b 6.66900
_cell_length_c 12.08500
_cell_angle_alpha 85.81000
_cell_angle_beta 89.38000
_cell_angle_gamma 88.57000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ho1 Ho 2 i -0.05210 0.33100 0.11666 1.00000
Ho2 Ho 2 i 0.88450 0.09080 0.35915 1.00000
Ho3 Ho 2 i 0.37050 0.77560 0.36947 1.00000
Ho4 Ho 2 i 0.66570 0.82800 0.10590 1.00000
O1 O 2 i 0.64080 0.48930 0.12580 1.00000
O2 O 2 i 0.62260 0.14010 0.20960 1.00000
O3 O 2 i 0.29680 0.29560 0.09480 1.00000
O4 O 2 i 0.40020 0.42550 0.30400 1.00000
O5 O 2 i 0.58790 0.17170 0.44540 1.00000
O6 O 2 i 0.22400 0.09500 0.37850 1.00000
O7 O 2 i 0.29370 0.42200 0.50990 1.00000
O8 O 2 i 0.28620 0.21790 0.68700 1.00000
O9 O 2 i -0.03740 0.22810 0.57140 1.00000
O10 O 2 i 0.07520 0.57890 0.68640 1.00000
O11 O 2 i 0.24140 -0.08590 -0.00100 1.00000
O12 O 2 i 0.33970 0.78410 0.19030 1.00000
O13 O 2 i -0.00410 0.67080 0.07970 1.00000
O14 O 2 i 0.00230 0.01370 0.18570 1.00000
Si1 Si 2 i 0.15390 0.85050 0.11680 1.00000
Si2 Si 2 i 0.48620 0.33530 0.17610 1.00000
Si3 Si 2 i 0.37810 0.27260 0.40510 1.00000
Si4 Si 2 i 0.14570 0.37190 0.61790 1.00000
```

α-$Ho_2Si_2O_7$: A2B7C2_aP44_2_4i_14i_4i - POSCAR

```
A2B7C2_aP44_2_4i_14i_4i & a,b/a,c/a,alpha,beta,gamma,x1,y1,z1,x2,y2,z2,
      ↪ x3,y3,z3,x4,y4,z4,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,
      ↪ x10,y10,z10,x11,y11,z11,x12,y12,z12,x13,y13,z13,x14,y14,z14,x15
      ↪ ,y15,z15,x16,y16,z16,x17,y17,z17,x18,y18,z18,x19,y19,z19,x20,
      ↪ y20,z20,x21,y21,z21,x22,y22,z22 --params=6.612,1.00862068966,
      ↪ 1.82773744707,85.81,89.38,88.57,-0.0521,0.331,0.11666,0.8845,
      ↪ 0.0908,0.35915,0.3705,0.7756,0.36947,0.6657,0.828,0.1059,0.6408
      ↪ ,0.4893,0.1258,0.6226,0.1401,0.2096,0.2968,0.2956,0.0948,0.4002
      ↪ ,0.4255,0.304,0.5879,0.1717,0.4454,0.224,0.095,0.3785,0.2937,
      ↪ 0.422,0.5099,0.2862,0.2179,0.687,-0.0374,0.2281,0.5714,0.0752,
      ↪ 0.5789,0.6864,0.2414,-0.0859,-0.001,0.3397,0.7841,0.1903,-
      ↪ 0.0041,0.6708,0.0797,0.0023,0.0137,0.1857,0.1539,0.8505,0.1168,
      ↪ 0.4862,0.3353,0.1761,0.3781,0.2726,0.4051,0.1457,0.3719,0.6179
      ↪ & P-1 C_{i}^{1} #2 (i^22) & aP44 & None & Ho2O7Si2 & Ho2O7Si2 &
      ↪ J. Felsche, Naturwissenschaften 59, 35-36 (1972)
   1.00000000000000
   6.61200000000000    0.00000000000000    0.00000000000000
   0.16642901147786    6.66692300721544    0.00000000000000
   0.13076973275040    0.87999065761810   12.05220895601720
     Ho      O     Si
      8     28      8
Direct
  -0.05210000000000    0.33100000000000    0.11666000000000   Ho   (2i)
   0.05210000000000   -0.33100000000000   -0.11666000000000   Ho   (2i)
   0.88450000000000    0.09080000000000    0.35915000000000   Ho   (2i)
  -0.88450000000000   -0.09080000000000   -0.35915000000000   Ho   (2i)
   0.37050000000000    0.77560000000000    0.36947000000000   Ho   (2i)
  -0.37050000000000   -0.77560000000000   -0.36947000000000   Ho   (2i)
   0.66570000000000    0.82800000000000    0.10590000000000   Ho   (2i)
  -0.66570000000000   -0.82800000000000   -0.10590000000000   Ho   (2i)
   0.64080000000000    0.48930000000000    0.12580000000000    O   (2i)
  -0.64080000000000   -0.48930000000000   -0.12580000000000    O   (2i)
   0.62260000000000    0.14010000000000    0.20960000000000    O   (2i)
  -0.62260000000000   -0.14010000000000   -0.20960000000000    O   (2i)
   0.29680000000000    0.29560000000000    0.09480000000000    O   (2i)
  -0.29680000000000   -0.29560000000000   -0.09480000000000    O   (2i)
   0.40020000000000    0.42550000000000    0.30400000000000    O   (2i)
  -0.40020000000000   -0.42550000000000   -0.30400000000000    O   (2i)
   0.58790000000000    0.17170000000000    0.44540000000000    O   (2i)
  -0.58790000000000   -0.17170000000000   -0.44540000000000    O   (2i)
   0.22400000000000    0.09500000000000    0.37850000000000    O   (2i)
  -0.22400000000000   -0.09500000000000   -0.37850000000000    O   (2i)
   0.29370000000000    0.42200000000000    0.50990000000000    O   (2i)
  -0.29370000000000   -0.42200000000000   -0.50990000000000    O   (2i)
   0.28620000000000    0.21790000000000    0.68700000000000    O   (2i)
  -0.28620000000000   -0.21790000000000   -0.68700000000000    O   (2i)
  -0.03740000000000    0.22810000000000    0.57140000000000    O   (2i)
   0.03740000000000   -0.22810000000000   -0.57140000000000    O   (2i)
   0.07520000000000    0.57890000000000    0.68640000000000    O   (2i)
  -0.07520000000000   -0.57890000000000   -0.68640000000000    O   (2i)
   0.24140000000000   -0.08590000000000   -0.00100000000000    O   (2i)
  -0.24140000000000    0.08590000000000    0.00100000000000    O   (2i)
   0.33970000000000    0.78410000000000    0.19030000000000    O   (2i)
  -0.33970000000000   -0.78410000000000   -0.19030000000000    O   (2i)
  -0.00410000000000    0.67080000000000    0.07970000000000    O   (2i)
   0.00410000000000   -0.67080000000000   -0.07970000000000    O   (2i)
   0.00230000000000    0.01370000000000    0.18570000000000    O   (2i)
  -0.00230000000000   -0.01370000000000   -0.18570000000000    O   (2i)
   0.15390000000000    0.85050000000000    0.11680000000000   Si   (2i)
  -0.15390000000000   -0.85050000000000   -0.11680000000000   Si   (2i)
   0.48620000000000    0.33530000000000    0.17610000000000   Si   (2i)
  -0.48620000000000   -0.33530000000000   -0.17610000000000   Si   (2i)
   0.37810000000000    0.27260000000000    0.40510000000000   Si   (2i)
  -0.37810000000000   -0.27260000000000   -0.40510000000000   Si   (2i)
   0.14570000000000    0.37190000000000    0.61790000000000   Si   (2i)
  -0.14570000000000   -0.37190000000000   -0.61790000000000   Si   (2i)
```

$Co_3(SeO_3)_3{\cdot}H_2O$: A3B2C10D3_aP36_2_ah2i_2i_10i_3i - CIF

```
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_chemical_name_mineral 'Co3H2O10Se3'
_chemical_formula_sum 'Co3 H2 O10 Se3'

loop_
_publ_author_name
 'M. Wildner'
_journal_name_full_name
;
 Monatshefte f{\"u}r Chemie - Chemical Monthly
;
_journal_volume 122
_journal_year 1991
_journal_page_first 585
_journal_page_last 594
_publ_Section_title
;
 Crystal structures of Co$_{3}$(SeO$_{3}$)$_{3}$\cdot$H$_{2}$O and Ni$_
      ↪ {3}$(SeO$_{3}$)$_{3}$\cdot$H$_{2}$O, two new isotypic
      ↪ compounds
;

# Found in K. M. Taddei et al., Tunable magnetic order in low-symmetry
      ↪ SeO$_{3}$ ligand linked $TM$$_{3}$(SeO$_{3}$)$_{3}$H$_{2}$O (
      ↪ $TM$ = Mn, Co and Ni) compounds, Phys. Rev. Mater. 4, 024410 (
      ↪ 2020).

_aflow_title 'Co$_{3}$(SeO$_{3}$)$_{3}$\cdot$H$_{2}$O Structure'
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      ↪ x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{10},y_{10},z_{10},x_{11
      ↪ },y_{11},z_{11},x_{12},y_{12},z_{12},x_{13},y_{13},z_{13},x_{14
      ↪ },y_{14},z_{14},x_{15},y_{15},z_{15},x_{16},y_{16},z_{16},x_{17
      ↪ },y_{17},z_{17},x_{18},y_{18},z_{18},x_{19},y_{19},z_{19}'
_aflow_params_values '8.102,1.0144408788,1.05801036781,69.15,62.88,67.23
      ↪ ,0.65388,0.03726,0.79895,0.71664,0.85636,0.39005,0.004,0.867,
      ↪ 0.704,-0.065,0.823,0.624,0.3023,0.7124,0.3858,0.5468,0.6953,
      ↪ 0.5643,0.2128,0.8964,0.7589,0.5375,-0.0553,0.6739,-0.0504,
      ↪ 0.7241,0.4515,0.7083,0.518,0.2467,0.5988,0.8268,0.0273,0.8667,
      ↪ 0.7852,0.1424,0.1894,0.8592,0.1263,-0.0821,0.8725,0.6893,
      ↪ 0.56457,0.23365,0.23877,0.79662,0.66737,0.06492,0.17624,0.66966
      ↪ ,0.29811'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'aP36'

_symmetry_space_group_name_H-M "P -1"
_symmetry_Int_Tables_number 2

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_cell_length_b 8.21900
_cell_length_c 8.57200
_cell_angle_alpha 69.15000
_cell_angle_beta 62.88000
_cell_angle_gamma 67.23000

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_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,-z
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_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
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Co2 Co 1 h 0.50000 0.50000 0.50000 1.00000
Co3 Co 2 i 0.65388 0.03726 0.79895 1.00000
Co4 Co 2 i 0.71664 0.85636 0.39005 1.00000
H1 H 2 i 0.00400 0.86700 0.70400 1.00000
H2 H 2 i −0.06500 0.82300 0.62400 1.00000
O1 O 2 i 0.30230 0.71240 0.38580 1.00000
O2 O 2 i 0.54680 0.69530 0.56430 1.00000
O3 O 2 i 0.21280 0.89640 0.75890 1.00000
O4 O 2 i 0.53750 −0.05530 0.67390 1.00000
O5 O 2 i −0.05040 0.72410 0.45150 1.00000
O6 O 2 i 0.70830 0.51800 0.24670 1.00000
O7 O 2 i 0.59880 0.82680 0.02730 1.00000
O8 O 2 i 0.86670 0.78520 0.14240 1.00000
O9 O 2 i 0.18940 0.85920 0.12630 1.00000
O10 O 2 i −0.08210 0.87250 0.68930 1.00000
Se1 Se 2 i 0.56457 0.23365 0.23877 1.00000
Se2 Se 2 i 0.79662 0.66737 0.06492 1.00000
Se3 Se 2 i 0.17624 0.66966 0.29811 1.00000
```

$Co_3(SeO_3)_3{\cdot}H_2O$: A3B2C10D3_aP36_2_ah2i_2i_10i_3i - POSCAR

```
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    ↪ ,y4,z4,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10
    ↪ ,x11,y11,z11,x12,y12,z12,x13,y13,z13,x14,y14,z14,x15,y15,z15,
    ↪ x16,y16,z16,x17,y17,z17,x18,y18,z18,x19,y19,z19 --params=8.102,
    ↪ 1.0144408788,1.05801036781,69.15,62.88,67.23,0.65388,0.03726,
    ↪ 0.79895,0.71664,0.85636,0.39005,0.004,0.867,0.704,−0.065,0.823,
    ↪ 0.624,0.3023,0.7124,0.3858,0.5468,0.6953,0.5643,0.2128,0.8964,
    ↪ 0.7589,0.5375,−0.0553,0.6739,−0.0504,0.7241,0.4515,0.7083,0.518
    ↪ ,0.2467,0.5988,0.8268,0.0273,0.8667,0.7852,0.1424,0.1894,0.8592
    ↪ ,0.1263,−0.0821,0.8725,0.6893,0.56457,0.23365,0.23877,0.79662,
    ↪ 0.66737,0.06492,0.17624,0.66966,0.29811 & P−1 C_{i}^{1} #2 (ahi
    ↪ ^17) & aP36 & None & Co3H2O10Se3 & Co3H2O10Se3 & M. Wildner,
    ↪ Monatshefte f{"u}r Chemie − Chemical Monthly 122, 585−594 (1991
    ↪ )
   1.00000000000000
   8.10200000000000    0.00000000000000    0.00000000000000
   3.18102296897454    7.57845986140036    0.00000000000000
   3.90759439489992    1.66864563346477    7.44483121332282
    Co   H   O   Se
     6   4  20    6
Direct
   0.00000000000000    0.00000000000000    0.00000000000000  Co   (1a)
   0.50000000000000    0.50000000000000    0.50000000000000  Co   (1h)
   0.65388000000000    0.03726000000000    0.79895000000000  Co   (2i)
  −0.65388000000000   −0.03726000000000   −0.79895000000000  Co   (2i)
   0.71664000000000    0.85636000000000    0.39005000000000  Co   (2i)
  −0.71664000000000   −0.85636000000000   −0.39005000000000  Co   (2i)
   0.00400000000000    0.86700000000000    0.70400000000000   H   (2i)
  −0.00400000000000   −0.86700000000000   −0.70400000000000   H   (2i)
  −0.06500000000000    0.82300000000000    0.62400000000000   H   (2i)
   0.06500000000000   −0.82300000000000   −0.62400000000000   H   (2i)
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   0.54680000000000    0.69530000000000    0.56430000000000   O   (2i)
  −0.54680000000000   −0.69530000000000   −0.56430000000000   O   (2i)
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  −0.21280000000000   −0.89640000000000   −0.75890000000000   O   (2i)
   0.53750000000000   −0.05530000000000    0.67390000000000   O   (2i)
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  −0.70830000000000   −0.51800000000000   −0.24670000000000   O   (2i)
   0.59880000000000    0.82680000000000    0.02730000000000   O   (2i)
  −0.59880000000000   −0.82680000000000   −0.02730000000000   O   (2i)
   0.86670000000000    0.78520000000000    0.14240000000000   O   (2i)
  −0.86670000000000   −0.78520000000000   −0.14240000000000   O   (2i)
   0.18940000000000    0.85920000000000    0.12630000000000   O   (2i)
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  −0.08210000000000    0.87250000000000    0.68930000000000   O   (2i)
   0.08210000000000   −0.87250000000000   −0.68930000000000   O   (2i)
   0.56457000000000    0.23365000000000    0.23877000000000  Se   (2i)
  −0.56457000000000   −0.23365000000000   −0.23877000000000  Se   (2i)
   0.79662000000000    0.66737000000000    0.06492000000000  Se   (2i)
  −0.79662000000000   −0.66737000000000   −0.06492000000000  Se   (2i)
   0.17624000000000    0.66966000000000    0.29811000000000  Se   (2i)
  −0.17624000000000   −0.66966000000000   −0.29811000000000  Se   (2i)
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δ-WO_3: A3B_aP32_2_12i_4i - CIF

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 'A. W. Sleight'
 'T. Vogt'
_journal_name_full_name
;
 Journal of Solid State Chemistry
;
_journal_volume 131
_journal_year 1997
_journal_page_first 9
_journal_page_last 17
_publ_Section_title
;
 Ferroelectric Tungsten Trioxide
;

_aflow_title '$\delta-WO$_{3}$ Structure'
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    ↪ x_{6},y_{6},z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{
    ↪ 9},z_{9},x_{10},y_{10},z_{10},x_{11},y_{11},z_{11},x_{12},y_{12
    ↪ },z_{12},x_{13},y_{13},z_{13},x_{14},y_{14},z_{14},x_{15},y_{15
    ↪ },z_{15},x_{16},y_{16},z_{16}'
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    ↪ 0.2884,0.4972,0.9638,0.2878,0.2851,0.2574,0.287,0.2204,0.763,
    ↪ 0.2232,0.2186,0.2627,0.7258,0.284,0.7583,0.7679,0.2943,0.0422,−
    ↪ 0.0002,0.2971,0.5446,0.4982,0.2096,0.482,−0.0072,0.2088,−0.017,
    ↪ 0.5051,0.2566,0.0259,0.285,0.2502,0.528,0.2158,0.2438,0.0313,
    ↪ 0.7817,0.2499,0.5338,0.719'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'aP32'

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_symmetry_Int_Tables_number 2

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_cell_angle_gamma  89.07000

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_space_group_symop_operation_xyz
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2 −x,−y,−z

loop_
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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
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O3 O 2 i 0.00760 0.46600 0.28840 1.00000
O4 O 2 i 0.49720 0.96380 0.28780 1.00000
O5 O 2 i 0.28510 0.25740 0.28700 1.00000
O6 O 2 i 0.22040 0.76300 0.22320 1.00000
O7 O 2 i 0.21860 0.26270 0.72580 1.00000
O8 O 2 i 0.28400 0.75830 0.76790 1.00000
O9 O 2 i 0.29430 0.04220 −0.00020 1.00000
O10 O 2 i 0.29710 0.54460 0.49820 1.00000
O11 O 2 i 0.20960 0.48200 −0.00720 1.00000
O12 O 2 i 0.20880 −0.01700 0.50510 1.00000
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W2 W 2 i 0.25020 0.52800 0.21580 1.00000
W3 W 2 i 0.24380 0.03130 0.78170 1.00000
W4 W 2 i 0.24990 0.53380 0.71900 1.00000
```

δ-WO_3: A3B_aP32_2_12i_4i - POSCAR

```
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    ↪ z3,x4,y4,z4,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,
    ↪ y10,z10,x11,y11,z11,x12,y12,z12,x13,y13,z13,x14,y14,z14,x15,y15
    ↪ ,z15,x16,y16,z16 --params=7.309,1.02914215351,1.05048570256,
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    ↪ 0.0076,0.466,0.2884,0.4972,0.9638,0.2878,0.2851,0.2574,0.287,
    ↪ 0.2204,0.763,0.2232,0.2186,0.2627,0.7258,0.284,0.7583,0.7679,
    ↪ 0.2943,0.0422,−0.0002,0.2971,0.5446,0.4982,0.2096,0.482,−0.0072
    ↪ ,0.2088,−0.017,0.5051,0.2566,0.0259,0.285,0.2502,0.528,0.2158,
    ↪ 0.2438,0.0313,0.7817,0.2499,0.5338,0.719 & P−1 C_{i}^{1} #2 (i^
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    ↪ Sleight and T. Vogt, J. Solid State Chem. 131, 9−17 (1997)
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   0.12208844854962    7.52100913513145    0.00000000000000
   0.12328057187699    0.15747592326685    7.67539492366282
     O   W
    24   8
Direct
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  −0.00070000000000   −0.03860000000000   −0.21000000000000   O   (2i)
   0.50380000000000    0.53610000000000    0.21810000000000   O   (2i)
  −0.50380000000000   −0.53610000000000   −0.21810000000000   O   (2i)
   0.00760000000000    0.46600000000000    0.28840000000000   O   (2i)
  −0.00760000000000   −0.46600000000000   −0.28840000000000   O   (2i)
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  −0.28510000000000   −0.25740000000000   −0.28700000000000   O   (2i)
   0.22040000000000    0.76300000000000    0.22320000000000   O   (2i)
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   0.21860000000000    0.26270000000000    0.72580000000000   O   (2i)
```

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 −0.28400000000000  −0.75830000000000  −0.76790000000000   O   (2i)
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 −0.29430000000000  −0.04220000000000   0.00020000000000   O   (2i)
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 −0.29710000000000  −0.54460000000000  −0.49820000000000   O   (2i)
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 −0.20960000000000  −0.48200000000000   0.00720000000000   O   (2i)
  0.20880000000000  −0.01700000000000   0.50510000000000   O   (2i)
 −0.20880000000000   0.01700000000000  −0.50510000000000   O   (2i)
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  0.25020000000000   0.52800000000000   0.21580000000000   W   (2i)
 −0.25020000000000  −0.52800000000000  −0.21580000000000   W   (2i)
  0.24380000000000   0.03130000000000   0.78170000000000   W   (2i)
 −0.24380000000000  −0.03130000000000  −0.78170000000000   W   (2i)
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 −0.24990000000000  −0.53380000000000  −0.71900000000000   W   (2i)
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Chalcanthite ($CuSO_4{\cdot}5H_2O$, $H4_{10}$): AB10C9D_aP42_2_ae_10i_9i_i - CIF

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 'G. E. Bacon'
 'D. H. Titterton'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 141
_journal_year 1975
_journal_page_first 330
_journal_page_last 341
_publ_Section_title
;
 Neutron-diffraction studies of CuSO$_{4}$$\cdot$5H$_{2}$O and CuSO$_{4}
      ↪ $$\cdot$5D$_{2}$O
;

# Found in The American Mineralogist Crystal Structure Database, 2003

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      ↪ x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{10},y_{10},z_{10},x_{11
      ↪ },y_{11},z_{11},x_{12},y_{12},z_{12},x_{13},y_{13},z_{13},x_{14
      ↪ },y_{14},z_{14},x_{15},y_{15},z_{15},x_{16},y_{16},z_{16},x_{17
      ↪ },y_{17},z_{17},x_{18},y_{18},z_{18},x_{19},y_{19},z_{19},x_{20
      ↪ },y_{20},z_{20},x_{21},y_{21},z_{21},x_{22},y_{22},z_{22}'
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      ↪ 0.6033,0.1321,0.6671,0.4108,0.1932,0.6922,0.9072,0.152,0.6734,
      ↪ 0.2442,0.3172,0.796,0.8061,0.3724,0.6363,0.0444,0.3022,0.3849,
      ↪ 0.8176,0.0737,0.1519,0.2887,0.1177,0.149,0.4654,0.4063,0.2975,
      ↪ 0.756,0.4161,0.0191,0.435,0.1263,0.6289,0.0133,0.2871,0.6253'
_aflow_Strukturbericht '$H4_{10}$'
_aflow_Pearson 'aP42'

_symmetry_space_group_name_H-M "P -1"
_symmetry_Int_Tables_number 2

_cell_length_a       6.14100
_cell_length_b      10.73600
_cell_length_c       5.98600
_cell_angle_alpha   82.26667
_cell_angle_beta   107.43333
_cell_angle_gamma  102.66667

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cu1 Cu 1 a 0.0000 0.0000 0.0000 1.00000
Cu2 Cu 1 e 0.5000 0.5000 0.0000 1.00000
H1 H 2 i 0.8980 0.1412 0.2547 1.00000
H2 H 2 i 0.7185 0.0126 0.2283 1.00000
H3 H 2 i 0.3010 0.2016 0.0667 1.00000
H4 H 2 i 0.3341 0.1270 0.3188 1.00000
H5 H 2 i 0.3231 0.3785 0.3406 1.00000
H6 H 2 i 0.6016 0.3937 0.4256 1.00000
H7 H 2 i 0.8012 0.4011 0.8847 1.00000
H8 H 2 i 0.8570 0.3845 0.1620 1.00000
H9 H 2 i 0.6033 0.1321 0.6671 1.00000
H10 H 2 i 0.4108 0.1932 0.6922 1.00000
O1 O 2 i 0.9072 0.1520 0.6734 1.00000
O2 O 2 i 0.2442 0.3172 0.7960 1.00000
O3 O 2 i 0.8061 0.3724 0.6363 1.00000
O4 O 2 i 0.0444 0.3022 0.3849 1.00000
O5 O 2 i 0.8176 0.0737 0.1519 1.00000
O6 O 2 i 0.2887 0.1177 0.1490 1.00000
O7 O 2 i 0.4654 0.4063 0.2975 1.00000
O8 O 2 i 0.7560 0.4161 0.0191 1.00000
O9 O 2 i 0.4350 0.1263 0.6289 1.00000
S1 S 2 i 0.0133 0.2871 0.6253 1.00000
```

Chalcanthite ($CuSO_4{\cdot}5H_2O$, $H4_{10}$): AB10C9D_aP42_2_ae_10i_9i_i - POSCAR

```
AB10C9D_aP42_2_ae_10i_9i_i & a,b/a,c/a,alpha,beta,gamma,x3,y3,z3,x4,y4,
      ↪ z4,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10,x11
      ↪ ,y11,z11,x12,y12,z12,x13,y13,z13,x14,y14,z14,x15,y15,z15,x16,
      ↪ y16,z16,x17,y17,z17,x18,y18,z18,x19,y19,z19,x20,y20,z20,x21,y21
      ↪ ,z21,x22,y22,z22 --params=6.141,1.74824947077,0.974759811106,
      ↪ 82.26667,107.43333,102.66667,0.898,0.1412,0.2547,0.7185,0.0126,
      ↪ 0.2283,0.301,0.2016,0.0667,0.3341,0.127,0.3188,0.3231,0.3785,
      ↪ 0.3406,0.6016,0.3937,0.4256,0.8012,0.4011,0.8847,0.857,0.3845,
      ↪ 0.162,0.6033,0.1321,0.6671,0.4108,0.1932,0.6922,0.9072,0.152,
      ↪ 0.6734,0.2442,0.3172,0.796,0.8061,0.3724,0.6363,0.0444,0.3022,
      ↪ 0.3849,0.8176,0.0737,0.1519,0.2887,0.1177,0.149,0.4654,0.4063,
      ↪ 0.2975,0.756,0.4161,0.0191,0.435,0.1263,0.6289,0.0133,0.2871,
      ↪ 0.6253 & P-1 C_{i}^{1} #2 (aei^20) & aP42 & $H4_{10}$ &
      ↪ CuH10O9S & Chalcanthite & G. E. Bacon and D. H. Titterton,
      ↪ Zeitschrift f{"u}r Kristallographie - Crystalline Materials 141
      ↪ , 330-341 (1975)
   1.00000000000000
   6.14100000000000   0.00000000000000   0.00000000000000
  -2.35417591867527  10.47471010309730   0.00000000000000
  -1.79338070202199   0.42252508515638   5.69538885503253
     Cu     H     O     S
      2    20    18     2
Direct
   0.00000000000000   0.00000000000000   0.00000000000000  Cu   (1a)
   0.50000000000000   0.50000000000000   0.00000000000000  Cu   (1e)
   0.89800000000000   0.14120000000000   0.25470000000000   H   (2i)
  -0.89800000000000  -0.14120000000000  -0.25470000000000   H   (2i)
   0.71850000000000   0.01260000000000   0.22830000000000   H   (2i)
  -0.71850000000000  -0.01260000000000  -0.22830000000000   H   (2i)
   0.30100000000000   0.20160000000000   0.06670000000000   H   (2i)
  -0.30100000000000  -0.20160000000000  -0.06670000000000   H   (2i)
   0.33410000000000   0.12700000000000   0.31880000000000   H   (2i)
  -0.33410000000000  -0.12700000000000  -0.31880000000000   H   (2i)
   0.32310000000000   0.37850000000000   0.34060000000000   H   (2i)
  -0.32310000000000  -0.37850000000000  -0.34060000000000   H   (2i)
   0.60160000000000   0.39370000000000   0.42560000000000   H   (2i)
  -0.60160000000000  -0.39370000000000  -0.42560000000000   H   (2i)
   0.80120000000000   0.40110000000000   0.88470000000000   H   (2i)
  -0.80120000000000  -0.40110000000000  -0.88470000000000   H   (2i)
   0.85700000000000   0.38450000000000   0.16200000000000   H   (2i)
  -0.85700000000000  -0.38450000000000  -0.16200000000000   H   (2i)
   0.60330000000000   0.13210000000000   0.66710000000000   H   (2i)
  -0.60330000000000  -0.13210000000000  -0.66710000000000   H   (2i)
   0.41080000000000   0.19320000000000   0.69220000000000   H   (2i)
  -0.41080000000000  -0.19320000000000  -0.69220000000000   H   (2i)
   0.90720000000000   0.15200000000000   0.67340000000000   O   (2i)
  -0.90720000000000  -0.15200000000000  -0.67340000000000   O   (2i)
   0.24420000000000   0.31720000000000   0.79600000000000   O   (2i)
  -0.24420000000000  -0.31720000000000  -0.79600000000000   O   (2i)
   0.80610000000000   0.37240000000000   0.63630000000000   O   (2i)
  -0.80610000000000  -0.37240000000000  -0.63630000000000   O   (2i)
   0.04440000000000   0.30220000000000   0.38490000000000   O   (2i)
  -0.04440000000000  -0.30220000000000  -0.38490000000000   O   (2i)
   0.81760000000000   0.07370000000000   0.15190000000000   O   (2i)
  -0.81760000000000  -0.07370000000000  -0.15190000000000   O   (2i)
   0.28870000000000   0.11770000000000   0.14900000000000   O   (2i)
  -0.28870000000000  -0.11770000000000  -0.14900000000000   O   (2i)
   0.46540000000000   0.40630000000000   0.29750000000000   O   (2i)
  -0.46540000000000  -0.40630000000000  -0.29750000000000   O   (2i)
   0.75600000000000   0.41610000000000   0.01910000000000   O   (2i)
  -0.75600000000000  -0.41610000000000  -0.01910000000000   O   (2i)
   0.43500000000000   0.12630000000000   0.62890000000000   O   (2i)
  -0.43500000000000  -0.12630000000000  -0.62890000000000   O   (2i)
   0.01330000000000   0.28710000000000   0.62530000000000   S   (2i)
  -0.01330000000000  -0.28710000000000  -0.62530000000000   S   (2i)
```

Boric Acid (H_3BO_3, $G5_1$): AB3C3_aP28_2_2i_6i_6i - CIF

```
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_chemical_formula_sum 'B H3 O3'

loop_
_publ_author_name
 'W. H. Zachariasen'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 88
_journal_year 1934
_journal_page_first 150
_journal_page_last 161
_publ_Section_title
;
 The Crystal Lattice of Boric Acid, BO$_{3}$H$_{3}$
;
```

```
# Found in Strukturbericht Band III 1933-1935, 1937

_aflow_title 'Boric Acid (H$_{3}$BO$_{3}$, $G5_{1}$) Structure'
_aflow_proto 'AB3C3_aP28_2_2i_6i_6i'
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    ↪ x_{6},y_{6},z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{
    ↪ 9},z_{9},x_{10},y_{10},z_{10},x_{11},y_{11},z_{11},x_{12},y_{12
    ↪ },z_{12},x_{13},y_{13},z_{13},x_{14},y_{14},z_{14}'
_aflow_params_values '6.56,1.07317073171,1.07317073171,120.0,92.5,
    ↪ 101.16667,0.25,0.653,0.431,0.25,0.31,0.764,0.25,0.319,0.431,
    ↪ 0.25,0.319,0.097,0.25,-0.01,0.431,0.25,0.653,0.097,0.25,-0.01,
    ↪ 0.764,0.25,0.65,0.764,0.25,0.431,0.319,0.25,0.764,0.319,0.25,
    ↪ 0.764,0.653,0.25,0.21,0.542,0.25,0.21,0.875,0.25,0.542,0.875'
_aflow_Strukturbericht '$G5_{1}$'
_aflow_Pearson 'aP28'

_symmetry_space_group_name_H-M "P -1"
_symmetry_Int_Tables_number 2

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_cell_length_b      7.04000
_cell_length_c      7.04000
_cell_angle_alpha  120.00000
_cell_angle_beta    92.50000
_cell_angle_gamma  101.16667

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
B1 B 2 i 0.25000 0.65300 0.43100 1.00000
B2 B 2 i 0.25000 0.31000 0.76400 1.00000
H1 H 2 i 0.25000 0.31900 0.43100 1.00000
H2 H 2 i 0.25000 0.31900 0.09700 1.00000
H3 H 2 i 0.25000 -0.01000 0.43100 1.00000
H4 H 2 i 0.25000 0.65300 0.09700 1.00000
H5 H 2 i 0.25000 -0.01000 0.76400 1.00000
H6 H 2 i 0.25000 0.65000 0.76400 1.00000
O1 O 2 i 0.25000 0.43100 0.31900 1.00000
O2 O 2 i 0.25000 0.76400 0.31900 1.00000
O3 O 2 i 0.25000 0.76400 0.65300 1.00000
O4 O 2 i 0.25000 0.21000 0.54200 1.00000
O5 O 2 i 0.25000 0.21000 0.87500 1.00000
O6 O 2 i 0.25000 0.54200 0.87500 1.00000
```

Boric Acid (H_3BO_3, $G5_1$): AB3C3_aP28_2_2i_6i_6i - POSCAR

```
AB3C3_aP28_2_2i_6i_6i & a,b/a,c/a,alpha,beta,gamma,x1,y1,z1,x2,y2,z2,x3,
    ↪ y3,z3,x4,y4,z4,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10
    ↪ ,y10,z10,x11,y11,z11,x12,y12,z12,x13,y13,z13,x14,y14,z14 --
    ↪ params=6.56,1.07317073171,1.07317073171,120.0,92.5,101.16667,
    ↪ 0.25,0.653,0.431,0.25,0.31,0.764,0.25,0.319,0.431,0.25,0.319,
    ↪ 0.097,0.25,-0.01,0.431,0.25,0.653,0.097,0.25,-0.01,0.764,0.25,
    ↪ 0.65,0.764,0.25,0.431,0.319,0.25,0.764,0.319,0.25,0.764,0.653,
    ↪ 0.25,0.21,0.542,0.25,0.21,0.875,0.25,0.542,0.875 & P-1 C_{i}^{1
    ↪ } #2 (i^14) & aP28 & $G5_{1}$ & BH3O3 & Boric acid & W. H.
    ↪ Zachariasen, Zeitschrift f{"u}r Kristallographie - Crystalline
    ↪ Materials 88, 150-161 (1934)
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   6.56000000000000    0.00000000000000    0.00000000000000
  -1.36339229932856    6.90671857238527    0.00000000000000
  -0.30708048705196   -3.64854466085737    6.01293800335586
     B     H     O
     4    12    12
Direct
   0.25000000000000    0.65300000000000    0.43100000000000    B   (2i)
  -0.25000000000000   -0.65300000000000   -0.43100000000000    B   (2i)
   0.25000000000000    0.31000000000000    0.76400000000000    B   (2i)
  -0.25000000000000   -0.31000000000000   -0.76400000000000    B   (2i)
   0.25000000000000    0.31900000000000    0.43100000000000    H   (2i)
  -0.25000000000000   -0.31900000000000   -0.43100000000000    H   (2i)
   0.25000000000000    0.31900000000000    0.09700000000000    H   (2i)
  -0.25000000000000   -0.31900000000000   -0.09700000000000    H   (2i)
   0.25000000000000   -0.01000000000000    0.43100000000000    H   (2i)
  -0.25000000000000    0.01000000000000   -0.43100000000000    H   (2i)
   0.25000000000000    0.65300000000000    0.09700000000000    H   (2i)
  -0.25000000000000   -0.65300000000000   -0.09700000000000    H   (2i)
   0.25000000000000   -0.01000000000000    0.76400000000000    H   (2i)
  -0.25000000000000    0.01000000000000   -0.76400000000000    H   (2i)
   0.25000000000000    0.65000000000000    0.76400000000000    H   (2i)
  -0.25000000000000   -0.65000000000000   -0.76400000000000    H   (2i)
   0.25000000000000    0.43100000000000    0.31900000000000    O   (2i)
  -0.25000000000000   -0.43100000000000   -0.31900000000000    O   (2i)
   0.25000000000000    0.76400000000000    0.31900000000000    O   (2i)
  -0.25000000000000   -0.76400000000000   -0.31900000000000    O   (2i)
   0.25000000000000    0.76400000000000    0.65300000000000    O   (2i)
  -0.25000000000000   -0.76400000000000   -0.65300000000000    O   (2i)
   0.25000000000000    0.21000000000000    0.54200000000000    O   (2i)
  -0.25000000000000   -0.21000000000000   -0.54200000000000    O   (2i)
   0.25000000000000    0.21000000000000    0.87500000000000    O   (2i)
  -0.25000000000000   -0.21000000000000   -0.87500000000000    O   (2i)
   0.25000000000000    0.54200000000000    0.87500000000000    O   (2i)
  -0.25000000000000   -0.54200000000000   -0.87500000000000    O   (2i)
```

Wollastonite ($CaSiO_3$): AB3C_aP30_2_3i_9i_3i - CIF

```
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_audit_creation_method FINDSYM

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_chemical_formula_sum 'Ca O3 Si'

loop_
_publ_author_name
 'M. J. Buerger'
 'C. T. Prewitt'
_journal_name_full_name
;
 Proceedings of the National Academy of Sciences of the United States of
    ↪ America
;
_journal_volume 47
_journal_year 1961
_journal_page_first 1884
_journal_page_last 1888
_publ_Section_title
;
 The Crystal Structures of Wollastonite and Pectolite
;

_aflow_title 'Wollastonite (CaSiO$_{3}$) Structure'
_aflow_proto 'AB3C_aP30_2_3i_9i_3i'
_aflow_params 'a,b/a,c/a,\alpha,\beta,\gamma,x_{1},y_{1},z_{1},x_{2},y_{
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    ↪ x_{6},y_{6},z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{
    ↪ 9},z_{9},x_{10},y_{10},z_{10},x_{11},y_{11},z_{11},x_{12},y_{12
    ↪ },z_{12},x_{13},y_{13},z_{13},x_{14},y_{14},z_{14},x_{15},y_{15
    ↪ },z_{15}'
_aflow_params_values '7.94,0.921914357683,0.890428211587,90.03333,
    ↪ 95.36667,103.43333,0.1985,0.4228,0.7608,0.2027,-0.0707,0.764,
    ↪ 0.4966,0.2495,0.472,0.4291,0.2314,0.8019,0.4008,0.7259,0.8302,
    ↪ 0.3037,0.4635,0.4641,0.3017,-0.0626,0.4655,0.0154,0.6254,0.7343
    ↪ ,0.0175,0.1319,0.7353,0.2732,0.5118,0.0919,0.2713,0.8717,0.094,
    ↪ 0.2188,0.1784,0.2228,0.1852,0.387,0.2687,0.1849,-0.0455,0.2692,
    ↪ 0.397,0.7235,0.056'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'aP30'

_symmetry_space_group_name_H-M "P -1"
_symmetry_Int_Tables_number 2

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_cell_length_c      7.07000
_cell_angle_alpha  90.03333
_cell_angle_beta   95.36667
_cell_angle_gamma  103.43333

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ca1 Ca 2 i 0.19850 0.42280 0.76080 1.00000
Ca2 Ca 2 i 0.20270 -0.07070 0.76400 1.00000
Ca3 Ca 2 i 0.49660 0.24950 0.47200 1.00000
O1 O 2 i 0.42910 0.23140 0.80190 1.00000
O2 O 2 i 0.40080 0.72590 0.83020 1.00000
O3 O 2 i 0.30370 0.46350 0.46410 1.00000
O4 O 2 i 0.30170 -0.06260 0.46550 1.00000
O5 O 2 i 0.01540 0.62540 0.73430 1.00000
O6 O 2 i 0.01750 0.13190 0.73530 1.00000
O7 O 2 i 0.27320 0.51180 0.09190 1.00000
O8 O 2 i 0.27130 0.87170 0.09400 1.00000
O9 O 2 i 0.21880 0.17840 0.22280 1.00000
Si1 Si 2 i 0.18520 0.38700 0.26870 1.00000
Si2 Si 2 i 0.18490 -0.04550 0.26920 1.00000
Si3 Si 2 i 0.39700 0.72350 0.05600 1.00000
```

Wollastonite ($CaSiO_3$): AB3C_aP30_2_3i_9i_3i - POSCAR

```
AB3C_aP30_2_3i_9i_3i & a,b/a,c/a,alpha,beta,gamma,x1,y1,z1,x2,y2,z2,x3,
    ↪ y3,z3,x4,y4,z4,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10
    ↪ ,y10,z10,x11,y11,z11,x12,y12,z12,x13,y13,z13,x14,y14,z14,x15,
    ↪ y15,z15 --params=7.94,0.921914357683,0.890428211587,90.03333,
    ↪ 95.36667,103.43333,0.1985,0.4228,0.7608,0.2027,-0.0707,0.764,
    ↪ 0.4966,0.2495,0.472,0.4291,0.2314,0.8019,0.4008,0.7259,0.8302,
    ↪ 0.3037,0.4635,0.4641,0.3017,-0.0626,0.4655,0.0154,0.6254,0.7343
    ↪ ,0.0175,0.1319,0.7353,0.2732,0.5118,0.0919,0.2713,0.8717,0.094,
    ↪ 0.2188,0.1784,0.2228,0.1852,0.387,0.2687,0.1849,-0.0455,0.2692,
    ↪ 0.397,0.7235,0.056 & P-1 C_{i}^{1} #2 (i^15) & aP30 & None &
    ↪ CaO3Si & Wollastonite & M. J. Buerger and C. T. Prewitt, Proc.
    ↪ Natl. Acad. Sci. 47, 1884-1888 (1961)
   1.00000000000000
   7.94000000000000    0.00000000000000    0.00000000000000
  -1.70053661868821    7.11973140002490    0.00000000000000
```

```
  −0.66125116738401   −0.16216723254653    7.03714066097314
     Ca      O      Si
      6     18      6
Direct
    0.19850000000000    0.42280000000000    0.76080000000000   Ca   (2i)
   −0.19850000000000   −0.42280000000000   −0.76080000000000   Ca   (2i)
    0.20270000000000   −0.07070000000000    0.76400000000000   Ca   (2i)
   −0.20270000000000    0.07070000000000   −0.76400000000000   Ca   (2i)
    0.49660000000000    0.24950000000000    0.47200000000000   Ca   (2i)
   −0.49660000000000   −0.24950000000000   −0.47200000000000   Ca   (2i)
    0.42910000000000    0.23140000000000    0.80190000000000   O    (2i)
   −0.42910000000000   −0.23140000000000   −0.80190000000000   O    (2i)
    0.40080000000000    0.72590000000000    0.83020000000000   O    (2i)
   −0.40080000000000   −0.72590000000000   −0.83020000000000   O    (2i)
    0.30370000000000    0.46350000000000    0.46410000000000   O    (2i)
   −0.30370000000000   −0.46350000000000   −0.46410000000000   O    (2i)
    0.30170000000000   −0.06260000000000    0.46550000000000   O    (2i)
   −0.30170000000000    0.06260000000000   −0.46550000000000   O    (2i)
    0.01540000000000    0.62540000000000    0.73430000000000   O    (2i)
   −0.01540000000000   −0.62540000000000   −0.73430000000000   O    (2i)
    0.01750000000000    0.13190000000000    0.73530000000000   O    (2i)
   −0.01750000000000   −0.13190000000000   −0.73530000000000   O    (2i)
    0.27320000000000    0.51180000000000    0.09190000000000   O    (2i)
   −0.27320000000000   −0.51180000000000   −0.09190000000000   O    (2i)
    0.27130000000000    0.87170000000000    0.09400000000000   O    (2i)
   −0.27130000000000   −0.87170000000000   −0.09400000000000   O    (2i)
    0.21880000000000    0.17840000000000    0.22280000000000   O    (2i)
   −0.21880000000000   −0.17840000000000   −0.22280000000000   O    (2i)
    0.18520000000000    0.38700000000000    0.26870000000000   Si   (2i)
   −0.18520000000000   −0.38700000000000   −0.26870000000000   Si   (2i)
    0.18490000000000   −0.04550000000000    0.26920000000000   Si   (2i)
   −0.18490000000000    0.04550000000000   −0.26920000000000   Si   (2i)
    0.39700000000000    0.72350000000000    0.05600000000000   Si   (2i)
   −0.39700000000000   −0.72350000000000   −0.05600000000000   Si   (2i)
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Albite ($NaAlSi_3O_8$, $S6_8$): ABC8D3_aP26_2_i_i_8i_3i - CIF

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_chemical_formula_sum 'Al Na O8 Si3'

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 'J. V. Smith'
 'G. Artioli'
 '{\AA}. Kvick'
_journal_name_full_name
;
 American Mineralogist
;
_journal_volume 71
_journal_year 1986
_journal_page_first 727
_journal_page_last 733
_publ_Section_title
;
 Low albite, NaAlSi$_{3}$O$_{8}$: Neutron diffraction study of crystal
      ↪ structure at 13~K
;

_aflow_title 'Albite (NaAlSi$_{3}$O$_{8}$, $S6_{8}$) Structure'
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      ↪ 107.37724,100.55864,0.20773,0.84122,0.1767,0.14441,0.27518,
      ↪ 0.25422,-0.03356,0.87658,0.13472,0.72111,0.40891,0.41381,
      ↪ 0.80981,0.2953,0.07942,0.74076,0.0321,0.32928,0.26814,0.70828,
      ↪ 0.31212,0.22592,0.32857,0.71731,0.38942,0.10112,0.31914,0.43708
      ↪ ,0.31682,0.05324,0.23701,0.18228,0.8249,0.68677,0.41955,0.19963
      ↪ ,0.6404,0.203,0.4392'
_aflow_Strukturbericht '$S6_{8}$'
_aflow_Pearson 'aP26'

_symmetry_space_group_name_H-M "P -1"
_symmetry_Int_Tables_number 2

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_cell_angle_alpha  115.11511
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_atom_site_Wyckoff_label
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Na1  Na  2  i  0.14441  0.27518  0.25422  1.00000
O1   O   2  i -0.03356  0.87658  0.13472  1.00000
O2   O   2  i  0.72111  0.40891  0.41381  1.00000
O3   O   2  i  0.80981  0.29530  0.07942  1.00000
O4   O   2  i  0.74076  0.03210  0.32928  1.00000
O5   O   2  i  0.26814  0.70828  0.31212  1.00000
O6   O   2  i  0.22592  0.32857  0.71731  1.00000
O7   O   2  i  0.38942  0.10112  0.31914  1.00000
O8   O   2  i  0.43708  0.31682  0.05324  1.00000
Si1  Si  2  i  0.23701  0.18228  0.82490  1.00000
Si2  Si  2  i  0.68677  0.41955  0.19963  1.00000
Si3  Si  2  i  0.64040  0.20300  0.43920  1.00000
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Albite ($NaAlSi_3O_8$, $S6_8$): ABC8D3_aP26_2_i_i_8i_3i - POSCAR

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ABC8D3_aP26_2_i_i_8i_3i & a,b/a,c/a,alpha,beta,gamma,x1,y1,z1,x2,y2,z2,
      ↪ x3,y3,z3,x4,y4,z4,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,
      ↪ x10,y10,z10,x11,y11,z11,x12,y12,z12,x13,y13,z13 --params=7.1576
      ↪ ,1.03716189784,1.07544707723,115.11511,107.37724,100.55864,
      ↪ 0.20773,0.84122,0.1767,0.14441,0.27518,0.25422,-0.03356,0.87658
      ↪ ,0.13472,0.72111,0.40891,0.41381,0.80981,0.2953,0.07942,0.74076
      ↪ ,0.0321,0.32928,0.26814,0.70828,0.31212,0.22592,0.32857,0.71731
      ↪ ,0.38942,0.10112,0.31914,0.43708,0.31682,0.05324,0.23701,
      ↪ 0.18228,0.8249,0.68677,0.41955,0.19963,0.6404,0.203,0.4392 &
      ↪ P-1 C_{i}^{1} #2 (i^13) & aP26 & $S6_{8}$ & AlNaO8Si3 & Albite
      ↪ & J. V. Smith and G. Artioli and {\AA}. Kvick, Am. Mineral. 71,
      ↪ 727-733 (1986)
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   7.15760000000000    0.00000000000000    0.00000000000000
  -1.36031164512511    7.29789289564029    0.00000000000000
  -2.29898434571175   -3.75196276024920    6.31591640922167
     Al      Na      O      Si
      2       2     16       6
Direct
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   -0.20773000000000   -0.84122000000000   -0.17670000000000   Al   (2i)
    0.14441000000000    0.27518000000000    0.25422000000000   Na   (2i)
   -0.14441000000000   -0.27518000000000   -0.25422000000000   Na   (2i)
   -0.03356000000000    0.87658000000000    0.13472000000000   O    (2i)
    0.03356000000000   -0.87658000000000   -0.13472000000000   O    (2i)
    0.72111000000000    0.40891000000000    0.41381000000000   O    (2i)
   -0.72111000000000   -0.40891000000000   -0.41381000000000   O    (2i)
    0.80981000000000    0.29530000000000    0.07942000000000   O    (2i)
   -0.80981000000000   -0.29530000000000   -0.07942000000000   O    (2i)
    0.74076000000000    0.03210000000000    0.32928000000000   O    (2i)
   -0.74076000000000   -0.03210000000000   -0.32928000000000   O    (2i)
    0.26814000000000    0.70828000000000    0.31212000000000   O    (2i)
   -0.26814000000000   -0.70828000000000   -0.31212000000000   O    (2i)
    0.22592000000000    0.32857000000000    0.71731000000000   O    (2i)
   -0.22592000000000   -0.32857000000000   -0.71731000000000   O    (2i)
    0.38942000000000    0.10112000000000    0.31914000000000   O    (2i)
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    0.43708000000000    0.31682000000000    0.05324000000000   O    (2i)
   -0.43708000000000   -0.31682000000000   -0.05324000000000   O    (2i)
    0.23701000000000    0.18228000000000    0.82490000000000   Si   (2i)
   -0.23701000000000   -0.18228000000000   -0.82490000000000   Si   (2i)
    0.68677000000000    0.41955000000000    0.19963000000000   Si   (2i)
   -0.68677000000000   -0.41955000000000   -0.19963000000000   Si   (2i)
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   -0.64040000000000   -0.20300000000000   -0.43920000000000   Si   (2i)
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TaTi (BCC SQS-16): AB_aP16_2_4i_4i - CIF

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# AFLOW.org Repositories
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      ↪ 0.875 SG=2 [ANRL doi: 10.1016/j.commatsci.2017.01.017 (part 1),
      ↪ doi: 10.1016/j.commatsci.2018.10.043 (part 2)]
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_chemical_name_mineral 'TaTi'
_chemical_formula_sum 'Ta Ti'

loop_
_publ_author_name
 'C. Jiang'
 'C. Wolverton'
 'J. Sofo'
 'L.-Q. Chen'
 'Z.-K. Liu'
_journal_name_full_name
;
 Physical Review B
;
_journal_volume 69
_journal_year 2004
_journal_page_first 214202
_journal_page_last 214202
_publ_Section_title
;
 First-principles study of binary bcc alloys using special quasirandom
      ↪ structures
;

_aflow_title 'TaTi (BCC SQS-16) Structure'
_aflow_proto 'AB_aP16_2_4i_4i'
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      ↪ ,0.0,0.375,0.375,0.5,0.875,0.625,0.75,0.625,0.875,0.5,0.375,
      ↪ 0.125,0.75,0.125,0.875,0.75,0.875,0.375,0.0,0.875'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'aP16'

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_cell_length_b 5.6656530442
_cell_length_c 9.6759446334
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_cell_angle_gamma 79.9750121379
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_symmetry_Int_Tables_Number 2
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 2 -x,-y,-z
loop_
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 _atom_site_occupancy
 _atom_site_fract_x
 _atom_site_fract_y
 _atom_site_fract_z
 _atom_site_thermal_displace_type
 _atom_site_B_iso_or_equiv
 _atom_site_type_symbol
 _atom_site_symmetry_multiplicity
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Ta2 1.0000000000 0.8750000000 0.0000000000 0.3750000000 Biso 1.0 Ta 2 i
Ta3 1.0000000000 0.3750000000 0.5000000000 0.8750000000 Biso 1.0 Ta 2 i
Ta4 1.0000000000 0.6250000000 0.7500000000 0.6250000000 Biso 1.0 Ta 2 i
Ti1 1.0000000000 0.8750000000 0.5000000000 0.3750000000 Biso 1.0 Ti 2 i
Ti2 1.0000000000 0.1250000000 0.7500000000 0.1250000000 Biso 1.0 Ti 2 i
Ti3 1.0000000000 0.8750000000 0.7500000000 0.8750000000 Biso 1.0 Ti 2 i
Ti4 1.0000000000 0.3750000000 0.0000000000 0.8750000000 Biso 1.0 Ti 2 i
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TaTi (BCC SQS-16): AB_aP16_2_4i_4i - POSCAR

```
AB_aP16_2_4i_4i & a,b/a,c/a,alpha,beta,gamma,x1,y1,z1,x2,y2,z2,x3,y3,z3,
      ↪ x4,y4,z4,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8 --params=
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      ↪ 87.0786711125,79.9750121379,0.375,0.75,0.375,0.875,0.0,0.375,
      ↪ 0.375,0.5,0.875,0.625,0.75,0.625,0.875,0.5,0.375,0.125,0.75,
      ↪ 0.125,0.875,0.75,0.875,0.375,0.0,0.875 & P-1 C_{i}^{1} #2 (i^8)
      ↪ & aP16 & None & TaTi & TaTi & C. Jiang et al., Phys. Rev. B 69
      ↪ , 214202(2004)
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   0.98626360139496    5.57914944465697    0.00000000000000
   0.49313180069476    2.78957472230357    9.25197267854001
     Ta      Ti
      8       8
Direct
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   -0.37500000000000   -0.75000000000000   -0.37500000000000   Ta   (2i)
    0.87500000000000    0.00000000000000    0.37500000000000   Ta   (2i)
   -0.87500000000000    0.00000000000000   -0.37500000000000   Ta   (2i)
    0.37500000000000    0.50000000000000    0.87500000000000   Ta   (2i)
   -0.37500000000000   -0.50000000000000   -0.87500000000000   Ta   (2i)
    0.62500000000000    0.75000000000000    0.62500000000000   Ta   (2i)
   -0.62500000000000   -0.75000000000000   -0.62500000000000   Ta   (2i)
    0.87500000000000    0.50000000000000    0.37500000000000   Ti   (2i)
   -0.87500000000000   -0.50000000000000   -0.37500000000000   Ti   (2i)
    0.12500000000000    0.75000000000000    0.12500000000000   Ti   (2i)
   -0.12500000000000   -0.75000000000000   -0.12500000000000   Ti   (2i)
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$W_2O_3(PO_4)_2$: A11B2C2_mP60_4_22a_4a_4a - CIF

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 'S. {\AA}sbrink'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 18
_journal_year 1964
_journal_page_first 2329
_journal_page_last 2338
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;
 The Crystal Structure of W$_{2}$O$_{3}$(PO$_{4}$)$_{2}$. Determination
      ↪ of a Superstructure by Means of Least-Squares Calculations
;

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      ↪ 0.722,0.652,0.239,0.082,0.326,0.287,-0.078,0.41,0.01,0.888,
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      ↪ 0.841,0.669,0.64,0.155,0.343,0.3226,0.0046,0.0737,0.8943,-
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_aflow_Pearson 'mP60'

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2 -x,y+1/2,-z

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_atom_site_fract_x
_atom_site_fract_y
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_atom_site_occupancy
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O2 O 2 a 0.79600 0.88900 0.81100 1.00000
O3 O 2 a 0.25800 -0.09600 0.12900 1.00000
O4 O 2 a 0.83900 0.07900 0.84200 1.00000
O5 O 2 a 0.19300 0.00900 0.44200 1.00000
O6 O 2 a 0.85200 -0.00500 0.56100 1.00000
O7 O 2 a 0.83400 0.01700 0.19700 1.00000
O8 O 2 a 0.08600 -0.02400 0.78800 1.00000
O9 O 2 a 0.51700 -0.01200 0.57000 1.00000
O10 O 2 a 0.99100 0.22900 0.99100 1.00000
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O12 O 2 a -0.08900 0.30000 0.66500 1.00000
O13 O 2 a 0.42100 0.26300 0.25700 1.00000
O14 O 2 a 0.58400 0.26700 0.72200 1.00000
O15 O 2 a 0.65200 0.23900 0.08200 1.00000
O16 O 2 a 0.32600 0.28700 -0.07800 1.00000
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O19 O 2 a 0.27900 0.16200 0.66400 1.00000
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O22 O 2 a 0.64000 0.15500 0.34300 1.00000
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P4 P 2 a 0.38200 0.26060 0.75140 1.00000
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W2 W 2 a 0.70240 -0.00710 0.35850 1.00000
W3 W 2 a 0.19130 0.25010 0.15330 1.00000
W4 W 2 a 0.80740 0.25570 0.84620 1.00000
```

$W_2O_3(PO_4)_2$: A11B2C2_mP60_4_22a_4a_4a - POSCAR

```
A11B2C2_mP60_4_22a_4a_4a & a,b/a,c/a,beta,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,
      ↪ y4,z4,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10,
      ↪ x11,y11,z11,x12,y12,z12,x13,y13,z13,x14,y14,z14,x15,y15,z15,x16
      ↪ ,y16,z16,x17,y17,z17,x18,y18,z18,x19,y19,z19,x20,y20,z20,x21,
      ↪ y21,z21,x22,y22,z22,x23,y23,z23,x24,y24,z24,x25,y25,z25,x26,y26
      ↪ ,z26,x27,y27,z27,x28,y28,z28,x29,y29,z29,x30,y30,z30 --params=
      ↪ 7.822,1.59805676298,0.991306571209,91.05,0.224,0.099,0.084,
      ↪ 0.796,0.889,0.811,0.258,-0.096,0.129,0.839,0.079,0.842,0.193,
      ↪ 0.009,0.442,0.852,-0.005,0.561,0.834,0.017,0.197,0.086,-0.024,
      ↪ 0.788,0.517,-0.012,0.57,0.991,0.229,0.991,0.088,0.243,0.331,-
      ↪ 0.089,0.3,0.665,0.421,0.263,0.257,0.584,0.267,0.722,0.652,0.239
      ↪ ,0.082,0.326,0.287,-0.078,0.41,0.01,0.888,0.497,0.007,0.208,
      ↪ 0.279,0.162,0.664,0.687,0.842,0.389,0.301,0.841,0.669,0.64,
      ↪ 0.155,0.343,0.3226,0.0046,0.0737,0.8943,-0.0084,0.7508,0.6178,
      ↪ 0.2582,0.2607,0.382,0.2606,0.7514,0.3082,-0.0024,0.6351,0.7024
      ↪ ,-0.0071,0.3585,0.1913,0.2501,0.1533,0.8074,0.2557,0.8462 & P2_
      ↪ {1} C_{2}^{2} #4 (a^30) & mP60 & None & O11P2W2 & O11P2W2 & P.
      ↪ Kierkegaard and S. {\AA}sbrink, Acta Chem. Scand. 18, 2329-2338
      ↪ (1964)
   1.00000000000000
   7.82200000000000    0.00000000000000    0.00000000000000
   0.00000000000000   12.50000000000000    0.00000000000000
  -0.14209151804538    0.00000000000000    7.75269798202533
     O       P       W
    44       8       8
Direct
    0.22400000000000    0.09900000000000    0.08400000000000     O     (2a)
```

```
 −0.22400000000000   0.59900000000000  −0.08400000000000   O   (2a)
  0.79600000000000   0.88900000000000   0.81100000000000   O   (2a)
 −0.79600000000000   1.38900000000000  −0.81100000000000   O   (2a)
  0.25800000000000  −0.09600000000000   0.12900000000000   O   (2a)
 −0.25800000000000   0.40400000000000  −0.12900000000000   O   (2a)
  0.83900000000000   0.07900000000000   0.84200000000000   O   (2a)
 −0.83900000000000   0.57900000000000  −0.84200000000000   O   (2a)
  0.19300000000000   0.00900000000000   0.44200000000000   O   (2a)
 −0.19300000000000   0.50900000000000  −0.44200000000000   O   (2a)
  0.85200000000000  −0.00500000000000   0.56100000000000   O   (2a)
 −0.85200000000000   0.49500000000000  −0.56100000000000   O   (2a)
  0.83400000000000   0.01700000000000   0.19700000000000   O   (2a)
 −0.83400000000000   0.51700000000000  −0.19700000000000   O   (2a)
  0.08600000000000  −0.02400000000000   0.78800000000000   O   (2a)
 −0.08600000000000   0.47600000000000  −0.78800000000000   O   (2a)
  0.51700000000000  −0.01200000000000   0.57000000000000   O   (2a)
 −0.51700000000000   0.48800000000000  −0.57000000000000   O   (2a)
  0.99100000000000   0.22900000000000   0.99100000000000   O   (2a)
 −0.99100000000000   0.72900000000000  −0.99100000000000   O   (2a)
  0.08800000000000   0.24300000000000   0.33100000000000   O   (2a)
 −0.08800000000000   0.74300000000000  −0.33100000000000   O   (2a)
 −0.08900000000000   0.30000000000000   0.66500000000000   O   (2a)
  0.08900000000000   0.80000000000000  −0.66500000000000   O   (2a)
  0.42100000000000   0.26300000000000   0.25700000000000   O   (2a)
 −0.42100000000000   0.76300000000000  −0.25700000000000   O   (2a)
  0.58400000000000   0.26700000000000   0.72200000000000   O   (2a)
 −0.58400000000000   0.76700000000000  −0.72200000000000   O   (2a)
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 −0.65200000000000   0.73900000000000  −0.08200000000000   O   (2a)
  0.32600000000000   0.28700000000000  −0.07800000000000   O   (2a)
 −0.32600000000000   0.78700000000000   0.07800000000000   O   (2a)
  0.41000000000000   0.01000000000000   0.88800000000000   O   (2a)
 −0.41000000000000   0.51000000000000  −0.88800000000000   O   (2a)
  0.49700000000000   0.00700000000000   0.20800000000000   O   (2a)
 −0.49700000000000   0.50700000000000  −0.20800000000000   O   (2a)
  0.27900000000000   0.16200000000000   0.66400000000000   O   (2a)
 −0.27900000000000   0.66200000000000  −0.66400000000000   O   (2a)
  0.68700000000000   0.84200000000000   0.38900000000000   O   (2a)
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  0.30100000000000   0.84100000000000   0.66900000000000   O   (2a)
 −0.30100000000000   1.34100000000000  −0.66900000000000   O   (2a)
  0.64000000000000   0.15500000000000   0.34300000000000   O   (2a)
 −0.64000000000000   0.65500000000000  −0.34300000000000   O   (2a)
  0.32260000000000   0.00460000000000   0.07370000000000   P   (2a)
 −0.32260000000000   0.50460000000000  −0.07370000000000   P   (2a)
  0.89430000000000  −0.00840000000000   0.75080000000000   P   (2a)
 −0.89430000000000   0.49160000000000  −0.75080000000000   P   (2a)
  0.61780000000000   0.25820000000000   0.26070000000000   P   (2a)
 −0.61780000000000   0.75820000000000  −0.26070000000000   P   (2a)
  0.38200000000000   0.26060000000000   0.75140000000000   P   (2a)
 −0.38200000000000   0.76060000000000  −0.75140000000000   P   (2a)
  0.30820000000000  −0.00240000000000   0.63510000000000   W   (2a)
 −0.30820000000000   0.49760000000000  −0.63510000000000   W   (2a)
  0.70240000000000  −0.00710000000000   0.35850000000000   W   (2a)
 −0.70240000000000   0.49290000000000  −0.35850000000000   W   (2a)
  0.19130000000000   0.25010000000000   0.15330000000000   W   (2a)
 −0.19130000000000   0.75010000000000  −0.15330000000000   W   (2a)
  0.80740000000000   0.25570000000000   0.84620000000000   W   (2a)
 −0.80740000000000   0.75570000000000  −0.84620000000000   W   (2a)
```

$Li_2SO_4{\cdot}H_2O$ ($H4_8$): A2B2C5D_mP20_4_2a_2a_5a_a - CIF

```
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_audit_creation_method FINDSYM

_chemical_name_mineral 'H2Li2O5S'
_chemical_formula_sum 'H2 Li2 O5 S'

loop_
_publ_author_name
 'J.-O. Lundgren'
 '{\AA}. Kvick'
 'M. Karppinen'
 'R. Liminga'
 'S. C. Abrahams'
_journal_name_full_name
;
 Journal of Chemical Physics
;
_journal_volume 80
_journal_year 1984
_journal_page_first 423
_journal_page_last 430
_publ_Section_title
;
 Neutron diffraction structural study of pyroelectric Li$_{2}$SO$_{4}$$$\
      ↪ cdot$H$_{2}$O at 293, 80, and 20 K
;

_aflow_title 'Li$_{2}$SO$_{4}$$$\cdot$H$_{2}$O ($H4_{8}$) Structure'
_aflow_proto 'A2B2C5D_mP20_4_2a_2a_5a_a'
_aflow_params 'a,b/a,c/a,\beta,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3
      ↪ },y_{3},z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},
      ↪ z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{
      ↪ 10},y_{10},z_{10}'
_aflow_params_values '5.449,0.886768214351,1.49330152322,107.19,-0.02933
      ↪ ,0.38995,0.31142,0.00673,0.63149,0.44573,0.3031,0.49638,-
      ↪ 0.00741,0.56046,0.48968,0.39524,0.02087,0.07409,0.17122,0.43724
      ↪ ,0.11111,0.37857,0.39807,0.12076,0.07634,0.32462,0.69627,
      ↪ 0.20975,-0.08635,0.46436,0.40379,0.29229,0.0,0.20805'
_aflow_Strukturbericht '$H4_{8}$'
_aflow_Pearson 'mP20'

_symmetry_space_group_name_H-M "P 1 21 1"
_symmetry_Int_Tables_number 4

_cell_length_a      5.44900
_cell_length_b      4.83200
_cell_length_c      8.13700
_cell_angle_alpha  90.00000
_cell_angle_beta   107.19000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
H1 H 2 a -0.02933 0.38995 0.31142 1.00000
H2 H 2 a 0.00673 0.63149 0.44573 1.00000
Li1 Li 2 a 0.30310 0.49638 -0.00741 1.00000
Li2 Li 2 a 0.56046 0.48968 0.39524 1.00000
O1 O 2 a 0.02087 0.07409 0.17122 1.00000
O2 O 2 a 0.43724 0.11111 0.37857 1.00000
O3 O 2 a 0.39807 0.12076 0.07634 1.00000
O4 O 2 a 0.32462 0.69627 0.20975 1.00000
O5 O 2 a -0.08635 0.46436 0.40379 1.00000
S1 S 2 a 0.29229 0.00000 0.20805 1.00000
```

$Li_2SO_4{\cdot}H_2O$ ($H4_8$): A2B2C5D_mP20_4_2a_2a_5a_a - POSCAR

```
A2B2C5D_mP20_4_2a_2a_5a_a & a,b/a,c/a,beta,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4
      ↪ ,y4,z4,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10
      ↪  --params=5.449,0.886768214351,1.49330152322,107.19,-0.02933,
      ↪ 0.38995,0.31142,0.00673,0.63149,0.44573,0.3031,0.49638,-0.00741
      ↪ ,0.56046,0.48968,0.39524,0.02087,0.07409,0.17122,0.43724,
      ↪ 0.11111,0.37857,0.39807,0.12076,0.07634,0.32462,0.69627,0.20975
      ↪ ,-0.08635,0.46436,0.40379,0.29229,0.0,0.20805 & P2_{1} C_{2}^{2
      ↪ } #4 (a^10) & mP20 & $H4_{8}$ & H2Li2O5S & H2Li2O5S & J.-O.
      ↪ Lundgren et al., J. Chem. Phys. 80, 423-430 (1984)
  1.00000000000000
  5.44900000000000   0.00000000000000   0.00000000000000
  0.00000000000000   4.83200000000000   0.00000000000000
 -2.40481970468058   0.00000000000000   7.77351987120249
     H     Li     O     S
     4      4     10     2
Direct
 -0.02933000000000   0.38995000000000   0.31142000000000   H   (2a)
  0.02933000000000   0.88995000000000  -0.31142000000000   H   (2a)
  0.00673000000000   0.63149000000000   0.44573000000000   H   (2a)
 -0.00673000000000   1.13149000000000  -0.44573000000000   H   (2a)
  0.30310000000000   0.49638000000000  -0.00741000000000   Li  (2a)
 -0.30310000000000   0.99638000000000   0.00741000000000   Li  (2a)
  0.56046000000000   0.48968000000000   0.39524000000000   Li  (2a)
 -0.56046000000000   0.98968000000000  -0.39524000000000   Li  (2a)
  0.02087000000000   0.07409000000000   0.17122000000000   O   (2a)
 -0.02087000000000   0.57409000000000  -0.17122000000000   O   (2a)
  0.43724000000000   0.11111000000000   0.37857000000000   O   (2a)
 -0.43724000000000   0.61111000000000  -0.37857000000000   O   (2a)
  0.39807000000000   0.12076000000000   0.07634000000000   O   (2a)
 -0.39807000000000   0.62076000000000  -0.07634000000000   O   (2a)
  0.32462000000000   0.69627000000000   0.20975000000000   O   (2a)
 -0.32462000000000   1.19627000000000  -0.20975000000000   O   (2a)
 -0.08635000000000   0.46436000000000   0.40379000000000   O   (2a)
  0.08635000000000   0.96436000000000  -0.40379000000000   O   (2a)
  0.29229000000000   0.00000000000000   0.20805000000000   S   (2a)
 -0.29229000000000   0.50000000000000  -0.20805000000000   S   (2a)
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Ca_3UO_6: A3B6C_mP20_4_3a_6a_a - CIF

```
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_audit_creation_method FINDSYM

_chemical_name_mineral 'Ca3O6U'
_chemical_formula_sum 'Ca3 O6 U'

loop_
_publ_author_name
 'B. O. Loopstra'
 'H. M. Rietveld'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 25
_journal_year 1969
_journal_page_first 787
_journal_page_last 791
_publ_Section_title
;
 The structure of some alkaline-earth metal uranates
;

_aflow_title 'Ca$_{3}$UO$_{6}$ Structure'
_aflow_proto 'A3B6C_mP20_4_3a_6a_a'
_aflow_params 'a,b/a,c/a,\beta,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3
      ↪ },y_{3},z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},
```

```
      ↪ z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{
      ↪ 10},y_{10},z_{10}'
_aflow_params_values '5.7275,1.03996508075,1.44883457006,90.568,0.2621,-
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      ↪ 0.41,0.5173,0.5605,0.1618,0.1938,0.0747,0.2744,0.3323,0.1165,
      ↪ 0.0484,0.0152,0.0327,0.3053,0.6942,0.574,0.1963,0.8252,0.2459,
      ↪ 0.0,0.2519'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mP20'

_symmetry_space_group_name_H-M "P 1 21 1"
_symmetry_Int_Tables_number 4

_cell_length_a      5.72750
_cell_length_b      5.95640
_cell_length_c      8.29820
_cell_angle_alpha  90.00000
_cell_angle_beta   90.56800
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ca1 Ca 2 a 0.26210 -0.00250 0.74880 1.00000
Ca2 Ca 2 a 0.27230 0.42230 0.01230 1.00000
Ca3 Ca 2 a 0.75320 0.02880 0.47610 1.00000
O1 O 2 a 0.63190 0.41000 0.51730 1.00000
O2 O 2 a 0.56050 0.16180 0.19380 1.00000
O3 O 2 a 0.07470 0.27440 0.33230 1.00000
O4 O 2 a 0.11650 0.04840 0.01520 1.00000
O5 O 2 a 0.03270 0.30530 0.69420 1.00000
O6 O 2 a 0.57400 0.19630 0.82520 1.00000
U1 U 2 a 0.24590 0.00000 0.25190 1.00000
```

Ca_3UO_6: A3B6C_mP20_4_3a_6a_a - POSCAR

```
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      ↪ z4,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10 --
      ↪ params=5.7275,1.03996508075,1.44883457006,90.568,0.2621,-0.0025
      ↪ ,0.7488,0.2723,0.4223,0.0123,0.7532,0.0288,0.4761,0.6319,0.41,
      ↪ 0.5173,0.5605,0.1618,0.1938,0.0747,0.2744,0.3323,0.1165,0.0484,
      ↪ 0.0152,0.0327,0.3053,0.6942,0.574,0.1963,0.8252,0.2459,0.0,
      ↪ 0.2519 & P2_{1} C_{2}^{2} #4 (a^10) & mP20 & None & Ca3O6U &
      ↪ Ca3O6U & B. O. Loopstra and H. M. Rietveld, Acta Crystallogr.
      ↪ Sect. B Struct. Sci. 25, 787-791 (1969)
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   5.72750000000000    0.00000000000000    0.00000000000000
   0.00000000000000    5.95640000000000    0.00000000000000
  -0.08226261057537    0.00000000000000    8.29779224269331
     Ca      O      U
      6     12      2
Direct
   0.26210000000000   -0.00250000000000    0.74880000000000    Ca    (2a)
  -0.26210000000000    0.49750000000000   -0.74880000000000    Ca    (2a)
   0.27230000000000    0.42230000000000    0.01230000000000    Ca    (2a)
  -0.27230000000000    0.92230000000000   -0.01230000000000    Ca    (2a)
   0.75320000000000    0.02880000000000    0.47610000000000    Ca    (2a)
  -0.75320000000000    0.52880000000000   -0.47610000000000    Ca    (2a)
   0.63190000000000    0.41000000000000    0.51730000000000    O     (2a)
  -0.63190000000000    0.91000000000000   -0.51730000000000    O     (2a)
   0.56050000000000    0.16180000000000    0.19380000000000    O     (2a)
  -0.56050000000000    0.66180000000000   -0.19380000000000    O     (2a)
   0.07470000000000    0.27440000000000    0.33230000000000    O     (2a)
  -0.07470000000000    0.77440000000000   -0.33230000000000    O     (2a)
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  -0.11650000000000    0.54840000000000   -0.01520000000000    O     (2a)
   0.03270000000000    0.30530000000000    0.69420000000000    O     (2a)
  -0.03270000000000    0.80530000000000   -0.69420000000000    O     (2a)
   0.57400000000000    0.19630000000000    0.82520000000000    O     (2a)
  -0.57400000000000    0.69630000000000   -0.82520000000000    O     (2a)
   0.24590000000000    0.00000000000000    0.25190000000000    U     (2a)
  -0.24590000000000    0.50000000000000   -0.25190000000000    U     (2a)
```

Bassanite $[CaSO_4(H_2O)_{0.5}, H4_7]$: A2B2C9D2_mC90_5_ab2c_3c_b13c_3c - CIF

```
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_audit_creation_method FINDSYM

_chemical_name_mineral 'Bassanite'
_chemical_formula_sum 'Ca2 H2 O9 S2'

loop_
_publ_author_name
 'W. Abriel'
 'R. Nesper'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 205
_journal_year 1993
_journal_page_first 99
_journal_page_last 113
_publ_Section_title
;
 Bestimmung der Kristallstruktur von CaSO$_{4}$(H$_{2}$O)$_{0.5}$ mit R
      ↪ \"{o}ntgenbeugungsmethoden und mit Potentialprofil-Rechnungen
;

# Found in The monoclinic $I2$ structure of bassanite, calcium sulphate
      ↪ hemihydrate (CaSO$_{4}$$\cdot$0.5H$_{2}$O), 2001

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      ↪ z_{8},x_{9},y_{9},z_{9},x_{10},y_{10},z_{10},x_{11},y_{11},z_{
      ↪ 11},x_{12},y_{12},z_{12},x_{13},y_{13},z_{13},x_{14},y_{14},z_{
      ↪ 14},x_{15},y_{15},z_{15},x_{16},y_{16},z_{16},x_{17},y_{17},z_{
      ↪ 17},x_{18},y_{18},z_{18},x_{19},y_{19},z_{19},x_{20},y_{20},z_{
      ↪ 20},x_{21},y_{21},z_{21},x_{22},y_{22},z_{22},x_{23},y_{23},z_{
      ↪ 23},x_{24},y_{24},z_{24}'
_aflow_params_values '17.45812,0.39701869388,0.688934432803,133.36555,
      ↪ 0.4553,-0.0447,0.5985,0.1667,0.2724,0.8943,0.3333,0.2724,0.6057
      ↪ ,0.5,0.171,0.572,0.333,0.307,0.783,0.833,0.023,0.712,0.5173,
      ↪ 0.226,0.8882,0.0173,0.226,0.3882,0.8507,0.4435,0.5522,0.3507,
      ↪ 0.4435,0.0522,0.184,0.3306,0.1115,0.684,0.3306,0.6115,0.6556,
      ↪ 0.1185,-0.0982,0.1556,0.1185,0.4018,-0.0111,0.3102,0.8065,
      ↪ 0.4889,0.3102,0.3065,0.3222,0.0714,0.7583,0.8222,0.0714,0.2583,
      ↪ 0.3333,0.4508,0.784,0.0833,0.2752,0.3585,0.25,0.4496,0.25,
      ↪ 0.4167,0.2752,0.1415'
_aflow_Strukturbericht '$H4_{7}$'
_aflow_Pearson 'mC90'

_symmetry_space_group_name_H-M "C 1 2 1"
_symmetry_Int_Tables_number 5

_cell_length_a      17.45812
_cell_length_b      6.93120
_cell_length_c      12.02750
_cell_angle_alpha  90.00000
_cell_angle_beta   133.36555
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z
3 x+1/2,y+1/2,z
4 -x+1/2,y+1/2,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ca1 Ca 2 a 0.00000 0.45530 0.00000 1.00000
Ca2 Ca 2 b 0.00000 -0.04470 0.50000 1.00000
O1 O 2 b 0.00000 0.59850 0.50000 1.00000
Ca3 Ca 4 c 0.16670 0.27240 0.89430 1.00000
Ca4 Ca 4 c 0.33330 0.27240 0.60570 1.00000
H1 H 4 c 0.50000 0.17100 0.57200 1.00000
H2 H 4 c 0.33300 0.30700 0.78300 1.00000
H3 H 4 c 0.83300 0.02300 0.71200 1.00000
O2 O 4 c 0.51730 0.22600 0.88820 1.00000
O3 O 4 c 0.01730 0.22600 0.38820 1.00000
O4 O 4 c 0.85070 0.44350 0.55220 1.00000
O5 O 4 c 0.35070 0.44350 0.05220 1.00000
O6 O 4 c 0.18400 0.33060 0.11150 1.00000
O7 O 4 c 0.68400 0.33060 0.61150 1.00000
O8 O 4 c 0.65560 0.11850 -0.09820 1.00000
O9 O 4 c 0.15560 0.11850 0.40180 1.00000
O10 O 4 c -0.01110 0.31020 0.80650 1.00000
O11 O 4 c 0.48890 0.31020 0.30650 1.00000
O12 O 4 c 0.32220 0.07140 0.75830 1.00000
O13 O 4 c 0.82220 0.07140 0.25830 1.00000
O14 O 4 c 0.33330 0.45080 0.78400 1.00000
S1 S 4 c 0.08330 0.27520 0.35850 1.00000
S2 S 4 c 0.25000 0.44960 0.25000 1.00000
S3 S 4 c 0.41670 0.27520 0.14150 1.00000
```

Bassanite $[CaSO_4(H_2O)_{0.5}, H4_7]$: A2B2C9D2_mC90_5_ab2c_3c_b13c_3c - POSCAR

```
A2B2C9D2_mC90_5_ab2c_3c_b13c_3c & a,b/a,c/a,beta,y1,y2,y3,x4,y4,z4,x5,y5
      ↪ ,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10,x11,y11,z11
      ↪ ,x12,y12,z12,x13,y13,z13,x14,y14,z14,x15,y15,z15,x16,y16,z16,
      ↪ x17,y17,z17,x18,y18,z18,x19,y19,z19,x20,y20,z20,x21,y21,z21,x22
      ↪ ,y22,z22,x23,y23,z23,x24,y24,z24 --params=17.45812,
      ↪ 0.39701869388,0.688934432803,133.36555,0.4553,-0.0447,0.5985,
      ↪ 0.1667,0.2724,0.8943,0.3333,0.2724,0.6057,0.5,0.171,0.572,0.333
      ↪ ,0.307,0.783,0.833,0.023,0.712,0.5173,0.226,0.8882,0.0173,0.226
      ↪ ,0.3882,0.8507,0.4435,0.5522,0.3507,0.4435,0.0522,0.184,0.3306,
      ↪ 0.1115,0.684,0.3306,0.6115,0.6556,0.1185,-0.0982,0.1556,0.1185,
      ↪ 0.4018,-0.0111,0.3102,0.8065,0.4889,0.3102,0.3065,0.3222,0.0714
      ↪ ,0.7583,0.8222,0.0714,0.2583,0.3333,0.4508,0.784,0.0833,0.2752,
      ↪ 0.3585,0.25,0.4496,0.25,0.4167,0.2752,0.1415 & C2 C_{2}^{3} #5
      ↪ (ab^2c^21) & mC90 & $H4_{7}$ & Ca2H2O9S2 & Bassanite & W.
      ↪ Abriel and R. Nesper, Zeitschrift f{"u}r Kristallographie -
      ↪ Crystalline Materials 205, 99-113 (1993)
   1.00000000000000
   8.72906000000000   -3.46560000000000    0.00000000000000
```

```
   8.72906000000000    3.46560000000000    0.00000000000000
  -8.25868915382723    0.00000000000000    8.74384410373701
     Ca     H     O     S
      6     6    27     6
Direct
  -0.45530000000000    0.45530000000000    0.00000000000000   Ca   (2a)
   0.04470000000000   -0.04470000000000    0.50000000000000   Ca   (2b)
  -0.10570000000000    0.43910000000000    0.89430000000000   Ca   (4c)
  -0.43910000000000    0.10570000000000   -0.89430000000000   Ca   (4c)
   0.06090000000000    0.60570000000000    0.60570000000000   Ca   (4c)
  -0.60570000000000   -0.06090000000000   -0.60570000000000   Ca   (4c)
   0.32900000000000    0.67100000000000    0.57200000000000   H    (4c)
  -0.67100000000000   -0.32900000000000   -0.57200000000000   H    (4c)
   0.02600000000000    0.64000000000000    0.78300000000000   H    (4c)
  -0.64000000000000   -0.02600000000000   -0.78300000000000   H    (4c)
   0.81000000000000    0.85600000000000    0.71200000000000   H    (4c)
  -0.85600000000000   -0.81000000000000   -0.71200000000000   H    (4c)
  -0.59850000000000    0.59850000000000    0.50000000000000   O    (2b)
   0.29130000000000    0.74330000000000    0.88820000000000   O    (4c)
  -0.74330000000000   -0.29130000000000   -0.88820000000000   O    (4c)
  -0.20870000000000    0.24330000000000    0.38820000000000   O    (4c)
  -0.24330000000000    0.20870000000000   -0.38820000000000   O    (4c)
   0.40720000000000    1.29420000000000    0.55220000000000   O    (4c)
  -1.29420000000000   -0.40720000000000   -0.55220000000000   O    (4c)
  -0.09280000000000    0.79420000000000    0.05220000000000   O    (4c)
  -0.79420000000000    0.09280000000000   -0.05220000000000   O    (4c)
  -0.14660000000000    0.51460000000000    0.11150000000000   O    (4c)
  -0.51460000000000    0.14660000000000   -0.11150000000000   O    (4c)
   0.35340000000000    1.01460000000000    0.61150000000000   O    (4c)
  -1.01460000000000   -0.35340000000000   -0.61150000000000   O    (4c)
   0.53710000000000    0.77410000000000   -0.09820000000000   O    (4c)
  -0.77410000000000   -0.53710000000000    0.09820000000000   O    (4c)
   0.03710000000000    0.27410000000000    0.40180000000000   O    (4c)
  -0.27410000000000   -0.03710000000000   -0.40180000000000   O    (4c)
  -0.32130000000000    0.29910000000000    0.80650000000000   O    (4c)
  -0.29910000000000    0.32130000000000   -0.80650000000000   O    (4c)
   0.17870000000000    0.79910000000000    0.30650000000000   O    (4c)
  -0.79910000000000   -0.17870000000000   -0.30650000000000   O    (4c)
   0.25080000000000    0.39360000000000    0.75830000000000   O    (4c)
  -0.39360000000000   -0.25080000000000   -0.75830000000000   O    (4c)
   0.75080000000000    0.89360000000000    0.25830000000000   O    (4c)
  -0.89360000000000   -0.75080000000000   -0.25830000000000   O    (4c)
  -0.11750000000000    0.78410000000000    0.78400000000000   O    (4c)
  -0.78410000000000    0.11750000000000   -0.78400000000000   O    (4c)
  -0.19190000000000    0.35850000000000    0.35850000000000   S    (4c)
  -0.35850000000000    0.19190000000000   -0.35850000000000   S    (4c)
  -0.19960000000000    0.69960000000000    0.25000000000000   S    (4c)
  -0.69960000000000    0.19960000000000   -0.25000000000000   S    (4c)
   0.14150000000000    0.69190000000000    0.14150000000000   S    (4c)
  -0.69190000000000   -0.14150000000000   -0.14150000000000   S    (4c)
```

NbAs$_2$: A2B_mC12_5_2c_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'As2Nb'
_chemical_formula_sum 'As2 Nb'

loop_
_publ_author_name
 'S. Furuseth'
 'A. Kjekshus'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 18
_journal_year 1965
_journal_page_first 320
_journal_page_last 324
_publ_Section_title
;
 The Crystal Structures of NbAs$_{2}$ and NbSb$_{2}$
;

_aflow_title 'NbAs$_{2}$ Structure'
_aflow_proto 'A2B_mC12_5_2c_c'
_aflow_params 'a,b/a,c/a,\beta,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3
    ↪ },y_{3},z_{3}'
_aflow_params_values '9.357,0.36147269424,0.8327455381,119.76667,0.0948,
    ↪ 0.488,0.3928,0.1399,0.067,0.0257,0.3444,0.5,0.3044'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mC12'

_symmetry_space_group_name_H-M "C 1 2 1"
_symmetry_Int_Tables_number 5

_cell_length_a    9.35700
_cell_length_b    3.38230
_cell_length_c    7.79200
_cell_angle_alpha 90.00000
_cell_angle_beta  119.76667
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z
3 x+1/2,y+1/2,z
4 -x+1/2,y+1/2,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
As1 As 4 c 0.09480 0.48800 0.39280 1.00000
As2 As 4 c 0.13990 0.06700 0.02570 1.00000
Nb1 Nb 4 c 0.34440 0.50000 0.30440 1.00000
```

NbAs$_2$: A2B_mC12_5_2c_c - POSCAR

```
A2B_mC12_5_2c_c & a,b/a,c/a,beta,x1,y1,z1,x2,y2,z2,x3,y3,z3 --params=
    ↪ 9.357,0.36147269424,0.8327455381,119.76667,0.0948,0.488,0.3928,
    ↪ 0.1399,0.067,0.0257,0.3444,0.5,0.3044 & C2 C_{2}^{3} #5 (c^3) &
    ↪ mC12 & None & As2Nb & As2Nb & S. Furuseth and A. Kjekshus,
    ↪ Acta Cryst. 18, 320-324 (1965)
   1.00000000000000
   4.67850000000000   -1.69115000000000    0.00000000000000
   4.67850000000000    1.69115000000000    0.00000000000000
  -3.86848708676481    0.00000000000000    6.76387992645744
     As    Nb
      4     2
Direct
  -0.39320000000000    0.58280000000000    0.39280000000000   As   (4c)
  -0.58280000000000    0.39320000000000   -0.39280000000000   As   (4c)
   0.07290000000000    0.20690000000000    0.02570000000000   As   (4c)
  -0.20690000000000   -0.07290000000000   -0.02570000000000   As   (4c)
  -0.15560000000000    0.84440000000000    0.30440000000000   Nb   (4c)
  -0.84440000000000    0.15560000000000   -0.30440000000000   Nb   (4c)
```

$D0_{15}$ ($AlCl_3$) (*obsolete*): AB3_mC16_5_c_3c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'AlCl3'
_chemical_formula_sum 'Al Cl3'

loop_
_publ_author_name
 'J. A. A. Ketelaar'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 90
_journal_year 1935
_journal_page_first 237
_journal_page_last 255
_publ_Section_title
;
 Die Kristallstruktur der Aluminiumhalogenide II
;

# Found in Strukturbericht Band III 1933-1935, 1937

_aflow_title '$D0_{15}$ (AlCl$_{3}$) ({\em{obsolete}}) Structure'
_aflow_proto 'AB3_mC16_5_c_3c'
_aflow_params 'a,b/a,c/a,\beta,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3
    ↪ },y_{3},z_{3},x_{4},y_{4},z_{4}'
_aflow_params_values '5.91,1.73205076142,1.04286294416,108.64073,0.54333
    ↪ ,0.27833,0.045,0.24,0.27778,0.22,0.74,0.44444,0.22,0.74,0.11111
    ↪ ,0.22'
_aflow_Strukturbericht '$D0_{15}$'
_aflow_Pearson 'mC16'

_symmetry_space_group_name_H-M "C 1 2 1"
_symmetry_Int_Tables_number 5

_cell_length_a    5.91000
_cell_length_b    10.23642
_cell_length_c    6.16332
_cell_angle_alpha 90.00000
_cell_angle_beta  108.64073
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z
3 x+1/2,y+1/2,z
4 -x+1/2,y+1/2,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Al1 Al 4 c 0.54333 0.27833 0.04500 1.00000
Cl1 Cl 4 c 0.24000 0.27778 0.22000 1.00000
Cl2 Cl 4 c 0.74000 0.44444 0.22000 1.00000
Cl3 Cl 4 c 0.74000 0.11111 0.22000 1.00000
```

$D0_{15}$ ($AlCl_3$) (*obsolete*): AB3_mC16_5_c_3c - POSCAR

```
AB3_mC16_5_c_3c & a,b/a,c/a,beta,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,z4 --
    ↪ params=5.91,1.73205076142,1.04286294416,108.64073,0.54333,
    ↪ 0.27833,0.045,0.24,0.27778,0.22,0.74,0.44444,0.22,0.74,0.11111,
    ↪ 0.22 & C2 C_{2}^{3} #5 (c^4) & mC16 & $D0_{15}$ & AlCl3 & AlCl3
    ↪ & J. A. A. Ketelaar, Zeitschrift f{"u}r Kristallographie -
    ↪ Crystalline Materials 90, 237-255 (1935)
  1.00000000000000
  2.95500000000000  -5.11821000000000   0.00000000000000
  2.95500000000000   5.11821000000000   0.00000000000000
 -1.97000028435120   0.00000000000000   5.84000105325814
   Al   Cl
    2    6
Direct
  0.26500000000000   0.82166000000000   0.04500000000000  Al   (4c)
 -0.82166000000000  -0.26500000000000  -0.04500000000000  Al   (4c)
 -0.03778000000000   0.51778000000000   0.22000000000000  Cl   (4c)
 -0.51778000000000   0.03778000000000  -0.22000000000000  Cl   (4c)
  0.29556000000000   1.18444000000000   0.22000000000000  Cl   (4c)
 -1.18444000000000  -0.29556000000000  -0.22000000000000  Cl   (4c)
  0.62889000000000   0.85111000000000   0.22000000000000  Cl   (4c)
 -0.85111000000000  -0.62889000000000  -0.22000000000000  Cl   (4c)
```

$Rb_2CaCu_6(PO_4)_4O_2$: AB6C18D4E2_mC62_5_a_2b2c_9c_2c_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'CaCu6O18P4Rb2'
_chemical_formula_sum 'Ca Cu6 O18 P4 Rb2'

loop_
_publ_author_name
 'S. M. Aksenov'
 'E. Y. Borovikova'
 'V. S. Mironov'
 'N. A. Yamnova'
 'A. S. Volkov'
 'D. A. Ksenofontov'
 'O. A. Gurbanova'
 'O. V. Dimitrova'
 'D. V. Deyneko'
 'E. A. Zvereva'
 'O. V. Maximova'
 'S. V. Krivovichev'
 'P. C. Burns'
 'A. N. Vasiliev'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 75
_journal_year 2019
_journal_page_first 903
_journal_page_last 913
_publ_Section_title
;
 Rb$_{2}$CaCu$_{6}$(PO$_{4}$)$_{4}$O$_{2}$, a novel oxophosphate with a
    ↪ shchurovskyite-type topology: synthesis, structure, magnetic
    ↪ properties and crystal chemistry of rubidium copper phosphates
;

_aflow_title 'Rb$_{2}$CaCu$_{6}$(PO$_{4}$)$_{4}$O$_{2}$ Structure'
_aflow_proto 'AB6C18D4E2_mC62_5_a_2b2c_9c_2c_c'
_aflow_params 'a,b/a,c/a,\beta,y_{1},y_{2},y_{3},x_{4},y_{4},z_{4},x_{5
    ↪ },y_{5},z_{5},x_{6},y_{6},z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},
    ↪ z_{8},x_{9},y_{9},z_{9},x_{10},y_{10},z_{10},x_{11},y_{11},z_{
    ↪ 11},x_{12},y_{12},z_{12},x_{13},y_{13},z_{13},x_{14},y_{14},z_{
    ↪ 14},x_{15},y_{15},z_{15},x_{16},y_{16},z_{16},x_{17},y_{17},z_{
    ↪ 17}'
_aflow_params_values '16.8913,0.33393522109,0.494876060457,93.919,
    ↪ 0.67531,-0.07889,0.413,0.32143,-0.25821,0.47267,0.08396,0.18216
    ↪ ,0.22561,0.33645,-0.2696,0.2451,0.5762,0.1756,0.4677,0.07301,
    ↪ 0.1655,0.4508,0.47637,-0.1144,0.2054,0.4428,-0.5435,0.1571,
    ↪ 0.3822,-0.2162,-0.0248,0.71142,0.3504,0.4381,0.67462,-0.0515,
    ↪ 0.3216,0.61381,0.3236,0.1953,0.64484,0.2009,0.35178,0.4101,-
    ↪ 0.29307,0.14827,0.29009,0.20623,0.15752'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mC62'

_symmetry_space_group_name_H-M "C 1 2 1"
_symmetry_Int_Tables_number 5

_cell_length_a    16.89130
_cell_length_b    5.64060
_cell_length_c    8.35910
_cell_angle_alpha 90.00000
_cell_angle_beta  93.91900
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z
3 x+1/2,y+1/2,z
4 -x+1/2,y+1/2,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ca1 Ca 2 a 0.00000 0.67531 0.00000 1.00000
Cu1 Cu 2 b 0.00000 -0.07889 0.50000 1.00000
Cu2 Cu 2 b 0.00000 0.41300 0.50000 1.00000
Cu3 Cu 4 c 0.32143 -0.25821 0.47267 1.00000
Cu4 Cu 4 c 0.08396 0.18216 0.22561 1.00000
O1 O 4 c 0.33645 -0.26960 0.24510 1.00000
O2 O 4 c 0.57620 0.17560 0.46770 1.00000
O3 O 4 c 0.07301 0.16550 0.45080 1.00000
O4 O 4 c 0.47637 -0.11440 0.20540 1.00000
O5 O 4 c 0.44280 -0.54350 0.15710 1.00000
O6 O 4 c 0.38220 -0.21620 -0.02480 1.00000
O7 O 4 c 0.71142 0.35040 0.43810 1.00000
O8 O 4 c 0.67462 -0.05150 0.32160 1.00000
O9 O 4 c 0.61381 0.32360 0.19530 1.00000
P1 P 4 c 0.64484 0.20090 0.35178 1.00000
P2 P 4 c 0.41010 -0.29307 0.14827 1.00000
Rb1 Rb 4 c 0.29009 0.20623 0.15752 1.00000
```

$Rb_2CaCu_6(PO_4)_4O_2$: AB6C18D4E2_mC62_5_a_2b2c_9c_2c_c - POSCAR

```
AB6C18D4E2_mC62_5_a_2b2c_9c_2c_c & a,b/a,c/a,beta,y1,y2,y3,x4,y4,z4,x5,
    ↪ y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10,x11,y11,
    ↪ z11,x12,y12,z12,x13,y13,z13,x14,y14,z14,x15,y15,z15,x16,y16,z16
    ↪ ,x17,y17,z17 --params=16.8913,0.33393522109,0.494876060457,
    ↪ 93.919,0.67531,-0.07889,0.413,0.32143,-0.25821,0.47267,0.08396,
    ↪ 0.18216,0.22561,0.33645,-0.2696,0.2451,0.5762,0.1756,0.4677,
    ↪ 0.07301,0.1655,0.4508,0.47637,-0.1144,0.2054,0.4428,-0.5435,
    ↪ 0.1571,0.3822,-0.2162,-0.0248,0.71142,0.3504,0.4381,0.67462,-
    ↪ 0.0515,0.3216,0.61381,0.3236,0.1953,0.64484,0.2009,0.35178,
    ↪ 0.4101,-0.29307,0.14827,0.29009,0.20623,0.15752 & C2 C_{2}^{3}
    ↪ #5 (ab^2c^14) & mC62 & None & CaCu6O18P4Rb2 & CaCu6O18P4Rb2 &
    ↪ S. M. Aksenov et al., Acta Crystallogr. Sect. B Struct. Sci. 75
    ↪ , 903-913 (2019)
  1.00000000000000
  8.44565000000000  -2.82030000000000   0.00000000000000
  8.44565000000000   2.82030000000000   0.00000000000000
 -0.57131214828249   0.00000000000000   8.33955365947272
   Ca   Cu   O   P   Rb
    1    6  18    4    2
Direct
 -0.67531000000000   0.67531000000000   0.00000000000000  Ca   (2a)
  0.07889000000000  -0.07889000000000   0.50000000000000  Cu   (2b)
 -0.41300000000000   0.41300000000000   0.50000000000000  Cu   (2b)
  0.57964000000000   0.06322000000000   0.47267000000000  Cu   (4c)
 -0.06322000000000  -0.57964000000000  -0.47267000000000  Cu   (4c)
 -0.09820000000000   0.26612000000000   0.22561000000000  Cu   (4c)
 -0.26612000000000   0.09820000000000  -0.22561000000000  Cu   (4c)
  0.60605000000000   0.06685000000000   0.24510000000000  O    (4c)
 -0.06685000000000  -0.60605000000000  -0.24510000000000  O    (4c)
  0.40060000000000   0.75180000000000   0.46770000000000  O    (4c)
 -0.75180000000000  -0.40060000000000  -0.46770000000000  O    (4c)
 -0.09249000000000   0.23851000000000   0.45080000000000  O    (4c)
 -0.23851000000000   0.09249000000000  -0.45080000000000  O    (4c)
  0.59077000000000   0.36197000000000   0.20540000000000  O    (4c)
 -0.36197000000000  -0.59077000000000  -0.20540000000000  O    (4c)
  0.98630000000000  -0.10070000000000   0.15710000000000  O    (4c)
  0.10070000000000  -0.98630000000000  -0.15710000000000  O    (4c)
  0.59840000000000   0.16600000000000  -0.02480000000000  O    (4c)
 -0.16600000000000  -0.59840000000000   0.02480000000000  O    (4c)
  0.36102000000000   1.06182000000000   0.43810000000000  O    (4c)
 -1.06182000000000  -0.36102000000000  -0.43810000000000  O    (4c)
  0.72612000000000   0.62312000000000   0.32160000000000  O    (4c)
 -0.62312000000000  -0.72612000000000  -0.32160000000000  O    (4c)
  0.29021000000000   0.93741000000000   0.19530000000000  O    (4c)
 -0.93741000000000  -0.29021000000000  -0.19530000000000  O    (4c)
  0.44394000000000   0.84574000000000   0.35178000000000  P    (4c)
 -0.84574000000000  -0.44394000000000  -0.35178000000000  P    (4c)
  0.70317000000000   0.11703000000000   0.14827000000000  P    (4c)
 -0.11703000000000  -0.70317000000000  -0.14827000000000  P    (4c)
  0.08386000000000   0.49632000000000   0.15752000000000  Rb   (4c)
 -0.49632000000000  -0.08386000000000  -0.15752000000000  Rb   (4c)
```

$C2$ $(Ba,Ca)CO_3$: ABC3_mC10_5_b_a_ac - CIF

```
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_audit_creation_method FINDSYM

_chemical_name_mineral '(Ba,Ca)CO3'
_chemical_formula_sum 'Ba C O3'

loop_
_publ_author_name
 'D. Spahr'
 'L. Bayarjargal'
 'V. Vinograd'
 'R. Luchitskaia'
 'V. Milman'
 'B. Winkler'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 75
_journal_year 2019
_journal_page_first 291
_journal_page_last 300
_publ_Section_title
;
 A new BaCa(CO$_{3}$)$_{2}$ polymorph
;

_aflow_title '$C2$ (Ba,Ca)CO$_{3}$ Structure'
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_aflow_proto 'ABC3_mC10_5_b_a_ac'
_aflow_params 'a,b/a,c/a,\beta,y_{1},y_{2},y_{3},x_{4},y_{4},z_{4}'
_aflow_params_values '6.7064,0.763241083144,0.627266491709,109.284,
      ↪ 0.6259,0.8722,0.103,0.8957,0.5037,0.7367'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mC10'

_symmetry_space_group_name_H-M "C 1 2 1"
_symmetry_Int_Tables_number 5

_cell_length_a 6.70640
_cell_length_b 5.11860
_cell_length_c 4.20670
_cell_angle_alpha 90.00000
_cell_angle_beta 109.28400
_cell_angle_gamma 90.00000

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_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z
3 x+1/2,y+1/2,z
4 -x+1/2,y+1/2,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
C1 C 2 a 0.00000 0.62590 0.00000 1.00000
O1 O 2 a 0.00000 0.87220 0.00000 1.00000
Ba1 Ba 2 b 0.00000 0.10300 0.50000 1.00000
O2 O 4 c 0.89570 0.50370 0.73670 1.00000
```

$C2$ (Ba,Ca)CO_3: ABC3_mC10_5_b_a_ac - POSCAR

```
ABC3_mC10_5_b_a_ac & a,b/a,c/a,beta,y1,y2,y3,x4,y4,z4 --params=6.7064,
      ↪ 0.763241083144,0.627266491709,109.284,0.6259,0.8722,0.103,
      ↪ 0.8957,0.5037,0.7367 & C2 C_{2}^{3} #5 (a^2bc) & mC10 & None &
      ↪ (Ba,Ca)CO3 & (Ba,Ca)CO3 & D. Spahr et al., Acta Crystallogr.
      ↪ Sect. B Struct. Sci. 75, 291-300 (2019)
   1.00000000000000
   3.35320000000000  -2.55930000000000   0.00000000000000
   3.35320000000000   2.55930000000000   0.00000000000000
  -1.38926612821329   0.00000000000000   3.97067557413084
     Ba     C     O
      1     1     3
Direct
  -0.10300000000000   0.10300000000000   0.50000000000000  Ba   (2b)
  -0.62590000000000   0.62590000000000   0.00000000000000  C    (2a)
  -0.87220000000000   0.87220000000000   0.00000000000000  O    (2a)
   0.39200000000000   1.39940000000000   0.73670000000000  O    (4c)
  -1.39940000000000  -0.39200000000000  -0.73670000000000  O    (4c)
```

Ta_5Ti_{11} (BCC SQS-16): A5B11_mP16_6_2abc_2a3b3c - CIF

```
# AFLOW.org Repositories
# TaTi/A5B11_mP16_6_2abc_2a3b3c-001.AB params=6.5421326204,1.0,1.0,90.0,
      ↪ 0.5,-0.0,0.5,0.5,0.0,-0.0,0.0,0.5,0.0,0.5,0.0,-0.0,0.5,-0.0,0.5
      ↪ ,0.5,0.25,0.25,0.25,0.75,0.25,0.25,0.75,0.25,0.75,0.25,0.25,
      ↪ 0.75 SG=6 [ANRL doi: 10.1016/j.commatsci.2017.01.017 (part 1),
      ↪ doi: 10.1016/j.commatsci.2018.10.043 (part 2)]
data_TaTi
_pd_phase_name A5B11_mP16_6_2abc_2a3b3c-001.AB

_chemical_name_mineral 'Ta5Ti11'
_chemical_formula_sum 'Ta5 Ti11'

loop_
_publ_author_name
 'T. Chakraborty'
 'J. Rogal'
 'R. Drautz'
_journal_name_full_name
;
 Physical Review B
;
_journal_volume 94
_journal_year 2016
_journal_page_first 224104
_journal_page_last 224104
_publ_Section_title
;
 Unraveling the composition dependence of the martensitic transformation
      ↪ temperature: A first-principles study of Ti-Ta alloys
;

_aflow_title 'Ta$_{5}$Ti$_{11}$ (BCC SQS-16) Structure'
_aflow_proto 'A5B11_mP16_6_2abc_2a3b3c'
_aflow_params 'a,b/a,c/a,\beta,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3},x_{4
      ↪ },z_{4},x_{5},z_{5},x_{6},z_{6},x_{7},z_{7},x_{8},z_{8},x_{9},
      ↪ y_{9},z_{9},x_{10},y_{10},z_{10},x_{11},y_{11},z_{11},x_{12},y_
      ↪ {12},z_{12}'
_aflow_params_values '6.5421326204,1.0,1.0,90.0,0.0,0.5,0.5,0.5,0.0,0.0,
      ↪ 0.5,0.0,0.5,0.0,0.0,0.0,0.0,0.5,0.5,0.5,0.25,0.25,0.75,0.25,
      ↪ 0.25,0.25,0.75,0.25,0.25,0.75'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mP16'

_cell_length_a 6.5421326204
_cell_length_b 6.5421326204
_cell_length_c 6.5421326204
_cell_angle_alpha 90.0000000000
_cell_angle_beta 90.0000000000
_cell_angle_gamma 90.0000000000
_symmetry_space_group_name_H-M 'Pm'
_symmetry_Int_Tables_Number 6
loop_
_symmetry_equiv_pos_site_id
_symmetry_equiv_pos_as_xyz
  1 x,y,z
  2 x,-y,z
loop_
 _atom_site_label
 _atom_site_occupancy
 _atom_site_fract_x
 _atom_site_fract_y
 _atom_site_fract_z
 _atom_site_thermal_displace_type
 _atom_site_B_iso_or_equiv
 _atom_site_type_symbol
 _atom_site_symmetry_multiplicity
 _atom_site_Wyckoff_label
Ta1 1.0000000000 0.0000000000 0.0000000000 0.5000000000 Biso 1.0 Ta 1 a
Ta2 1.0000000000 0.5000000000 0.0000000000 0.5000000000 Biso 1.0 Ta 1 a
Ti1 1.0000000000 0.0000000000 0.0000000000 0.0000000000 Biso 1.0 Ti 1 a
Ti2 1.0000000000 0.5000000000 0.0000000000 0.0000000000 Biso 1.0 Ti 1 a
Ta3 1.0000000000 0.5000000000 0.5000000000 0.0000000000 Biso 1.0 Ta 1 b
Ti3 1.0000000000 0.0000000000 0.5000000000 0.0000000000 Biso 1.0 Ti 1 b
Ti4 1.0000000000 0.0000000000 0.5000000000 0.5000000000 Biso 1.0 Ti 1 b
Ti5 1.0000000000 0.5000000000 0.5000000000 0.5000000000 Biso 1.0 Ti 1 b
Ta4 1.0000000000 0.2500000000 0.2500000000 0.7500000000 Biso 1.0 Ta 2 c
Ti6 1.0000000000 0.2500000000 0.2500000000 0.2500000000 Biso 1.0 Ti 2 c
Ti7 1.0000000000 0.7500000000 0.2500000000 0.2500000000 Biso 1.0 Ti 2 c
Ti8 1.0000000000 0.7500000000 0.2500000000 0.7500000000 Biso 1.0 Ti 2 c
```

Ta_5Ti_{11} (BCC SQS-16): A5B11_mP16_6_2abc_2a3b3c - POSCAR

```
A5B11_mP16_6_2abc_2a3b3c & a,b/a,c/a,beta,x1,z1,x2,z2,x3,z3,x4,z4,x5,z5,
      ↪ x6,z6,x7,z7,x8,z8,x9,y9,z9,x10,y10,z10,x11,y11,z11,x12,y12,z12
      ↪ --params=6.5421326204,1.0,1.0,90.0,0.0,0.5,0.5,0.5,0.0,0.0,0.5,
      ↪ 0.0,0.5,0.0,0.0,0.0,0.0,0.5,0.5,0.5,0.25,0.25,0.75,0.25,0.25,
      ↪ 0.25,0.75,0.25,0.25,0.75,0.25,0.75 & Pm C_{s}^{1} #6 (a^4b^4c^4
      ↪ ) & mP16 & None & Ta5Ti11 & Ta5Ti11 & T. Chakraborty and J.
      ↪ Rogal and R. Drautz, Phys. Rev. B 94, 224104(2016)
   1.00000000000000
   6.54213262040000   0.00000000000000   0.00000000000000
   0.00000000000000   6.54213262040000   0.00000000000000
   0.00000000000000   0.00000000000000   6.54213262040000
    Ta    Ti
     5    11
Direct
   0.00000000000000   0.00000000000000   0.50000000000000  Ta   (1a)
   0.50000000000000   0.00000000000000   0.50000000000000  Ta   (1a)
   0.50000000000000   0.50000000000000   0.00000000000000  Ta   (1b)
   0.25000000000000   0.25000000000000   0.75000000000000  Ta   (2c)
   0.25000000000000  -0.25000000000000   0.75000000000000  Ta   (2c)
   0.00000000000000   0.00000000000000   0.00000000000000  Ti   (1a)
   0.50000000000000   0.00000000000000   0.00000000000000  Ti   (1a)
   0.00000000000000   0.50000000000000   0.00000000000000  Ti   (1b)
   0.00000000000000   0.50000000000000   0.50000000000000  Ti   (1b)
   0.50000000000000   0.50000000000000   0.50000000000000  Ti   (1b)
   0.25000000000000   0.25000000000000   0.25000000000000  Ti   (2c)
   0.25000000000000  -0.25000000000000   0.25000000000000  Ti   (2c)
   0.75000000000000   0.25000000000000   0.25000000000000  Ti   (2c)
   0.75000000000000  -0.25000000000000   0.25000000000000  Ti   (2c)
   0.75000000000000   0.25000000000000   0.75000000000000  Ti   (2c)
   0.75000000000000  -0.25000000000000   0.75000000000000  Ti   (2c)
```

$Na_2Ca_6Si_4O_{15}$: A6B2C15D4_mP54_7_6a_2a_15a_4a - CIF

```
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_chemical_formula_sum 'Ca6 Na2 O15 Si4'

loop_
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 'V. Kahlenberg'
 'M. Maier'
_journal_name_full_name
;
 Mineralogy and Petrology
;
_journal_volume 110
_journal_year 2016
_journal_page_first 905
_journal_page_last 915
_publ_Section_title
;
 On the existence of a high-temperature polymorph of Na$_{2}$Ca$_{6}
      ↪ $Si$_{4}$O$_{15}$-implications for the phase equilibria in
      ↪ the system Na$_{2}$O-CaO-SiO$_{2}$
;

_aflow_title 'Na$_{2}$Ca$_{6}$Si$_{4}$O$_{15}$ Structure'
_aflow_proto 'A6B2C15D4_mP54_7_6a_2a_15a_4a'
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      ↪ z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{
      ↪ 10},y_{10},z_{10},x_{11},y_{11},z_{11},x_{12},y_{12},z_{12},x_{
      ↪ 13},y_{13},z_{13},x_{14},y_{14},z_{14},x_{15},y_{15},z_{15},x_{
      ↪ 16},y_{16},z_{16},x_{17},y_{17},z_{17},x_{18},y_{18},z_{18},x_{
```

```
      ↪ 19},y_{19},z_{19},x_{20},y_{20},z_{20},x_{21},y_{21},z_{21},x_{
      ↪ 22},y_{22},z_{22},x_{23},y_{23},z_{23},x_{24},y_{24},z_{24},x_{
      ↪ 25},y_{25},z_{25},x_{26},y_{26},z_{26},x_{27},y_{27},z_{27}'
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      ↪ 0.2658,0.07662,0.6782,0.50471,0.1731,0.3148,0.0044,0.3329,-
      ↪ 0.0022,0.2541,0.27358,-0.0033,0.7531,0.31251,0.6554,0.2291,-
      ↪ 0.09514,0.3179,0.718,0.0741,0.6731,0.2144,0.4123,0.8207,0.4832,
      ↪ 0.8963,0.2666,0.3166,0.3293,0.2631,0.6918,0.3234,0.0643,-0.0059
      ↪ ,0.1862,0.7163,0.8148,0.1874,-0.0723,0.5119,0.1535,0.497,0.4776
      ↪ ,-0.022,0.5282,-0.0241,0.5029,0.1732,0.0159,-0.0177,0.0011,
      ↪ 0.7477,-0.0003,0.4988,0.0147,0.234,0.456,0.5016,0.2321,0.7307,
      ↪ 0.1924,0.191,0.8718,0.0697,-0.0515,0.1138,0.4252,0.0116,0.3664,
      ↪ 0.5049,0.3381,0.6271,-0.0087,0.6556,0.0263,-0.0323,0.0347,-
      ↪ 0.0333,0.5304,0.0206'
_aflow_Strukturbericht  'None'
_aflow_Pearson  'mP54'

_symmetry_space_group_name_H-M  "P 1 c 1"
_symmetry_Int_Tables_number  7

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_cell_length_b      7.31710
_cell_length_c      10.97230
_cell_angle_alpha  90.00000
_cell_angle_beta   107.72000
_cell_angle_gamma  90.00000

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_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
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Ca2  Ca  2  a  0.67820  0.50471  0.17310  1.00000
Ca3  Ca  2  a  0.31480  0.00440  0.33290  1.00000
Ca4  Ca  2  a -0.00220  0.25410  0.27358  1.00000
Ca5  Ca  2  a -0.00330  0.75310  0.31251  1.00000
Ca6  Ca  2  a  0.65540  0.22910 -0.09514  1.00000
Na1  Na  2  a  0.31790  0.71800  0.07410  1.00000
Na2  Na  2  a  0.67310  0.21440  0.41230  1.00000
O1  O  2  a  0.82070  0.48320  0.89630  1.00000
O2  O  2  a  0.26660  0.31660  0.32930  1.00000
O3  O  2  a  0.26310  0.69180  0.32340  1.00000
O4  O  2  a  0.06430 -0.00590  0.18620  1.00000
O5  O  2  a  0.71630  0.81480  0.18740  1.00000
O6  O  2  a -0.07230  0.51190  0.15350  1.00000
O7  O  2  a  0.49700  0.47760 -0.02200  1.00000
O8  O  2  a  0.52820 -0.02410  0.50290  1.00000
O9  O  2  a  0.17320  0.01590 -0.01770  1.00000
O10  O  2  a  0.00110  0.74770 -0.00030  1.00000
O11  O  2  a  0.49880  0.01470  0.23400  1.00000
O12  O  2  a  0.45600  0.50160  0.23210  1.00000
O13  O  2  a  0.73070  0.19240  0.19100  1.00000
O14  O  2  a  0.87180  0.06970 -0.05150  1.00000
O15  O  2  a  0.11380  0.42520  0.01160  1.00000
Si1  Si  2  a  0.36640  0.50490  0.33810  1.00000
Si2  Si  2  a  0.62710 -0.00870  0.65560  1.00000
Si3  Si  2  a  0.02630 -0.03230  0.03470  1.00000
Si4  Si  2  a -0.03330  0.53040  0.02060  1.00000
```

$Na_2Ca_6Si_4O_{15}$: A6B2C15D4_mP54_7_6a_2a_15a_4a - POSCAR

```
A6B2C15D4_mP54_7_6a_2a_15a_4a & a,b/a,c/a,beta,x1,y1,z1,x2,y2,z2,x3,y3,
      ↪ z3,x4,y4,z4,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,
      ↪ y10,z10,x11,y11,z11,x12,y12,z12,x13,y13,z13,x14,y14,z14,x15,y15
      ↪ ,z15,x16,y16,z16,x17,y17,z17,x18,y18,z18,x19,y19,z19,x20,y20,
      ↪ z20,x21,y21,z21,x22,y22,z22,x23,y23,z23,x24,y24,z24,x25,y25,z25
      ↪ ,x26,y26,z26,x27,y27,z27 --params=9.0112,0.812000621449,
      ↪ 1.21762917259,107.72,0.3409,0.2658,0.07662,0.6782,0.50471,
      ↪ 0.1731,0.3148,0.0044,0.3329,-0.0022,0.2541,0.27358,-0.0033,
      ↪ 0.7531,0.31251,0.6554,0.2291,-0.09514,0.3179,0.718,0.0741,
      ↪ 0.6731,0.2144,0.4123,0.8207,0.4832,0.8963,0.2666,0.3166,0.3293,
      ↪ 0.2631,0.6918,0.3234,0.0643,-0.0059,0.1862,0.7163,0.8148,0.1874
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      ↪ Ca6Na2O15Si4 & V. Kahlenberg and M. Maier, Mineral. Petrol. 110
      ↪ , 905-915 (2016)
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Low-Temperature Mo_8O_{23}: A8B23_mP124_7_16a_46a - CIF

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 Journal of Solid State Chemistry
;
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_journal_year 1987
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;
 Structure Determination of low-dimensional conductor Mo$_{8}$O$_{23}$
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Low-Temperature Mo_8O_{23}: A8B23_mP124_7_16a_46a - POSCAR

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  0.81760000000000   -0.20600000000000    0.75430000000000  Mo  (2a)
  0.18620000000000    0.70700000000000    0.24480000000000  Mo  (2a)
  0.18620000000000   -0.70700000000000    0.74480000000000  Mo  (2a)
  0.81560000000000    0.70490000000000    0.25630000000000  Mo  (2a)
  0.81560000000000   -0.70490000000000    0.75630000000000  Mo  (2a)
  0.31540000000000    0.28990000000000    0.07960000000000  Mo  (2a)
  0.31540000000000   -0.28990000000000    0.57960000000000  Mo  (2a)
  0.68510000000000    0.29540000000000    0.41870000000000  Mo  (2a)
  0.68510000000000   -0.29540000000000    0.91870000000000  Mo  (2a)
  0.31560000000000    0.79360000000000    0.07740000000000  Mo  (2a)
  0.31560000000000   -0.79360000000000    0.57740000000000  Mo  (2a)
  0.68450000000000    0.79970000000000    0.42100000000000  Mo  (2a)
  0.68450000000000   -0.79970000000000    0.92100000000000  Mo  (2a)
  0.44720000000000    0.20430000000000    0.40150000000000  Mo  (2a)
  0.44720000000000   -0.20430000000000    0.90150000000000  Mo  (2a)
  0.55370000000000    0.20600000000000    0.09700000000000  Mo  (2a)
  0.55370000000000   -0.20600000000000    0.59700000000000  Mo  (2a)
  0.44700000000000    0.70360000000000    0.40120000000000  Mo  (2a)
  0.44700000000000   -0.70360000000000    0.90120000000000  Mo  (2a)
  0.55460000000000    0.70490000000000    0.09810000000000  Mo  (2a)
  0.55460000000000   -0.70490000000000    0.59810000000000  Mo  (2a)
  0.07790000000000   -0.00110000000000    0.40500000000000   O  (2a)
  0.07790000000000    0.00110000000000    0.90500000000000   O  (2a)
 -0.07280000000000    0.00120000000000    0.08910000000000   O  (2a)
 -0.07280000000000   -0.00120000000000    0.58910000000000   O  (2a)
  0.06070000000000    0.49900000000000    0.42960000000000   O  (2a)
  0.06070000000000   -0.49900000000000    0.92960000000000   O  (2a)
 -0.04860000000000    0.50480000000000    0.07830000000000   O  (2a)
 -0.04860000000000   -0.50480000000000    0.57830000000000   O  (2a)
  0.19290000000000    0.49120000000000    0.23390000000000   O  (2a)
  0.19290000000000   -0.49120000000000    0.73390000000000   O  (2a)
  0.79760000000000    0.50400000000000    0.26130000000000   O  (2a)
  0.79760000000000   -0.50400000000000    0.76130000000000   O  (2a)
  0.19000000000000   -0.00290000000000    0.25740000000000   O  (2a)
  0.19000000000000    0.00290000000000    0.75740000000000   O  (2a)
  0.82160000000000   -0.00400000000000    0.24780000000000   O  (2a)
  0.82160000000000    0.00400000000000    0.74780000000000   O  (2a)
  0.32140000000000    0.00820000000000    0.06700000000000   O  (2a)
  0.32140000000000   -0.00820000000000    0.56700000000000   O  (2a)
  0.68130000000000    0.00600000000000    0.43030000000000   O  (2a)
  0.68130000000000   -0.00600000000000    0.93030000000000   O  (2a)
  0.31520000000000    0.50000000000000    0.08820000000000   O  (2a)
  0.31520000000000   -0.50000000000000    0.58820000000000   O  (2a)
  0.67750000000000    0.50670000000000    0.41520000000000   O  (2a)
  0.67750000000000   -0.50670000000000    0.91520000000000   O  (2a)
  0.44360000000000    0.49780000000000    0.40740000000000   O  (2a)
  0.44360000000000   -0.49780000000000    0.90740000000000   O  (2a)
```

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   0.54600000000000   0.48850000000000   0.09950000000000   O   (2a)
   0.54600000000000  -0.48850000000000   0.59950000000000   O   (2a)
   0.45210000000000  -0.01090000000000   0.41520000000000   O   (2a)
   0.45210000000000   0.01090000000000   0.91520000000000   O   (2a)
   0.55800000000000  -0.00260000000000   0.09350000000000   O   (2a)
   0.55800000000000   0.00260000000000   0.59350000000000   O   (2a)
  -0.00020000000000   0.25050000000000  -0.00150000000000   O   (2a)
  -0.00020000000000  -0.25050000000000   0.49850000000000   O   (2a)
   0.00240000000000   0.79000000000000  -0.00150000000000   O   (2a)
   0.00240000000000  -0.79000000000000   0.49850000000000   O   (2a)
   0.06060000000000   0.24670000000000   0.16010000000000   O   (2a)
   0.06060000000000  -0.24670000000000   0.66010000000000   O   (2a)
  -0.06720000000000   0.24290000000000   0.32820000000000   O   (2a)
  -0.06720000000000  -0.24290000000000   0.82820000000000   O   (2a)
   0.06010000000000   0.75420000000000   0.15890000000000   O   (2a)
   0.06010000000000  -0.75420000000000   0.65890000000000   O   (2a)
  -0.06730000000000   0.74280000000000   0.33170000000000   O   (2a)
  -0.06730000000000  -0.74280000000000   0.83170000000000   O   (2a)
   0.12650000000000   0.27070000000000   0.32740000000000   O   (2a)
   0.12650000000000  -0.27070000000000   0.82740000000000   O   (2a)
   0.86680000000000   0.27760000000000   0.16590000000000   O   (2a)
   0.86680000000000  -0.27760000000000   0.66590000000000   O   (2a)
   0.13410000000000   0.71860000000000   0.32880000000000   O   (2a)
   0.13410000000000  -0.71860000000000   0.82880000000000   O   (2a)
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   0.87560000000000  -0.71300000000000   0.66920000000000   O   (2a)
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   0.20110000000000  -0.24640000000000   0.99290000000000   O   (2a)
   0.80065000000000   0.25820000000000  -0.00100000000000   O   (2a)
   0.80065000000000  -0.25820000000000   0.49900000000000   O   (2a)
   0.19430000000000   0.73510000000000   0.49210000000000   O   (2a)
   0.19430000000000  -0.73510000000000   0.99210000000000   O   (2a)
   0.79940000000000   0.74760000000000   0.00180000000000   O   (2a)
   0.79940000000000  -0.74760000000000   0.50180000000000   O   (2a)
   0.25930000000000   0.24370000000000   0.15780000000000   O   (2a)
   0.25930000000000  -0.24370000000000   0.65780000000000   O   (2a)
   0.74170000000000   0.22900000000000   0.33360000000000   O   (2a)
   0.74170000000000  -0.22900000000000   0.83360000000000   O   (2a)
   0.25790000000000   0.77180000000000   0.16050000000000   O   (2a)
   0.25790000000000  -0.77180000000000   0.66050000000000   O   (2a)
   0.73510000000000   0.78230000000000   0.33000000000000   O   (2a)
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   0.32260000000000   0.25450000000000   0.32430000000000   O   (2a)
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   0.67080000000000  -0.23160000000000   0.67060000000000   O   (2a)
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   0.40850000000000  -0.25800000000000   0.50330000000000   O   (2a)
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   0.40240000000000  -0.74510000000000   0.49870000000000   O   (2a)
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   0.45420000000000   0.24770000000000   0.14920000000000   O   (2a)
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```

Calaverite ($AuTe_2$): AB2_mP12_7_2a_4a - CIF

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_chemical_formula_sum 'Au Te2'

loop_
_publ_author_name
 'F. Pertlik'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 169
_journal_year 1984
_journal_page_first 227
_journal_page_last 236
_publ_Section_title
;
 Kristallchemie nat\"{u}licher Telluride III: Die Kristallstruktur des
      ↪ Minerals Calaverit, AuTe$_2$
;

_aflow_title 'Calaverite (AuTe$_{2}$) Structure'
_aflow_proto 'AB2_mP12_7_2a_4a'
_aflow_params 'a,b/a,c/a,\beta,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3
      ↪ },y_{3},z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},
      ↪ z_{6}'
_aflow_params_values '8.76,0.503424657534,1.15867579909,125.2,0.7483,
      ↪ 0.4938,0.1243,0.246,-0.0059,0.8731,-0.0649,-0.0138,0.5728,
      ↪ 0.5585,0.052,0.6747,0.0575,0.4464,0.4209,0.4363,0.5585,0.3235'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mP12'

_symmetry_space_group_name_H-M "P 1 c 1"
_symmetry_Int_Tables_number 7

_cell_length_a      8.76000
_cell_length_b      4.41000
_cell_length_c      10.15000
_cell_angle_alpha  90.00000
_cell_angle_beta   125.20000
_cell_angle_gamma  90.00000

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_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Au1 Au 2 a 0.74830 0.49380 0.12430 1.00000
Au2 Au 2 a 0.24600 -0.00590 0.87310 1.00000
Te1 Te 2 a -0.06490 -0.01380 0.57280 1.00000
Te2 Te 2 a 0.55850 0.05200 0.67470 1.00000
Te3 Te 2 a 0.05750 0.44640 0.42090 1.00000
Te4 Te 2 a 0.43630 0.55850 0.32350 1.00000
```

Calaverite ($AuTe_2$): AB2_mP12_7_2a_4a - POSCAR

```
AB2_mP12_7_2a_4a & a,b/a,c/a,beta,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,z4,x5
      ↪ ,y5,z5,x6,y6,z6 --params=8.76,0.503424657534,1.15867579909,
      ↪ 125.2,0.7483,0.4938,0.1243,0.246,-0.0059,0.8731,-0.0649,-0.0138
      ↪ ,0.5728,0.5585,0.052,0.6747,0.0575,0.4464,0.4209,0.4363,0.5585,
      ↪ 0.3235 & Pc C_{s}^{2} #7 (a^6) & mP12 & None & AuTe2 & AuTe2 &
      ↪ F. Pertlik, Zeitschrift f{"u}r Kristallographie - Crystalline
      ↪ Materials 169, 227-236 (1984)
   1.00000000000000
     8.76000000000000   0.00000000000000   0.00000000000000
     0.00000000000000   4.41000000000000   0.00000000000000
    -5.85078800912340   0.00000000000000   8.29402071810155
      Au    Te
       4     8
Direct
     0.74830000000000   0.49380000000000   0.12430000000000   Au   (2a)
     0.74830000000000  -0.49380000000000   0.62430000000000   Au   (2a)
     0.24600000000000  -0.00590000000000   0.87310000000000   Au   (2a)
     0.24600000000000   0.00590000000000   1.37310000000000   Au   (2a)
    -0.06490000000000  -0.01380000000000   0.57280000000000   Te   (2a)
    -0.06490000000000   0.01380000000000   1.07280000000000   Te   (2a)
     0.55850000000000   0.05200000000000   0.67470000000000   Te   (2a)
     0.55850000000000  -0.05200000000000   1.17470000000000   Te   (2a)
     0.05750000000000   0.44640000000000   0.42090000000000   Te   (2a)
     0.05750000000000  -0.44640000000000   0.92090000000000   Te   (2a)
     0.43630000000000   0.55850000000000   0.32350000000000   Te   (2a)
     0.43630000000000  -0.55850000000000   0.82350000000000   Te   (2a)
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Monoclinic Co_4Al_{13}: A13B4_mC102_8_17a11b_8a2b - CIF

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loop_
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 'R. C. Hudd'
 'W. H. Taylor'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 15
_journal_year 1962
_journal_page_first 441
_journal_page_last 442
_publ_Section_title
;
 The Structure of Co$_{4}$Al$_{13}$
;

# Found in Structure investigation of the (100) surface of the
      ↪ orthorhombic Al$_{13}$Co$_{4}$ crystal, 2009

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_aflow_proto 'A13B4_mC102_8_17a11b_8a2b'
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      ↪ z_{9},x_{10},z_{10},x_{11},z_{11},x_{12},z_{12},x_{13},z_{13},
      ↪ x_{14},z_{14},x_{15},z_{15},x_{16},z_{16},x_{17},z_{17},x_{18},
      ↪ z_{18},x_{19},z_{19},x_{20},z_{20},x_{21},z_{21},x_{22},z_{22},
      ↪ x_{23},z_{23},x_{24},z_{24},x_{25},z_{25},x_{26},y_{26},z_{26},
      ↪ x_{27},y_{27},z_{27},x_{28},y_{28},z_{28},x_{29},y_{29},z_{29},
      ↪ x_{30},y_{30},z_{30},x_{31},y_{31},z_{31},x_{32},y_{32},z_{32},
      ↪ x_{33},y_{33},z_{33},x_{34},y_{34},z_{34},x_{35},y_{35},z_{35},
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_aflow_params_values '15.183,0.534940393862,0.812751103208,107.9,0.055,
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      ↪ 0.169,0.502,0.499,0.702,0.23,-0.095,0.212,-0.055,0.825,0.67,
      ↪ 0.72,0.232,0.532,0.077,0.576,0.245,-0.034,0.477,0.831,0.297,
```

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      ↪ 0.769,0.095,0.788,0.086,0.382,0.604,0.373,-0.091,0.012,0.597,
      ↪ 0.018,-0.087,0.617,0.395,0.626,0.09,-0.013,0.402,-0.018,0.185,
      ↪ 0.217,0.111,0.364,0.21,0.111,0.176,0.219,0.334,0.491,0.224,
      ↪ 0.331,0.367,0.21,0.477,0.498,0.248,0.0,0.314,0.282,0.888,0.124,
      ↪ 0.271,0.888,0.322,0.28,0.666,0.008,0.279,0.669,0.131,0.289,
      ↪ 0.532,0.318,0.292,0.277,0.182,0.229,0.722'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mC102'

_symmetry_space_group_name_H-M "C 1 m 1"
_symmetry_Int_Tables_number 8

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_cell_length_b 8.12200
_cell_length_c 12.34000
_cell_angle_alpha 90.00000
_cell_angle_beta 107.90000
_cell_angle_gamma 90.00000

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_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,z
3 x+1/2,y+1/2,z
4 x+1/2,-y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
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Al2 Al 2 a 0.32900 0.00000 0.28000 1.00000
Al3 Al 2 a 0.76700 0.00000 0.46900 1.00000
Al4 Al 2 a -0.07700 0.00000 0.42400 0.30000
Al5 Al 2 a 0.75400 0.00000 0.03100 0.70000
Al6 Al 2 a 0.52200 0.00000 0.16900 1.00000
Al7 Al 2 a 0.50200 0.00000 0.49900 1.00000
Al8 Al 2 a 0.70200 0.00000 0.23000 1.00000
Al9 Al 2 a -0.09500 0.00000 0.21200 0.70000
Al10 Al 2 a -0.05500 0.00000 0.82500 1.00000
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Al12 Al 2 a 0.23200 0.00000 0.53200 1.00000
Al13 Al 2 a 0.07700 0.00000 0.57600 0.30000
Al14 Al 2 a 0.24500 0.00000 -0.03400 0.70000
Al15 Al 2 a 0.47700 0.00000 0.83100 1.00000
Al16 Al 2 a 0.29700 0.00000 0.76900 1.00000
Al17 Al 2 a 0.09500 0.00000 0.78800 0.30000
Co1 Co 2 a 0.08600 0.00000 0.38200 1.00000
Co2 Co 2 a 0.60400 0.00000 0.37300 1.00000
Co3 Co 2 a -0.09100 0.00000 0.01200 1.00000
Co4 Co 2 a 0.59700 0.00000 0.01800 1.00000
Co5 Co 2 a -0.08700 0.00000 0.61700 1.00000
Co6 Co 2 a 0.39500 0.00000 0.62600 1.00000
Co7 Co 2 a 0.09000 0.00000 -0.01300 1.00000
Co8 Co 2 a 0.40200 0.00000 -0.01800 1.00000
Al18 Al 4 b 0.18500 0.21700 0.11100 1.00000
Al19 Al 4 b 0.36400 0.21000 0.11100 1.00000
Al20 Al 4 b 0.17600 0.21900 0.33400 1.00000
Al21 Al 4 b 0.49100 0.22400 0.33100 1.00000
Al22 Al 4 b 0.36700 0.21000 0.47700 1.00000
Al23 Al 4 b 0.49800 0.24800 0.00000 1.00000
Al24 Al 4 b 0.31400 0.28200 0.88800 1.00000
Al25 Al 4 b 0.12400 0.27100 0.88800 1.00000
Al26 Al 4 b 0.32200 0.28000 0.66600 1.00000
Al27 Al 4 b 0.00800 0.27900 0.66900 1.00000
Al28 Al 4 b 0.13100 0.28900 0.53200 1.00000
Co9 Co 4 b 0.31800 0.29200 0.27700 1.00000
Co10 Co 4 b 0.18200 0.22900 0.72200 1.00000
```

Monoclinic Co_4Al_{13}: A13B4_mC102_8_17a11b_8a2b - POSCAR

```
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      ↪ ,x6,z6,x7,z7,x8,z8,x9,z9,x10,z10,x11,z11,x12,z12,x13,z13,x14,
      ↪ z14,x15,z15,x16,z16,x17,z17,x18,z18,x19,z19,x20,z20,x21,z21,x22
      ↪ ,z22,x23,z23,x24,z24,x25,z25,x26,y26,z26,x27,y27,z27,x28,y28,
      ↪ z28,x29,y29,z29,x30,y30,z30,x31,y31,z31,x32,y32,z32,x33,y33,z33
      ↪ ,x34,y34,z34,x35,y35,z35,x36,y36,z36,x37,y37,z37,x38,y38,z38 --
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      ↪ 0.232,0.532,0.077,0.576,0.245,-0.034,0.477,0.831,0.297,0.769,
      ↪ 0.095,0.788,0.086,0.382,0.604,0.373,-0.091,0.012,0.597,0.018,-
      ↪ 0.087,0.617,0.395,0.626,0.09,-0.013,0.402,-0.018,0.185,0.217,
      ↪ 0.111,0.364,0.21,0.111,0.176,0.219,0.334,0.491,0.224,0.331,
      ↪ 0.367,0.21,0.477,0.498,0.248,0.0,0.314,0.282,0.888,0.124,0.271,
      ↪ 0.888,0.322,0.28,0.666,0.008,0.279,0.669,0.131,0.289,0.532,
      ↪ 0.318,0.292,0.277,0.182,0.229,0.722 & Cm C_{s}^{3} #8 (a^25b^13
      ↪ ) & mC102 & None & Al13Co4 & Al13Co4 & R. C. Hudd and W. H.
      ↪ Taylor, Acta Cryst. 15, 441-442 (1962)
   1.00000000000000
   7.59150000000000  -4.06100000000000   0.00000000000000
   7.59150000000000   4.06100000000000   0.00000000000000
  -3.79278066364962   0.00000000000000  11.74267494387230
     Al    Co
     39    12
Direct
   0.05500000000000   0.05500000000000   0.17400000000000   Al   (2a)
   0.32900000000000   0.32900000000000   0.28000000000000   Al   (2a)
   0.76700000000000   0.76700000000000   0.46900000000000   Al   (2a)
  -0.07700000000000  -0.07700000000000   0.42400000000000   Al   (2a)
   0.75400000000000   0.75400000000000   0.03100000000000   Al   (2a)
   0.52200000000000   0.52200000000000   0.16900000000000   Al   (2a)
   0.50200000000000   0.50200000000000   0.49900000000000   Al   (2a)
   0.70200000000000   0.70200000000000   0.23000000000000   Al   (2a)
  -0.09500000000000  -0.09500000000000   0.21200000000000   Al   (2a)
  -0.05500000000000  -0.05500000000000   0.82500000000000   Al   (2a)
   0.67000000000000   0.67000000000000   0.72000000000000   Al   (2a)
   0.23200000000000   0.23200000000000   0.53200000000000   Al   (2a)
   0.07700000000000   0.07700000000000   0.57600000000000   Al   (2a)
   0.24500000000000   0.24500000000000  -0.03400000000000   Al   (2a)
   0.47700000000000   0.47700000000000   0.83100000000000   Al   (2a)
   0.29700000000000   0.29700000000000   0.76900000000000   Al   (2a)
   0.09500000000000   0.09500000000000   0.78800000000000   Al   (2a)
  -0.03200000000000   0.04200000000000   0.11100000000000   Al   (4b)
   0.40200000000000  -0.03200000000000   0.11100000000000   Al   (4b)
   0.15400000000000   0.57400000000000   0.11100000000000   Al   (4b)
   0.57400000000000   0.15400000000000   0.11100000000000   Al   (4b)
  -0.04300000000000   0.39500000000000   0.33400000000000   Al   (4b)
   0.39500000000000  -0.04300000000000   0.33400000000000   Al   (4b)
   0.26700000000000   0.71500000000000   0.33100000000000   Al   (4b)
   0.71500000000000   0.26700000000000   0.33100000000000   Al   (4b)
   0.15700000000000   0.57700000000000   0.47700000000000   Al   (4b)
   0.57700000000000   0.15700000000000   0.47700000000000   Al   (4b)
   0.25000000000000   0.74600000000000   0.00000000000000   Al   (4b)
   0.74600000000000   0.25000000000000   0.00000000000000   Al   (4b)
   0.03200000000000   0.59600000000000   0.88800000000000   Al   (4b)
   0.59600000000000   0.03200000000000   0.88800000000000   Al   (4b)
  -0.14700000000000   0.39500000000000   0.88800000000000   Al   (4b)
   0.39500000000000  -0.14700000000000   0.88800000000000   Al   (4b)
   0.04200000000000   0.60200000000000   0.66600000000000   Al   (4b)
   0.60200000000000   0.04200000000000   0.66600000000000   Al   (4b)
  -0.27100000000000   0.28700000000000   0.66900000000000   Al   (4b)
   0.28700000000000  -0.27100000000000   0.66900000000000   Al   (4b)
  -0.15800000000000   0.42000000000000   0.53200000000000   Al   (4b)
   0.42000000000000  -0.15800000000000   0.53200000000000   Al   (4b)
   0.08600000000000   0.08600000000000   0.38200000000000   Co   (2a)
   0.60400000000000   0.60400000000000   0.37300000000000   Co   (2a)
  -0.09100000000000  -0.09100000000000   0.01200000000000   Co   (2a)
   0.59700000000000   0.59700000000000   0.01800000000000   Co   (2a)
  -0.08700000000000  -0.08700000000000   0.61700000000000   Co   (2a)
   0.39500000000000   0.39500000000000   0.62600000000000   Co   (2a)
   0.09000000000000   0.09000000000000  -0.01300000000000   Co   (2a)
   0.40200000000000   0.40200000000000  -0.01800000000000   Co   (2a)
   0.02600000000000   0.61000000000000   0.27700000000000   Co   (4b)
   0.61000000000000   0.02600000000000   0.27700000000000   Co   (4b)
  -0.04700000000000   0.41100000000000   0.72200000000000   Co   (4b)
   0.41100000000000  -0.04700000000000   0.72200000000000   Co   (4b)
```

$TaTi_3$ (BCC SQS-16): AB3_mC32_8_4a_12a - CIF

```
# AFLOW.org Repositories
# TaTi/AB3_mC32_8_4a_12a-001.AB params=13.8779590179,0.333333333333,
      ↪ 0.666666666667,109.471220634,-0.0,0.5,0.5,0.75,0.375,0.9375,
      ↪ 0.125,0.8125,0.875,0.1875,0.625,0.0625,0.75,0.375,0.875,0.6875
      ↪ ,-0.0,-0.0,0.5,0.25,0.625,0.5625,0.75,0.875,0.25,0.125,0.375,
      ↪ 0.4375,0.125,0.3125,0.25,0.625 SG=8 [ANRL doi: 10.1016/
      ↪ j.commatsci.2017.01.017 (part 1), doi: 10.1016/
      ↪ j.commatsci.2018.10.043 (part 2)]
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_pd_phase_name AB3_mC32_8_4a_12a-001.AB

_chemical_name_mineral 'TaTi3'
_chemical_formula_sum 'Ta Ti3'

loop_
_publ_author_name
 'C. Jiang'
 'C. Wolverton'
 'J. Sofo'
 'L.-Q. Chen'
 'Z.-K. Liu'
_journal_name_full_name
;
 Physical Review B
;
_journal_volume 69
_journal_year 2004
_journal_page_first 214202
_journal_page_last 214202
_publ_Section_title
;
 First-principles study of binary bcc alloys using special quasirandom
      ↪ structures
;

_aflow_title 'TaTi$_{3}$ (BCC SQS-16) Structure'
_aflow_proto 'AB3_mC32_8_4a_12a'
_aflow_params 'a,b/a,c/a,\beta,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3},x_{4
      ↪ },z_{4},x_{5},z_{5},x_{6},z_{6},x_{7},z_{7},x_{8},z_{8},x_{9},
      ↪ z_{9},x_{10},z_{10},x_{11},z_{11},x_{12},z_{12},x_{13},z_{13},
      ↪ x_{14},z_{14},x_{15},z_{15},x_{16},z_{16}'
_aflow_params_values '13.8779590179,0.333333333333,0.666666666667,
      ↪ 109.471220634,0.0,0.5,0.5,0.25,0.625,0.0625,0.875,0.1875,0.125,
      ↪ 0.8125,0.375,0.9375,0.25,0.625,0.125,0.3125,0.0,0.0,0.5,0.75,
      ↪ 0.375,0.4375,0.25,0.125,0.75,0.875,0.625,0.5625,0.875,0.6875,
      ↪ 0.75,0.375'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mC32'

_cell_length_a 13.8779590179
_cell_length_b 4.6259863393
_cell_length_c 9.2519726786
_cell_angle_alpha 90.0000000000
_cell_angle_beta 109.4712206340
```

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_symmetry_Int_Tables_Number   8
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_symmetry_equiv_pos_site_id
_symmetry_equiv_pos_as_xyz
  1  x,y,z
  2  x,-y,z
  3  x+1/2,y+1/2,z
  4  x+1/2,-y+1/2,z
loop_
 _atom_site_label
 _atom_site_occupancy
 _atom_site_fract_x
 _atom_site_fract_y
 _atom_site_fract_z
 _atom_site_thermal_displace_type
 _atom_site_B_iso_or_equiv
 _atom_site_type_symbol
 _atom_site_symmetry_multiplicity
 _atom_site_Wyckoff_label
Ta1  1.0000000000  0.0000000000  0.0000000000  0.5000000000  Biso  1.0  Ta  2  a
Ta2  1.0000000000  0.5000000000  -0.0000000000  0.2500000000  Biso  1.0  Ta  2  a
Ta3  1.0000000000  0.6250000000  0.0000000000  0.0625000000  Biso  1.0  Ta  2  a
Ta4  1.0000000000  0.8750000000  0.0000000000  0.1875000000  Biso  1.0  Ta  2  a
Ti1  1.0000000000  0.1250000000  -0.0000000000  0.8125000000  Biso  1.0  Ti  2  a
Ti2  1.0000000000  0.3750000000  -0.0000000000  0.9375000000  Biso  1.0  Ti  2  a
Ti3  1.0000000000  0.2500000000  -0.0000000000  0.6250000000  Biso  1.0  Ti  2  a
Ti4  1.0000000000  0.1250000000  -0.0000000000  0.3125000000  Biso  1.0  Ti  2  a
Ti5  1.0000000000  0.0000000000  0.0000000000  0.0000000000  Biso  1.0  Ti  2  a
Ti6  1.0000000000  0.5000000000  -0.0000000000  0.7500000000  Biso  1.0  Ti  2  a
Ti7  1.0000000000  0.3750000000  -0.0000000000  0.4375000000  Biso  1.0  Ti  2  a
Ti8  1.0000000000  0.2500000000  -0.0000000000  0.1250000000  Biso  1.0  Ti  2  a
Ti9  1.0000000000  0.7500000000  0.0000000000  0.8750000000  Biso  1.0  Ti  2  a
Ti10  1.0000000000  0.6250000000  0.0000000000  0.5625000000  Biso  1.0  Ti  2  a
Ti11  1.0000000000  0.8750000000  0.0000000000  0.6875000000  Biso  1.0  Ti  2  a
Ti12  1.0000000000  0.7500000000  0.0000000000  0.3750000000  Biso  1.0  Ti  2  a
```

$TaTi_3$ (BCC SQS-16): AB3_mC32_8_4a_12a - POSCAR

```
AB3_mC32_8_4a_12a & a,b/a,c/a,beta,x1,z1,x2,z2,x3,z3,x4,z4,x5,z5,x6,z6,
    ↪ x7,z7,x8,z8,x9,z9,x10,z10,x11,z11,x12,z12,x13,z13,x14,z14,x15,
    ↪ z15,x16,z16 --params=13.8779590179,0.333333333333,
    ↪ 0.666666666667,109.471220634,0.0,0.5,0.5,0.25,0.625,0.0625,
    ↪ 0.875,0.1875,0.125,0.8125,0.375,0.9375,0.25,0.625,0.125,0.3125,
    ↪ 0.0,0.0,0.5,0.75,0.375,0.4375,0.25,0.125,0.75,0.875,0.625,
    ↪ 0.5625,0.875,0.6875,0.75,0.375 & Cm C_{s}^{3} #8 (a^16) & mC32
    ↪ & None & TaTi3 & TaTi3 & C. Jiang et al., Phys. Rev. B 69,
    ↪ 214202(2004)
   1.00000000000000
   6.93897950895000   -2.31299316965000    0.00000000000000
   6.93897950895000    2.31299316965000    0.00000000000000
  -3.08399089279196    0.00000000000000    8.72284349388071
     Ta      Ti
      4      12
Direct
    0.00000000000000    0.00000000000000    0.50000000000000    Ta    (2a)
    0.50000000000000    0.50000000000000    0.25000000000000    Ta    (2a)
    0.62500000000000    0.62500000000000    0.06250000000000    Ta    (2a)
    0.87500000000000    0.87500000000000    0.18750000000000    Ta    (2a)
    0.12500000000000    0.12500000000000    0.81250000000000    Ti    (2a)
    0.37500000000000    0.37500000000000    0.93750000000000    Ti    (2a)
    0.25000000000000    0.25000000000000    0.62500000000000    Ti    (2a)
    0.12500000000000    0.12500000000000    0.31250000000000    Ti    (2a)
    0.00000000000000    0.00000000000000    0.00000000000000    Ti    (2a)
    0.50000000000000    0.50000000000000    0.75000000000000    Ti    (2a)
    0.37500000000000    0.37500000000000    0.43750000000000    Ti    (2a)
    0.25000000000000    0.25000000000000    0.12500000000000    Ti    (2a)
    0.75000000000000    0.75000000000000    0.87500000000000    Ti    (2a)
    0.62500000000000    0.62500000000000    0.56250000000000    Ti    (2a)
    0.87500000000000    0.87500000000000    0.68750000000000    Ti    (2a)
    0.75000000000000    0.75000000000000    0.37500000000000    Ti    (2a)
```

$TaTi_3$ (BCC SQS-16): AB3_mC32_8_4a_4a4b - CIF

```
# AFLOW.org Repositories
# TaTi/AB3_mC32_8_4a_4a4b-001.AB params=9.2519726786,1.0,0.707106781188,
    ↪ 90.0,0.0,0.0,0.5,0.25,0.25,0.5,0.5,0.25,0.75,0.5,0.0,0.75,
    ↪ 0.75,0.25,0.75,0.25,0.0,0.75,0.25,0.5,0.0,0.25,0.75,0.0,
    ↪ 0.25,0.25 SG=8 [ANRL doi: 10.1016/j.commatsci.2017.01.017 (part
    ↪ 1), doi: 10.1016/j.commatsci.2018.10.043 (part 2)]
data_TaTi
_pd_phase_name AB3_mC32_8_4a_4a4b-001.AB

_chemical_name_mineral 'TaTi3'
_chemical_formula_sum 'Ta Ti3'

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 'T. Chakraborty'
 'J. Rogal'
 'R. Drautz'
_journal_name_full_name
;
 Physical Review B
;
_journal_volume 94
_journal_year 2016
_journal_page_first 224104
_journal_page_last 224104
_publ_Section_title
;
 Unraveling the composition dependence of the martensitic transformation
    ↪ temperature: A first-principles study of Ti-Ta alloys
;
_aflow_title 'TaTi${_}3$ (BCC SQS-16) Structure'
_aflow_proto 'AB3_mC32_8_4a_4a4b'
_aflow_params 'a,b/a,c/a,\beta,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3},x_{4
    ↪ },z_{4},x_{5},z_{5},x_{6},z_{6},x_{7},z_{7},x_{8},z_{8},x_{9},
    ↪ y_{9},z_{9},x_{10},y_{10},z_{10},x_{11},y_{11},z_{11},x_{12},y_
    ↪ {12},z_{12}'
_aflow_params_values '9.2519726786,1.0,0.707106781188,90.0,-0.0,0.0,-0.0
    ↪ ,0.5,0.75,0.75,0.5,0.5,0.75,0.25,0.5,0.0,0.25,0.25,0.25,0.75,
    ↪ 0.75,0.75,0.0,0.75,0.75,0.5,0.5,0.75,0.25,0.5,0.75,0.75'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mC32'

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_cell_length_c   6.5421326204
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_cell_angle_gamma   90.0000000000
_symmetry_space_group_name_H-M   'Cm'
_symmetry_Int_Tables_Number   8
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_symmetry_equiv_pos_as_xyz
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  2  x,-y,z
  3  x+1/2,y+1/2,z
  4  x+1/2,-y+1/2,z
loop_
 _atom_site_label
 _atom_site_occupancy
 _atom_site_fract_x
 _atom_site_fract_y
 _atom_site_fract_z
 _atom_site_thermal_displace_type
 _atom_site_B_iso_or_equiv
 _atom_site_type_symbol
 _atom_site_symmetry_multiplicity
 _atom_site_Wyckoff_label
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Ta2  1.0000000000  -0.0000000000  0.0000000000  0.5000000000  Biso  1.0  Ta  2  a
Ta3  1.0000000000  0.7500000000  0.0000000000  0.7500000000  Biso  1.0  Ta  2  a
Ta4  1.0000000000  0.5000000000  -0.0000000000  0.5000000000  Biso  1.0  Ta  2  a
Ti1  1.0000000000  0.7500000000  -0.0000000000  0.2500000000  Biso  1.0  Ti  2  a
Ti2  1.0000000000  0.5000000000  -0.0000000000  0.0000000000  Biso  1.0  Ti  2  a
Ti3  1.0000000000  0.2500000000  -0.0000000000  0.2500000000  Biso  1.0  Ti  2  a
Ti4  1.0000000000  0.2500000000  -0.0000000000  0.7500000000  Biso  1.0  Ti  2  a
Ti5  1.0000000000  0.7500000000  0.7500000000  0.0000000000  Biso  1.0  Ti  4  b
Ti6  1.0000000000  0.7500000000  0.7500000000  0.5000000000  Biso  1.0  Ti  4  b
Ti7  1.0000000000  0.5000000000  0.7500000000  0.2500000000  Biso  1.0  Ti  4  b
Ti8  1.0000000000  0.5000000000  0.7500000000  0.7500000000  Biso  1.0  Ti  4  b
```

$TaTi_3$ (BCC SQS-16): AB3_mC32_8_4a_4a4b - POSCAR

```
AB3_mC32_8_4a_4a4b & a,b/a,c/a,beta,x1,z1,x2,z2,x3,z3,x4,z4,x5,z5,x6,z6,
    ↪ x7,z7,x8,z8,x9,y9,z9,x10,y10,z10,x11,y11,z11,x12,y12,z12 --
    ↪ params=9.2519726786,1.0,0.707106781188,90.0,-0.0,0.0,-0.0,0.5,
    ↪ 0.75,0.75,0.5,0.5,0.75,0.25,0.5,0.0,0.25,0.25,0.25,0.75,0.75,
    ↪ 0.75,0.0,0.75,0.75,0.5,0.5,0.75,0.25,0.5,0.75,0.75 & Cm C_{s}^{
    ↪ 3} #8 (a^8b^4) & mC32 & None & TaTi3 & TaTi3 & T. Chakraborty
    ↪ and J. Rogal and R. Drautz, Phys. Rev. B 94, 224104(2016)
   1.00000000000000
   4.62598633930000   -4.62598633930000    0.00000000000000
   4.62598633930000    4.62598633930000    0.00000000000000
   0.00000000000000    0.00000000000000    6.54213262040000
     Ta      Ti
      4      12
Direct
    0.00000000000000    0.00000000000000    0.00000000000000    Ta    (2a)
    0.00000000000000    0.00000000000000    0.50000000000000    Ta    (2a)
    0.75000000000000    0.75000000000000    0.75000000000000    Ta    (2a)
    0.50000000000000    0.50000000000000    0.50000000000000    Ta    (2a)
    0.75000000000000    0.75000000000000    0.25000000000000    Ti    (2a)
    0.50000000000000    0.50000000000000    0.00000000000000    Ti    (2a)
    0.25000000000000    0.25000000000000    0.25000000000000    Ti    (2a)
    0.25000000000000    0.25000000000000    0.75000000000000    Ti    (2a)
    0.00000000000000    1.50000000000000    0.00000000000000    Ti    (4b)
    1.50000000000000    0.00000000000000    0.00000000000000    Ti    (4b)
    0.00000000000000    1.50000000000000    0.50000000000000    Ti    (4b)
    1.50000000000000    0.00000000000000    0.50000000000000    Ti    (4b)
   -0.25000000000000    1.25000000000000    0.25000000000000    Ti    (4b)
    1.25000000000000   -0.25000000000000    0.25000000000000    Ti    (4b)
   -0.25000000000000    1.25000000000000    0.75000000000000    Ti    (4b)
    1.25000000000000   -0.25000000000000    0.75000000000000    Ti    (4b)
```

$F5_{11}$ (KNO_2) (*obsolete*): ABC2_mC8_8_a_a_b - CIF

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 'G. E. Ziegler'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 94
_journal_year 1936
_journal_page_first 491
_journal_page_last 499
```

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_publ_Section_title
;
 The Crystal Structure of Potassium Nitrite, KNO${2}$
;

_aflow_title 'F5_{11}$ (KNO${2}$) ({\em{obsolete}}) Structure'
_aflow_proto 'ABC2_mC8_8_a_a_b'
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    ↪ }'
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_aflow_Strukturbericht '$F5_{11}$'
_aflow_Pearson 'mC8'

_symmetry_space_group_name_H-M "C 1 m 1"
_symmetry_Int_Tables_number 8

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_cell_length_b      4.99000
_cell_length_c      4.45000
_cell_angle_alpha   90.00000
_cell_angle_beta    114.83330
_cell_angle_gamma   90.00000

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_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,z
3 x+1/2,y+1/2,z
4 x+1/2,-y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
K1 K 2 a 0.00000 0.00000 0.00000 1.00000
N1 N 2 a 0.48600 0.00000 0.50000 1.00000
O1 O 4 b -0.08300 0.30600 0.44400 1.00000
```

$F5_{11}$ (KNO_2) (*obsolete*): ABC2_mC8_8_a_a_b - POSCAR

```
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    ↪ 0.682626538988,0.608755129959,114.8333,0.0,0.0,0.486,0.5,-0.083
    ↪ ,0.306,0.444 & Cm C_{s}^{3} #8 (a^2b) & mC8 & $F5_{11}$ & KNO2
    ↪ & Potassium nitrite & G. E. Ziegler, Zeitschrift f{"u}r
    ↪ Kristallographie - Crystalline Materials 94, 491-499 (1936)
   1.00000000000000
   3.65500000000000   -2.49500000000000    0.00000000000000
   3.65500000000000    2.49500000000000    0.00000000000000
  -1.86890925105819    0.00000000000000    4.03852426157243
     K     N     O
     1     1     2
Direct
   0.00000000000000    0.00000000000000    0.00000000000000    K    (2a)
   0.48600000000000    0.48600000000000    0.50000000000000    N    (2a)
  -0.38900000000000    0.22300000000000    0.44400000000000    O    (4b)
   0.22300000000000   -0.38900000000000    0.44400000000000    O    (4b)
```

Nacrite [$Al_2Si_2O_5(OH)_4$, $S5_4$]: A2B4C9D2_mC68_9_2a_4a_9a_2a - CIF

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loop_
_publ_author_name
 'A. P. Zhukhlistov'
_journal_name_full_name
;
 Crystallography Reports
;
_journal_volume 53
_journal_year 2008
_journal_page_first 76
_journal_page_last 82
_publ_Section_title
;
 Crystal structure of nacrite from the electron diffraction data
;

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    ↪ Structure'
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    ↪ z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{
    ↪ 10},y_{10},z_{10},x_{11},y_{11},z_{11},x_{12},y_{12},z_{12},x_{
    ↪ 13},y_{13},z_{13},x_{14},y_{14},z_{14},x_{15},y_{15},z_{15},x_{
    ↪ 16},y_{16},z_{16},x_{17},y_{17},z_{17}'
_aflow_params_values '8.91,0.577328843996,1.63782267116,100.5,0.8275,
    ↪ 0.4297,-0.0002,0.1628,0.4245,0.0004,0.426,0.318,0.096,0.599,
    ↪ 0.41,0.378,0.3,0.346,0.38,-0.001,0.386,0.38,0.1349,0.0011,
    ↪ 0.2215,0.6809,0.0027,0.2362,-0.0881,0.3075,0.2371,0.0214,0.3111
    ↪ ,0.0788,0.7143,0.238,0.0794,0.3271,0.2379,0.0748,0.164,0.1153,-
    ↪ 0.066,0.7853,0.1207,-0.065,0.4772,0.0432,-0.0651,0.0632,0.2796,
    ↪ 0.1897,0.7342,0.2656,0.1907'
_aflow_Strukturbericht '$S5_{4}$'
_aflow_Pearson 'mC68'

_symmetry_space_group_name_H-M "C 1 c 1"
_symmetry_Int_Tables_number 9

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_cell_length_c      14.59300
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_cell_angle_gamma   90.00000

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_space_group_symop_operation_xyz
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2 x,-y,z+1/2
3 x+1/2,y+1/2,z
4 x+1/2,-y+1/2,z+1/2

loop_
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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
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Al2 Al 4 a 0.16280 0.42450 0.00040 1.00000
H1 H 4 a 0.42600 0.31800 0.09600 1.00000
H2 H 4 a 0.59900 0.41000 0.37800 1.00000
H3 H 4 a 0.30000 0.34600 0.38000 1.00000
H4 H 4 a -0.00100 0.38600 0.38000 1.00000
O1 O 4 a 0.13490 0.00110 0.22150 1.00000
O2 O 4 a 0.68090 0.00270 0.23620 1.00000
O3 O 4 a -0.08810 0.30750 0.23710 1.00000
O4 O 4 a 0.02140 0.31110 0.07880 1.00000
O5 O 4 a 0.71430 0.23800 0.07940 1.00000
O6 O 4 a 0.32710 0.23790 0.07480 1.00000
O7 O 4 a 0.16400 0.11530 -0.06600 1.00000
O8 O 4 a 0.78530 0.12070 -0.06500 1.00000
O9 O 4 a 0.47720 0.04320 -0.06510 1.00000
Si1 Si 4 a 0.06320 0.27960 0.18970 1.00000
Si2 Si 4 a 0.73420 0.26560 0.19070 1.00000
```

Nacrite [$Al_2Si_2O_5(OH)_4$, $S5_4$]: A2B4C9D2_mC68_9_2a_4a_9a_2a - POSCAR

```
A2B4C9D2_mC68_9_2a_4a_9a_2a & a,b/a,c/a,beta,x1,y1,z1,x2,y2,z2,x3,y3,z3,
    ↪ x4,y4,z4,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,
    ↪ z10,x11,y11,z11,x12,y12,z12,x13,y13,z13,x14,y14,z14,x15,y15,z15
    ↪ ,x16,y16,z16,x17,y17,z17 --params=8.91,0.577328843996,
    ↪ 1.63782267116,100.5,0.8275,0.4297,-0.0002,0.1628,0.4245,0.0004,
    ↪ 0.426,0.318,0.096,0.599,0.41,0.378,0.3,0.346,0.38,-0.001,0.386,
    ↪ 0.38,0.1349,0.0011,0.2215,0.6809,0.0027,0.2362,-0.0881,0.3075,
    ↪ 0.2371,0.0214,0.3111,0.0788,0.7143,0.238,0.0794,0.3271,0.2379,
    ↪ 0.0748,0.164,0.1153,-0.066,0.7853,0.1207,-0.065,0.4772,0.0432,-
    ↪ 0.0651,0.0632,0.2796,0.1897,0.7342,0.2656,0.1907 & Cc C_{s}^{4}
    ↪ #9 (a^17) & mC68 & $S5_{4}$ & Al2H4O9Si2 & Nacrite & A. P.
    ↪ Zhukhlistov, Crystal. Rep. 53, 76-82 (2008)
   1.00000000000000
   4.45500000000000   -2.57200000000000    0.00000000000000
   4.45500000000000    2.57200000000000    0.00000000000000
  -2.65936302350691    0.00000000000000   14.34863886608080
     Al    H    O    Si
      4    8   18     4
Direct
   0.39780000000000    1.25720000000000   -0.00020000000000   Al    (4a)
   1.25720000000000    0.39780000000000    0.49980000000000   Al    (4a)
  -0.26170000000000    0.58730000000000    0.00040000000000   Al    (4a)
   0.58730000000000   -0.26170000000000    0.50040000000000   Al    (4a)
   0.10800000000000    0.74400000000000    0.09600000000000    H    (4a)
   0.74400000000000    0.10800000000000    0.59600000000000    H    (4a)
   0.18900000000000    1.00900000000000    0.37800000000000    H    (4a)
   1.00900000000000    0.18900000000000    0.87800000000000    H    (4a)
  -0.04600000000000    0.64600000000000    0.38000000000000    H    (4a)
   0.64600000000000   -0.04600000000000    0.88000000000000    H    (4a)
  -0.38700000000000    0.38500000000000    0.38000000000000    H    (4a)
   0.38500000000000   -0.38700000000000    0.88000000000000    H    (4a)
   0.13380000000000    0.13600000000000    0.22150000000000    O    (4a)
   0.13600000000000    0.13380000000000    0.72150000000000    O    (4a)
   0.67820000000000    0.68360000000000    0.23620000000000    O    (4a)
   0.68360000000000    0.67820000000000    0.73620000000000    O    (4a)
  -0.39560000000000    0.21940000000000    0.23710000000000    O    (4a)
   0.21940000000000   -0.39560000000000    0.73710000000000    O    (4a)
  -0.28970000000000    0.33250000000000    0.07880000000000    O    (4a)
   0.33250000000000   -0.28970000000000    0.57880000000000    O    (4a)
   0.47630000000000    0.95230000000000    0.07940000000000    O    (4a)
   0.95230000000000    0.47630000000000    0.57940000000000    O    (4a)
   0.08920000000000    0.56500000000000    0.07480000000000    O    (4a)
   0.56500000000000    0.08920000000000    0.57480000000000    O    (4a)
   0.04870000000000    0.27930000000000   -0.06600000000000    O    (4a)
   0.27930000000000    0.04870000000000    0.43400000000000    O    (4a)
   0.66460000000000    0.90600000000000   -0.06500000000000    O    (4a)
   0.90600000000000    0.66460000000000    0.43500000000000    O    (4a)
   0.43400000000000    0.52040000000000   -0.06510000000000    O    (4a)
   0.52040000000000    0.43400000000000    0.43490000000000    O    (4a)
  -0.21640000000000    0.34280000000000    0.18970000000000   Si    (4a)
   0.34280000000000   -0.21640000000000    0.68970000000000   Si    (4a)
   0.46860000000000    0.99980000000000    0.19070000000000   Si    (4a)
   0.99980000000000    0.46860000000000    0.69070000000000   Si    (4a)
```

Chrysotile ($Mg_3Si_2O_5(OH)_4$): A3B5C4D2_mC56_9_3a_5a_4a_2a - CIF

```
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_chemical_name_mineral 'Chrysotile'
_chemical_formula_sum 'Mg3 O5 (OH)4 Si2'

loop_
_publ_author_name
 'G. Falini'
 'E. Foresti'
 'M. Gazzano'
 'A. F. Gualtieri'
 'M. Leoni'
 'I G. Lesci'
 'N. Roveri'
_journal_name_full_name
;
 Chemistry - A European Journal
;
_journal_volume 10
_journal_year 2004
_journal_page_first 3043
_journal_page_last 3049
_publ_Section_title
;
 Tubular-Shaped Stoichiometric Chrysotile Nanocrystals
;

_aflow_title 'Chrysotile (Mg${_}{3}$Si${_}{2}$O${_}{5}$(OH)${_}{4}$) Structure'
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      ↪ z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{
      ↪ 10},y_{10},z_{10},x_{11},y_{11},z_{11},x_{12},y_{12},z_{12},x_{
      ↪ 13},y_{13},z_{13},x_{14},y_{14},z_{14}'
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      ↪ 0.198,0.2303,0.3747,0.3771,0.2327,0.8818,0.5235,0.2204,0.0726,
      ↪ 0.1978,0.0,0.2387,0.475,0.0023,0.7395,0.4214,0.0063,0.0438,
      ↪ 0.3596,0.156,0.532,0.537,0.155,0.7299,0.363,0.2878,0.2395,0.197
      ↪ ,0.2791,0.5334,0.2053,0.1807,0.2025,0.5416,0.296,0.0257,0.3641,
      ↪ 0.041,0.5208,0.5335,0.04'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mC56'

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_symmetry_Int_Tables_number 9

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_cell_angle_alpha 90.00000
_cell_angle_beta  93.66000
_cell_angle_gamma 90.00000

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_space_group_symop_id
_space_group_symop_operation_xyz
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2 x,-y,z+1/2
3 x+1/2,y+1/2,z
4 x+1/2,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
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Mg2 Mg 4 a 0.37470 0.37710 0.23270 1.00000
Mg3 Mg 4 a 0.88180 0.52350 0.22040 1.00000
O1 O 4 a 0.07260 0.19780 0.00000 1.00000
O2 O 4 a 0.23870 0.47500 0.00230 1.00000
O3 O 4 a 0.73950 0.42140 0.00630 1.00000
O4 O 4 a 0.04380 0.35960 0.15600 1.00000
O5 O 4 a 0.53200 0.53700 0.15500 1.00000
OH1 OH 4 a 0.72990 0.36300 0.28780 1.00000
OH2 OH 4 a 0.23950 0.19700 0.27910 1.00000
OH3 OH 4 a 0.53340 0.20530 0.18070 1.00000
OH4 OH 4 a 0.20250 0.54160 0.29600 1.00000
Si1 Si 4 a 0.02570 0.36410 0.04100 1.00000
Si2 Si 4 a 0.52080 0.53350 0.04000 1.00000
```

Chrysotile ($Mg_3Si_2O_5(OH)_4$): A3B5C4D2_mC56_9_3a_5a_4a_2a - POSCAR

```
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      ↪ x4,y4,z4,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,
      ↪ z10,x11,y11,z11,x12,y12,z12,x13,y13,z13,x14,y14,z14 --params=
      ↪ 5.34,1.73052434457,2.75074906367,93.66,0.8852,0.198,0.2303,
      ↪ 0.3747,0.3771,0.2327,0.8818,0.5235,0.2204,0.0726,0.1978,0.0,
      ↪ 0.2387,0.475,0.0023,0.7395,0.4214,0.0063,0.0438,0.3596,0.156,
      ↪ 0.532,0.537,0.155,0.7299,0.363,0.2878,0.2395,0.197,0.2791,
      ↪ 0.5334,0.2053,0.1807,0.2025,0.5416,0.296,0.0257,0.3641,0.041,
      ↪ 0.5208,0.5335,0.04 & Cc C_{s}^{4} #9 (a^14) & mC56 & None &
      ↪ Mg3O5(OH)4Si2 & Chrysotile & G. Falini et al., Chem. Eur. J. 10
      ↪ , 3043-3049 (2004)
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   2.67000000000000    4.62050000000000    0.00000000000000
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     Mg      O     OH     Si
     6     10      8      4
Direct
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   1.08320000000000   0.68720000000000   0.73030000000000  Mg   (4a)
  -0.00240000000000   0.75180000000000   0.23270000000000  Mg   (4a)
   0.75180000000000  -0.00240000000000   0.73270000000000  Mg   (4a)
   0.35830000000000   1.40530000000000   0.22040000000000  Mg   (4a)
   1.40530000000000   0.35830000000000   0.72040000000000  Mg   (4a)
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   0.27040000000000  -0.12520000000000   0.50000000000000  O    (4a)
  -0.23630000000000   0.71370000000000   0.00230000000000  O    (4a)
   0.71370000000000  -0.23630000000000   0.50230000000000  O    (4a)
   0.31810000000000   1.16090000000000   0.00630000000000  O    (4a)
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  -0.31580000000000   0.40340000000000   0.15600000000000  O    (4a)
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  -0.00500000000000   1.06900000000000   0.15500000000000  O    (4a)
   1.06900000000000  -0.00500000000000   0.65500000000000  O    (4a)
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   0.04250000000000   0.43650000000000   0.27910000000000  OH   (4a)
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   0.73870000000000   0.32810000000000   0.68070000000000  OH   (4a)
  -0.33910000000000   0.74410000000000   0.29600000000000  OH   (4a)
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  -0.33840000000000   0.38980000000000   0.04100000000000  Si   (4a)
   0.38980000000000  -0.33840000000000   0.54100000000000  Si   (4a)
  -0.01270000000000   1.05430000000000   0.04000000000000  Si   (4a)
   1.05430000000000  -0.01270000000000   0.54000000000000  Si   (4a)
```

$Cs_6W_{11}O_{36}$: A6B36C11_mC212_9_6a_36a_11a - CIF

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 'F. Marumo'
 'S. Iwai'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 34
_journal_year 1978
_journal_page_first 50
_journal_page_last 54
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;
 The Crystal Structure of Cs${_}{6}$W${_}{11}$O${_}{36}$
;

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      ↪ z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{
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      ↪ 16},y_{16},z_{16},x_{17},y_{17},z_{17},x_{18},y_{18},z_{18},x_{
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      ↪ 31},y_{31},z_{31},x_{32},y_{32},z_{32},x_{33},y_{33},z_{33},x_{
      ↪ 34},y_{34},z_{34},x_{35},y_{35},z_{35},x_{36},y_{36},z_{36},x_{
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      ↪ 40},y_{40},z_{40},x_{41},y_{41},z_{41},x_{42},y_{42},z_{42},x_{
      ↪ 43},y_{43},z_{43},x_{44},y_{44},z_{44},x_{45},y_{45},z_{45},x_{
      ↪ 46},y_{46},z_{46},x_{47},y_{47},z_{47},x_{48},y_{48},z_{48},x_{
      ↪ 49},y_{49},z_{49},x_{50},y_{50},z_{50},x_{51},y_{51},z_{51},x_{
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      ↪ 0.0025,0.8992,0.4748,0.4942,-0.0638,0.058,0.0037,0.0625,0.528,
      ↪ 0.01,-0.044,0.082,0.48,0.667,0.731,0.007,0.892,0.749,0.02,0.791
      ↪ ,0.814,0.009,0.007,0.207,0.324,-0.048,0.701,0.203,-0.048,0.053,
      ↪ 0.311,0.894,0.561,0.202,0.888,0.152,0.318,0.832,0.618,0.168,
      ↪ 0.836,0.372,0.326,0.852,0.859,0.191,0.847,0.432,0.303,0.791,-
      ↪ 0.047,0.202,0.784,0.64,0.2,0.012,0.127,0.314,0.006,0.414,0.164
      ↪ ,-0.014,-0.093,0.313,-0.013,0.411,0.187,0.667,0.896,0.315,0.668
      ↪ ,0.185,0.174,0.652,0.689,0.312,0.646,0.091,0.215,0.713,0.579,
      ↪ 0.307,0.707,0.804,0.323,0.044,0.317,0.172,0.048,0.487,0.188,
      ↪ 0.611,-0.013,0.33,0.607,0.219,0.004,-0.014,-0.057,0.478,0.832,
      ↪ 0.537,0.467,0.846,0.797,0.497,0.712,0.292,0.016,0.609,0.495,
      ↪ 0.498,0.649,0.502,0.013,0.05,0.25,0.2342,0.0,0.7808,0.2354,0.0,
      ↪ 0.483,0.2476,0.8347,0.5381,0.2682,0.6654,0.0553,0.2539,0.6648,-
      ↪ 0.0022,0.2447,0.8349,0.6325,0.0051,-0.0841,-0.0894,0.495,0.0835
      ↪ ,0.7428,0.0229,0.8341,0.7913,0.4872,0.6644,0.5166,0.0298,0.0'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mC212'

_symmetry_space_group_name_H-M "C 1 c 1"
_symmetry_Int_Tables_number 9

_cell_length_a    12.57700
_cell_length_b    7.26100
_cell_length_c    37.80000
_cell_angle_alpha 90.00000
```

```
_cell_angle_beta   102.81000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,z+1/2
3 x+1/2,y+1/2,z
4 x+1/2,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cs1 Cs 4 a 0.83100 0.00010 0.72060 1.00000
Cs2 Cs 4 a 0.20290 0.00170 0.77860 1.00000
Cs3 Cs 4 a 0.24820 0.49740 0.09870 1.00000
Cs4 Cs 4 a 0.28170 0.00250 0.89920 1.00000
Cs5 Cs 4 a 0.47480 0.49420 -0.06380 1.00000
Cs6 Cs 4 a 0.05800 0.00370 0.06250 1.00000
O1 O 4 a 0.52800 0.01000 -0.04400 1.00000
O2 O 4 a 0.08200 0.48000 0.66700 1.00000
O3 O 4 a 0.73100 0.00700 0.89200 1.00000
O4 O 4 a 0.74900 0.02000 0.79100 1.00000
O5 O 4 a 0.81400 0.00900 0.00700 1.00000
O6 O 4 a 0.20700 0.32400 -0.04800 1.00000
O7 O 4 a 0.70100 0.20300 -0.04800 1.00000
O8 O 4 a 0.05300 0.31100 0.89400 1.00000
O9 O 4 a 0.56100 0.20200 0.88800 1.00000
O10 O 4 a 0.15200 0.31800 0.83200 1.00000
O11 O 4 a 0.61800 0.16800 0.83600 1.00000
O12 O 4 a 0.37200 0.32600 0.85200 1.00000
O13 O 4 a 0.85900 0.19100 0.84700 1.00000
O14 O 4 a 0.43200 0.30300 0.79100 1.00000
O15 O 4 a -0.04700 0.20200 0.78400 1.00000
O16 O 4 a 0.64000 0.20000 0.01200 1.00000
O17 O 4 a 0.12700 0.31400 0.00600 1.00000
O18 O 4 a 0.41400 0.16400 -0.01400 1.00000
O19 O 4 a -0.09300 0.31300 -0.01300 1.00000
O20 O 4 a 0.41100 0.18700 0.66700 1.00000
O21 O 4 a 0.89600 0.31500 0.66800 1.00000
O22 O 4 a 0.18500 0.17400 0.65200 1.00000
O23 O 4 a 0.68900 0.31200 0.64600 1.00000
O24 O 4 a 0.09100 0.21500 0.71300 1.00000
O25 O 4 a 0.57900 0.30700 0.70700 1.00000
O26 O 4 a 0.80400 0.32300 0.04400 1.00000
O27 O 4 a 0.31700 0.17200 0.04800 1.00000
O28 O 4 a 0.48700 0.18800 0.61100 1.00000
O29 O 4 a -0.01300 0.33000 0.60700 1.00000
O30 O 4 a 0.21900 0.00400 -0.01400 1.00000
O31 O 4 a -0.05700 0.47800 0.83200 1.00000
O32 O 4 a 0.53700 0.46700 0.84600 1.00000
O33 O 4 a 0.79700 0.49700 0.71200 1.00000
O34 O 4 a 0.29200 0.01600 0.60900 1.00000
O35 O 4 a 0.49500 0.49800 0.64900 1.00000
O36 O 4 a 0.50200 0.01300 0.05000 1.00000
W1 W 4 a 0.25000 0.23420 0.00000 1.00000
W2 W 4 a 0.78080 0.23540 0.00000 1.00000
W3 W 4 a 0.48300 0.24760 0.83470 1.00000
W4 W 4 a 0.53810 0.26820 0.66540 1.00000
W5 W 4 a 0.05530 0.25390 0.66480 1.00000
W6 W 4 a -0.00220 0.24470 0.83490 1.00000
W7 W 4 a 0.63250 0.00510 -0.08410 1.00000
W8 W 4 a -0.08940 0.49500 0.08350 1.00000
W9 W 4 a 0.74280 0.02290 0.83410 1.00000
W10 W 4 a 0.79130 0.48720 0.66440 1.00000
W11 W 4 a 0.51660 0.02980 0.00000 1.00000
```

$Cs_6W_{11}O_{36}$: A6B36C11_mC212_9_6a_36a_11a - POSCAR

```
A6B36C11_mC212_9_6a_36a_11a & a,b/a,c/a,beta,x1,y1,z1,x2,y2,z2,x3,y3,z3,
    ↪ x4,y4,z4,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,
    ↪ z10,x11,y11,z11,x12,y12,z12,x13,y13,z13,x14,y14,z14,x15,y15,z15
    ↪ ,x16,y16,z16,x17,y17,z17,x18,y18,z18,x19,y19,z19,x20,y20,z20,
    ↪ x21,y21,z21,x22,y22,z22,x23,y23,z23,x24,y24,z24,x25,y25,z25,x26
    ↪ ,y26,z26,x27,y27,z27,x28,y28,z28,x29,y29,z29,x30,y30,z30,x31,
    ↪ y31,z31,x32,y32,z32,x33,y33,z33,x34,y34,z34,x35,y35,z35,x36,y36
    ↪ ,z36,x37,y37,z37,x38,y38,z38,x39,y39,z39,x40,y40,z40,x41,y41,
    ↪ z41,x42,y42,z42,x43,y43,z43,x44,y44,z44,x45,y45,z45,x46,y46,z46
    ↪ ,x47,y47,z47,x48,y48,z48,x49,y49,z49,x50,y50,z50,x51,y51,z51,
    ↪ x52,y52,z52,x53,y53,z53 --params=12.577,0.577323686094,
    ↪ 3.00548620498,102.81,0.831,0.0001,0.7206,0.2029,0.0017,0.7786,
    ↪ 0.2482,0.4974,0.0987,0.2817,0.0025,0.8992,0.4748,0.4942,-0.0638
    ↪ ,0.058,0.0037,0.0625,0.528,0.01,-0.044,0.082,0.48,0.667,0.731,
    ↪ 0.007,0.892,0.749,0.02,0.791,0.814,0.009,0.007,0.207,0.324,-
    ↪ 0.048,0.701,0.203,-0.048,0.053,0.311,0.894,0.561,0.202,0.888,
    ↪ 0.152,0.318,0.832,0.618,0.168,0.836,0.372,0.326,0.852,0.859,
    ↪ 0.191,0.847,0.432,0.303,0.791,-0.047,0.202,0.784,0.64,0.2,0.012
    ↪ ,0.127,0.314,0.006,0.414,0.164,-0.014,-0.093,0.313,-0.013,0.411
    ↪ ,0.187,0.667,0.896,0.315,0.668,0.185,0.174,0.652,0.689,0.312,
    ↪ 0.646,0.091,0.215,0.713,0.579,0.307,0.707,0.804,0.323,0.044,
    ↪ 0.317,0.172,0.048,0.487,0.188,0.611,-0.013,0.33,0.607,0.219,
    ↪ 0.004,-0.014,-0.057,0.478,0.832,0.537,0.467,0.846,0.797,0.497,
    ↪ 0.712,0.292,0.016,0.609,0.495,0.498,0.649,0.502,0.013,0.05,0.25
    ↪ ,0.2342,0.0,0.7808,0.2354,0.0,0.483,0.2476,0.8347,0.5381,0.2682
    ↪ ,0.6654,0.0553,0.2539,0.6648,-0.0022,0.2447,0.8349,0.6325,
    ↪ 0.0051,-0.0841,-0.0894,0.495,0.0835,0.7428,0.0229,0.8341,0.7913
    ↪ ,0.4872,0.6644,0.5166,0.0298,0.0 & Cc C_{s}^{4} #9 (a^53) &
    ↪ mC212 & None & Cs6O36W11 & Cs6O36W11 & K. Okada and F. Marumo
    ↪ and S. Iwai, Acta Crystallogr. Sect. B Struct. Sci. 34, 50-54 (
    ↪ 1978)
  1.00000000000000
  6.28850000000000  -3.63050000000000   0.00000000000000
  6.28850000000000   3.63050000000000   0.00000000000000
 -8.38096647873181   0.00000000000000  36.85918339955960
   Cs   O   W
   12  72  22
Direct
  0.83090000000000  0.83110000000000  0.72060000000000  Cs  (4a)
  0.83110000000000  0.83090000000000  1.22060000000000  Cs  (4a)
  0.20120000000000  0.20460000000000  0.77860000000000  Cs  (4a)
  0.20460000000000  0.20120000000000  1.27860000000000  Cs  (4a)
 -0.24920000000000  0.74560000000000  0.09870000000000  Cs  (4a)
  0.74560000000000 -0.24920000000000  0.59870000000000  Cs  (4a)
  0.27920000000000  0.28420000000000  0.89920000000000  Cs  (4a)
  0.28420000000000  0.27920000000000  1.39920000000000  Cs  (4a)
 -0.01940000000000  0.96900000000000 -0.06380000000000  Cs  (4a)
  0.96900000000000 -0.01940000000000  0.43620000000000  Cs  (4a)
  0.05430000000000  0.06170000000000  0.06250000000000  Cs  (4a)
  0.06170000000000  0.05430000000000  0.56250000000000  Cs  (4a)
  0.51800000000000  0.53800000000000 -0.04400000000000  O  (4a)
  0.53800000000000  0.51800000000000  0.45600000000000  O  (4a)
 -0.39800000000000  0.56200000000000  0.66700000000000  O  (4a)
  0.56200000000000 -0.39800000000000  1.16700000000000  O  (4a)
  0.72400000000000  0.73800000000000  0.89200000000000  O  (4a)
  0.73800000000000  0.72400000000000  1.39200000000000  O  (4a)
  0.72900000000000  0.76900000000000  0.79100000000000  O  (4a)
  0.76900000000000  0.72900000000000  1.29100000000000  O  (4a)
  0.80500000000000  0.82300000000000  0.00700000000000  O  (4a)
  0.82300000000000  0.80500000000000  0.50700000000000  O  (4a)
 -0.11700000000000  0.53100000000000 -0.04800000000000  O  (4a)
  0.53100000000000 -0.11700000000000  0.45200000000000  O  (4a)
  0.49800000000000  0.90400000000000 -0.04800000000000  O  (4a)
  0.90400000000000  0.49800000000000  0.45200000000000  O  (4a)
 -0.25800000000000  0.36400000000000  0.89400000000000  O  (4a)
  0.36400000000000 -0.25800000000000  1.39400000000000  O  (4a)
  0.35900000000000  0.76300000000000  0.88800000000000  O  (4a)
  0.76300000000000  0.35900000000000  1.38800000000000  O  (4a)
 -0.16600000000000  0.47000000000000  0.83200000000000  O  (4a)
  0.47000000000000 -0.16600000000000  1.33200000000000  O  (4a)
  0.45000000000000  0.78600000000000  0.83600000000000  O  (4a)
  0.78600000000000  0.45000000000000  1.33600000000000  O  (4a)
  0.04600000000000  0.69800000000000  0.85200000000000  O  (4a)
  0.69800000000000  0.04600000000000  1.35200000000000  O  (4a)
  0.66800000000000  1.05000000000000  0.84700000000000  O  (4a)
  1.05000000000000  0.66800000000000  1.34700000000000  O  (4a)
  0.12900000000000  0.73500000000000  0.79100000000000  O  (4a)
  0.73500000000000  0.12900000000000  1.29100000000000  O  (4a)
 -0.24900000000000  0.15500000000000  0.78400000000000  O  (4a)
  0.15500000000000 -0.24900000000000  1.28400000000000  O  (4a)
  0.44000000000000  0.84000000000000  0.01200000000000  O  (4a)
  0.84000000000000  0.44000000000000  0.51200000000000  O  (4a)
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  0.44100000000000 -0.18700000000000  0.50600000000000  O  (4a)
  0.25000000000000  0.57800000000000 -0.01400000000000  O  (4a)
  0.57800000000000  0.25000000000000  0.48600000000000  O  (4a)
 -0.40600000000000  0.22000000000000 -0.01300000000000  O  (4a)
  0.22000000000000 -0.40600000000000  0.48700000000000  O  (4a)
  0.22400000000000  0.59800000000000  0.66700000000000  O  (4a)
  0.59800000000000  0.22400000000000  1.16700000000000  O  (4a)
  0.58100000000000  1.21100000000000  0.66800000000000  O  (4a)
  1.21100000000000  0.58100000000000  1.16800000000000  O  (4a)
  0.01100000000000  0.35900000000000  0.65200000000000  O  (4a)
  0.35900000000000  0.01100000000000  1.15200000000000  O  (4a)
  0.37700000000000  1.00100000000000  0.64600000000000  O  (4a)
  1.00100000000000  0.37700000000000  1.14600000000000  O  (4a)
 -0.12400000000000  0.30600000000000  0.71300000000000  O  (4a)
  0.30600000000000 -0.12400000000000  1.21300000000000  O  (4a)
  0.27200000000000  0.88600000000000  0.70700000000000  O  (4a)
  0.88600000000000  0.27200000000000  1.20700000000000  O  (4a)
  0.48100000000000  1.12700000000000  0.04400000000000  O  (4a)
  1.12700000000000  0.48100000000000  0.54400000000000  O  (4a)
  0.14500000000000  0.48900000000000  0.04800000000000  O  (4a)
  0.48900000000000  0.14500000000000  0.54800000000000  O  (4a)
  0.29900000000000  0.67500000000000  0.61100000000000  O  (4a)
  0.67500000000000  0.29900000000000  1.11100000000000  O  (4a)
 -0.34300000000000  0.31700000000000  0.60700000000000  O  (4a)
  0.31700000000000 -0.34300000000000  1.10700000000000  O  (4a)
  0.21500000000000  0.22300000000000 -0.01400000000000  O  (4a)
  0.22300000000000  0.21500000000000  0.48600000000000  O  (4a)
 -0.53500000000000  0.42100000000000  0.83200000000000  O  (4a)
  0.42100000000000 -0.53500000000000  1.33200000000000  O  (4a)
  0.07000000000000  1.00400000000000  0.84600000000000  O  (4a)
  1.00400000000000  0.07000000000000  1.34600000000000  O  (4a)
  0.30000000000000  1.29400000000000  0.71200000000000  O  (4a)
  1.29400000000000  0.30000000000000  1.21200000000000  O  (4a)
  0.27600000000000  0.30800000000000  0.60900000000000  O  (4a)
  0.30800000000000  0.27600000000000  1.10900000000000  O  (4a)
 -0.00300000000000  0.99300000000000  0.64900000000000  O  (4a)
  0.99300000000000 -0.00300000000000  1.14900000000000  O  (4a)
  0.48900000000000  0.51500000000000  0.05000000000000  O  (4a)
  0.51500000000000  0.48900000000000  0.55000000000000  O  (4a)
  0.01580000000000  0.48420000000000  0.00000000000000  W  (4a)
  0.48420000000000  0.01580000000000  0.50000000000000  W  (4a)
  0.54540000000000  1.01620000000000  0.00000000000000  W  (4a)
  1.01620000000000  0.54540000000000  0.50000000000000  W  (4a)
  0.23540000000000  0.73060000000000  0.83470000000000  W  (4a)
  0.73060000000000  0.23540000000000  1.33470000000000  W  (4a)
  0.26990000000000  0.80630000000000  0.66540000000000  W  (4a)
  0.80630000000000  0.26990000000000  1.16540000000000  W  (4a)
 -0.19860000000000  0.30920000000000  0.66480000000000  W  (4a)
  0.30920000000000 -0.19860000000000  1.16480000000000  W  (4a)
 -0.24690000000000  0.24250000000000  0.83490000000000  W  (4a)
  0.24250000000000 -0.24690000000000  1.33490000000000  W  (4a)
```

```
   0.62740000000000   0.63760000000000  −0.08410000000000   W   (4a)
   0.63760000000000   0.62740000000000   0.41590000000000   W   (4a)
  −0.58440000000000   0.40560000000000   0.08350000000000   W   (4a)
   0.40560000000000  −0.58440000000000   0.58350000000000   W   (4a)
   0.71990000000000   0.76570000000000   0.83410000000000   W   (4a)
   0.76570000000000   0.71990000000000   1.33410000000000   W   (4a)
   0.30410000000000   1.27850000000000   0.66440000000000   W   (4a)
   1.27850000000000   0.30410000000000   1.16440000000000   W   (4a)
   0.48680000000000   0.54640000000000   0.00000000000000   W   (4a)
   0.54640000000000   0.48680000000000   0.50000000000000   W   (4a)
```

$K_2S_2O_5$ ($K0_1$): A2B5C2_mP18_11_2e_e2f_2e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'K2O5S2'
_chemical_formula_sum 'K2 O5 S2'

loop_
_publ_author_name
 'I-C. Chen'
 'Y. Wang'
_journal_name_full_name
;
 Acta Crystallographica Section C: Structural Chemistry
;
_journal_volume 40
_journal_year 1984
_journal_page_first 1780
_journal_page_last 1781
_publ_Section_title
;
 Reinvestigation of potassium pyrosulfite, K$_{2}$S$_{2}$O$_{5}$
;

_aflow_title 'K$_{2}$S$_{2}$O$_{5}$ ($K0_{1}$) Structure'
_aflow_proto 'A2B5C2_mP18_11_2e_e2f_2e'
_aflow_params 'a,b/a,c/a,\beta,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3},x_{4
      ↪ },z_{4},x_{5},z_{5},x_{6},y_{6},z_{6},x_{7},y_{7},z_{7}'
_aflow_params_values '6.921,0.890044791215,1.08900447912,102.79,0.231,-
      ↪ 0.0647,0.6398,0.67278,0.6555,0.0407,0.0291,0.33,0.7014,0.2384,
      ↪ 0.0741,0.0515,0.2343,0.635,0.053,0.3147'
_aflow_Strukturbericht '$K0_{1}$'
_aflow_Pearson 'mP18'

_symmetry_space_group_name_H-M "P 1 21/m 1"
_symmetry_Int_Tables_number 11

_cell_length_a     6.92100
_cell_length_b     6.16000
_cell_length_c     7.53700
_cell_angle_alpha  90.00000
_cell_angle_beta   102.79000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z
3 -x,-y,-z
4 x,-y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
K1 K 2 e 0.23100 0.25000 -0.06470 1.00000
K2 K 2 e 0.63980 0.25000 0.67278 1.00000
O1 O 2 e 0.65550 0.25000 0.04070 1.00000
S1 S 2 e 0.02910 0.25000 0.33000 1.00000
S2 S 2 e 0.70140 0.25000 0.23840 1.00000
O2 O 4 f 0.07410 0.05150 0.23430 1.00000
O3 O 4 f 0.63500 0.05300 0.31470 1.00000
```

$K_2S_2O_5$ ($K0_1$): A2B5C2_mP18_11_2e_e2f_2e - POSCAR

```
A2B5C2_mP18_11_2e_e2f_2e & a,b/a,c/a,beta,x1,z1,x2,z2,x3,z3,x4,z4,x5,z5,
      ↪ x6,y6,z6,x7,y7,z7 --params=6.921,0.890044791215,1.08900447912,
      ↪ 102.79,0.231,-0.0647,0.6398,0.67278,0.6555,0.0407,0.0291,0.33,
      ↪ 0.7014,0.2384,0.0741,0.0515,0.2343,0.635,0.053,0.3147 & P2_{1}/
      ↪ m C_{2h}^{2} #11 (e^5f^2) & mP18 & $K0_{1}$ & K2O5S2 & K2O5S2 &
      ↪  I-C. Chen and Y. Wang, Acta Crystallogr. C 40, 1780-1781 (1984
      ↪ )
   1.00000000000000
   6.92100000000000   0.00000000000000   0.00000000000000
   0.00000000000000   6.16000000000000   0.00000000000000
  -1.66852823637215   0.00000000000000   7.34999200846020
     K     O     S
     4    10     4
Direct
   0.23100000000000   0.25000000000000  -0.06470000000000   K   (2e)
  -0.23100000000000   0.75000000000000   0.06470000000000   K   (2e)
   0.63980000000000   0.25000000000000   0.67278000000000   K   (2e)
  -0.63980000000000   0.75000000000000  -0.67278000000000   K   (2e)
   0.65550000000000   0.25000000000000   0.04070000000000   O   (2e)
  -0.65550000000000   0.75000000000000  -0.04070000000000   O   (2e)
   0.07410000000000   0.05150000000000   0.23430000000000   O   (4f)
  -0.07410000000000   0.55150000000000  -0.23430000000000   O   (4f)
  -0.07410000000000  -0.05150000000000  -0.23430000000000   O   (4f)
   0.07410000000000   0.44850000000000   0.23430000000000   O   (4f)
   0.63500000000000   0.05300000000000   0.31470000000000   O   (4f)
  -0.63500000000000   0.55300000000000  -0.31470000000000   O   (4f)
  -0.63500000000000  -0.05300000000000  -0.31470000000000   O   (4f)
   0.63500000000000   0.44700000000000   0.31470000000000   O   (4f)
   0.02910000000000   0.25000000000000   0.33000000000000   S   (2e)
  -0.02910000000000   0.75000000000000  -0.33000000000000   S   (2e)
   0.70140000000000   0.25000000000000   0.23840000000000   S   (2e)
  -0.70140000000000   0.75000000000000  -0.23840000000000   S   (2e)
```

$ZrSe_3$: A3B_mP8_11_3e_e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Se3Zr'
_chemical_formula_sum 'Se3 Zr'

loop_
_publ_author_name
 'S. Furuseth'
 'L. Bratt{\aa}s'
 'A. Kjekshus'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 29a
_journal_year 1975
_journal_page_first 623
_journal_page_last 631
_publ_Section_title
;
 On the Crystal Structures of TiS$_{3}$, ZrS$_{3}$, ZrSe$_{3}$, ZrTe$_{3
      ↪ }$, HfS$_{3}$, and HfSe$_{3}$
;

_aflow_title 'ZrSe$_{3}$ Structure'
_aflow_proto 'A3B_mP8_11_3e_e'
_aflow_params 'a,b/a,c/a,\beta,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3},x_{4
      ↪ },z_{4}'
_aflow_params_values '5.4109,0.692823744664,1.74536583563,97.48,0.762,
      ↪ 0.554,0.456,0.174,0.888,0.169,0.285,0.656'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mP8'

_symmetry_space_group_name_H-M "P 1 21/m 1"
_symmetry_Int_Tables_number 11

_cell_length_a     5.41090
_cell_length_b     3.74880
_cell_length_c     9.44400
_cell_angle_alpha  90.00000
_cell_angle_beta   97.48000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z
3 -x,-y,-z
4 x,-y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Se1 Se 2 e 0.76200 0.25000 0.55400 1.00000
Se2 Se 2 e 0.45600 0.25000 0.17400 1.00000
Se3 Se 2 e 0.88800 0.25000 0.16900 1.00000
Zr1 Zr 2 e 0.28500 0.25000 0.65600 1.00000
```

$ZrSe_3$: A3B_mP8_11_3e_e - POSCAR

```
A3B_mP8_11_3e_e & a,b/a,c/a,beta,x1,z1,x2,z2,x3,z3,x4,z4 --params=5.4109
      ↪ ,0.692823744664,1.74536583563,97.48,0.762,0.554,0.456,0.174,
      ↪ 0.888,0.169,0.285,0.656 & P2_{1}/m C_{2h}^{2} #11 (e^4) & mP8 &
      ↪  None & Se3Zr & Se3Zr & S. Furuseth and L. Bratt{\aa}s and A.
      ↪ Kjekshus, Acta Chem. Scand. 29a, 623-631 (1975)
   1.00000000000000
   5.41090000000000   0.00000000000000   0.00000000000000
   0.00000000000000   3.74880000000000   0.00000000000000
  -1.22942090908250   0.00000000000000   9.36363499012594
     Se    Zr
      6     2
Direct
   0.76200000000000   0.25000000000000   0.55400000000000   Se   (2e)
  -0.76200000000000   0.75000000000000  -0.55400000000000   Se   (2e)
   0.45600000000000   0.25000000000000   0.17400000000000   Se   (2e)
  -0.45600000000000   0.75000000000000  -0.17400000000000   Se   (2e)
   0.88800000000000   0.25000000000000   0.16900000000000   Se   (2e)
  -0.88800000000000   0.75000000000000  -0.16900000000000   Se   (2e)
   0.28500000000000   0.25000000000000   0.65600000000000   Zr   (2e)
  -0.28500000000000   0.75000000000000  -0.65600000000000   Zr   (2e)
```

γ-$Y_2Si_2O_7$: A7B2C2_mP22_11_3e2f_2e_ab - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'O7Si2Y2'
_chemical_formula_sum 'O7 Si2 Y2'

loop_
_publ_author_name
 'N. G. Batalieva'
 'Y. A. Pyatenko'
_journal_name_full_name
;
 Soviet Physics Crystallography
;
_journal_volume 16
_journal_year 1972
_journal_page_first 786
_journal_page_last 789
_publ_Section_title
;
 Artificial yttrialite (\'\'y-phase\'\') - a representative of a new
      ↪ structure type in the rare earth diorthosilicate series
;

# Found in Revision of the crystallographic data of polymorphic Y$_{2}
      ↪ $Si$_{2}$O$_{7}$ and Y$_{2}$SiO$_{5}$ compounds, 2004

_aflow_title '$y$-Y$_{2}$Si$_{2}$O$_{7}$ Structure'
_aflow_proto 'A7B2C2_mP22_11_3e2f_2e_ab'
_aflow_params 'a,b/a,c/a,\beta,x_{3},z_{3},x_{4},z_{4},x_{5},z_{5},x_{6
      ↪ },z_{6},x_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9}'
_aflow_params_values '7.5,1.07466666667,0.669333333333,112.0,0.19,0.31,
      ↪ 0.508,0.264,0.88,0.43,0.12,0.588,0.709,0.548,0.19,0.09,0.8,
      ↪ 0.688,0.09,0.736'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mP22'

_symmetry_space_group_name_H-M "P 1 21/m 1"
_symmetry_Int_Tables_number 11

_cell_length_a    7.50
_cell_length_b    8.06000
_cell_length_c    5.02
_cell_angle_alpha 90.00000
_cell_angle_beta  112.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z
3 -x,-y,-z
4 x,-y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Y1 Y 2 a 0.00000 0.00000 0.00000 1.00000
Y2 Y 2 b 0.50000 0.00000 0.00000 1.00000
O1 O 2 e 0.19000 0.25000 0.31000 1.00000
O2 O 2 e 0.50800 0.25000 0.26400 1.00000
O3 O 2 e 0.88000 0.25000 0.43000 1.00000
Si1 Si 2 e 0.12000 0.25000 0.58800 1.00000
Si2 Si 2 e 0.70900 0.25000 0.54800 1.00000
O4 O 4 f 0.19000 0.09000 0.80000 1.00000
O5 O 4 f 0.68800 0.09000 0.73600 1.00000
```

y-$Y_2Si_2O_7$: A7B2C2_mP22_11_3e2f_2e_ab - POSCAR

```
A7B2C2_mP22_11_3e2f_2e_ab & a,b/a,c/a,beta,x3,z3,x4,z4,x5,z5,x6,z6,x7,z7
      ↪ ,x8,y8,z8,x9,y9,z9 --params=7.5,1.07466666667,0.669333333333,
      ↪ 112.0,0.19,0.31,0.508,0.264,0.88,0.43,0.12,0.588,0.709,0.548,
      ↪ 0.19,0.09,0.8,0.688,0.09,0.736 & P2_{1}/m C_{2h}^{2} #11 (abe^
      ↪ 5f^2) & mP22 & None & O7Si2Y2 & O7Si2Y2 & N. G. Batalieva and
      ↪ Y. A. Pyatenko, Sov. Phys. Crystallogr. 16, 786-789 (1972)
   1.00000000000000
   7.50000000000000    0.00000000000000    0.00000000000000
   0.00000000000000    8.06000000000000    0.00000000000000
  -1.88052509894788    0.00000000000000    4.65446294992527
     O     Si     Y
    14      4      4
Direct
   0.19000000000000    0.25000000000000    0.31000000000000    O    (2e)
  -0.19000000000000    0.75000000000000   -0.31000000000000    O    (2e)
   0.50800000000000    0.25000000000000    0.26400000000000    O    (2e)
  -0.50800000000000    0.75000000000000   -0.26400000000000    O    (2e)
   0.88000000000000    0.25000000000000    0.43000000000000    O    (2e)
  -0.88000000000000    0.75000000000000   -0.43000000000000    O    (2e)
   0.19000000000000    0.09000000000000    0.80000000000000    O    (4f)
  -0.19000000000000    0.59000000000000   -0.80000000000000    O    (4f)
  -0.19000000000000   -0.09000000000000   -0.80000000000000    O    (4f)
   0.19000000000000    0.41000000000000    0.80000000000000    O    (4f)
   0.68800000000000    0.09000000000000    0.73600000000000    O    (4f)
  -0.68800000000000    0.59000000000000   -0.73600000000000    O    (4f)
  -0.68800000000000   -0.09000000000000   -0.73600000000000    O    (4f)
   0.68800000000000    0.41000000000000    0.73600000000000    O    (4f)
   0.12000000000000    0.25000000000000    0.58800000000000   Si    (2e)
  -0.12000000000000    0.75000000000000   -0.58800000000000   Si    (2e)
   0.70900000000000    0.25000000000000    0.54800000000000   Si    (2e)
  -0.70900000000000    0.75000000000000   -0.54800000000000   Si    (2e)
   0.00000000000000    0.00000000000000    0.00000000000000    Y    (2a)
   0.00000000000000    0.50000000000000    0.00000000000000    Y    (2a)
   0.50000000000000    0.00000000000000    0.00000000000000    Y    (2b)
   0.50000000000000    0.50000000000000    0.00000000000000    Y    (2b)
```

Barytocalcite ($BaCa(CO_3)_2$): AB2CD6_mP20_11_e_2e_e_2e2f - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Barytocalcite'
_chemical_formula_sum 'Ba C2 Ca O6'

loop_
_publ_author_name
 'B. Dickens'
 'J. S. Bowen'
_journal_name_full_name
;
 Journal of Research of the National Bureau of Standards, Section A:
      ↪ Physics and Chemistry
;
_journal_volume 75
_journal_year 1971
_journal_page_first 197
_journal_page_last 203
_publ_Section_title
;
 The Crystal Structure of BaCa(CO$_{3}$)$_{2}$ (barytocalcite)
;

# Found in A new BaCa(CO$_{3}$)$_{2}$ polymorph, 2019

_aflow_title 'Barytocalcite (BaCa(CO$_{3}$)$_{2}$) Structure'
_aflow_proto 'AB2CD6_mP20_11_e_2e_e_2e2f'
_aflow_params 'a,b/a,c/a,\beta,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3},x_{4
      ↪ },z_{4},x_{5},z_{5},x_{6},z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},
      ↪ z_{8}'
_aflow_params_values '5.2344,1.54592694483,1.25019104386,106.05,0.1474,
      ↪ 0.28824,0.1028,0.7517,0.6149,0.7468,0.6232,0.19855,-0.0057,
      ↪ 0.8607,0.6383,0.5644,0.8457,-0.038,0.3089,0.6066,0.4604,0.8474'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mP20'

_symmetry_space_group_name_H-M "P 1 21/m 1"
_symmetry_Int_Tables_number 11

_cell_length_a    5.23440
_cell_length_b    8.09200
_cell_length_c    6.54400
_cell_angle_alpha 90.00000
_cell_angle_beta  106.05000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z
3 -x,-y,-z
4 x,-y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ba1 Ba 2 e 0.14740 0.25000 0.28824 1.00000
C1 C 2 e 0.10280 0.25000 0.75170 1.00000
C2 C 2 e 0.61490 0.25000 0.74680 1.00000
Ca1 Ca 2 e 0.62320 0.25000 0.19855 1.00000
O1 O 2 e -0.00570 0.25000 0.86070 1.00000
O2 O 2 e 0.63830 0.25000 0.56440 1.00000
O3 O 4 f 0.84570 -0.03800 0.30890 1.00000
O4 O 4 f 0.60660 0.46040 0.84740 1.00000
```

Barytocalcite ($BaCa(CO_3)_2$): AB2CD6_mP20_11_e_2e_e_2e2f - POSCAR

```
AB2CD6_mP20_11_e_2e_e_2e2f & a,b/a,c/a,beta,x1,z1,x2,z2,x3,z3,x4,z4,x5,
      ↪ z5,x6,z6,x7,y7,z7,x8,y8,z8 --params=5.2344,1.54592694483,
      ↪ 1.25019104386,106.05,0.1474,0.28824,0.1028,0.7517,0.6149,0.7468
      ↪ ,0.6232,0.19855,-0.0057,0.8607,0.6383,0.5644,0.8457,-0.038,
      ↪ 0.3089,0.6066,0.4604,0.8474 & P2_{1}/m C_{2h}^{2} #11 (e^6f^2)
      ↪ & mP20 & None & BaC2CaO6 & Barytocalcite & B. Dickens and J. S.
      ↪ Bowen, J. Res. Nat. Stand. Sec. A 75, 197-203 (1971)
   1.00000000000000
   5.23440000000000    0.00000000000000    0.00000000000000
   0.00000000000000    8.09200000000000    0.00000000000000
  -1.80925966275357    0.00000000000000    6.28892005615693
     Ba     C     Ca     O
      2     4      2     12
Direct
   0.14740000000000    0.25000000000000    0.28824000000000   Ba    (2e)
  -0.14740000000000    0.75000000000000   -0.28824000000000   Ba    (2e)
   0.10280000000000    0.25000000000000    0.75170000000000    C    (2e)
  -0.10280000000000    0.75000000000000   -0.75170000000000    C    (2e)
   0.61490000000000    0.25000000000000    0.74680000000000    C    (2e)
```

```
  −0.61490000000000    0.75000000000000   −0.74680000000000    C    (2e)
   0.62320000000000    0.25000000000000    0.19855000000000    Ca   (2e)
  −0.62320000000000    0.75000000000000   −0.19855000000000    Ca   (2e)
  −0.00570000000000    0.25000000000000    0.86070000000000    O    (2e)
   0.00570000000000    0.75000000000000   −0.86070000000000    O    (2e)
   0.63830000000000    0.25000000000000    0.56440000000000    O    (2e)
  −0.63830000000000    0.75000000000000   −0.56440000000000    O    (2e)
   0.84570000000000   −0.03800000000000    0.30890000000000    O    (4f)
  −0.84570000000000    0.46200000000000   −0.30890000000000    O    (4f)
  −0.84570000000000    0.03800000000000   −0.30890000000000    O    (4f)
   0.84570000000000    0.53800000000000    0.30890000000000    O    (4f)
   0.60660000000000    0.46040000000000    0.84740000000000    O    (4f)
  −0.60660000000000    0.96040000000000   −0.84740000000000    O    (4f)
  −0.60660000000000   −0.46040000000000   −0.84740000000000    O    (4f)
   0.60660000000000    0.03960000000000    0.84740000000000    O    (4f)
```

O(OH)Y: ABC_mP6_11_e_e_e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'HO2Y'
_chemical_formula_sum 'O (OH) Y'

loop_
_publ_author_name
 'R. F. Klevtsova'
 'P. V. Klevtsov'
_journal_year 1965
_publ_Section_title
;
 The Crystal Structure of YOOH
;

_aflow_title 'O(OH)Y Structure'
_aflow_proto 'ABC_mP6_11_e_e_e'
_aflow_params 'a,b/a,c/a,\beta,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3}'
_aflow_params_values '4.28,0.848130841121,1.41355140187,112.5,0.219,
      ↪ 0.448,0.69,-0.058,0.643,0.312'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mP6'

_symmetry_space_group_name_H-M "P 1 21/m 1"
_symmetry_Int_Tables_number 11

_cell_length_a     4.28000
_cell_length_b     3.63000
_cell_length_c     6.05000
_cell_angle_alpha  90.00000
_cell_angle_beta   112.50000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z
3 -x,-y,-z
4 x,-y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 2 e 0.21900 0.25000 0.44800 1.00000
OH1 OH 2 e 0.69000 0.25000 -0.05800 1.00000
Y1 Y 2 e 0.64300 0.25000 0.31200 1.00000
```

O(OH)Y: ABC_mP6_11_e_e_e - POSCAR

```
ABC_mP6_11_e_e_e & a,b/a,c/a,beta,x1,z1,x2,z2,x3,z3 --params=4.28,
      ↪ 0.848130841121,1.41355140187,112.5,0.219,0.448,0.69,-0.058,
      ↪ 0.643,0.312 & P2_{1}/m C_{2h}^{2} #11 (e^3) & mP6 & None & HO2Y
      ↪ & HO2Y & R. F. Klevtsova and P. V. Klevtsov, (1965)
   1.00000000000000
   4.28000000000000    0.00000000000000    0.00000000000000
   0.00000000000000    3.63000000000000    0.00000000000000
  -2.31523476580879    0.00000000000000    5.58947117169329
     O     OH      Y
     2      2      2
Direct
   0.21900000000000    0.25000000000000    0.44800000000000    O    (2e)
  -0.21900000000000    0.75000000000000   -0.44800000000000    O    (2e)
   0.69000000000000    0.25000000000000   -0.05800000000000    OH   (2e)
  -0.69000000000000    0.75000000000000    0.05800000000000    OH   (2e)
   0.64300000000000    0.25000000000000    0.31200000000000    Y    (2e)
  -0.64300000000000    0.75000000000000   -0.31200000000000    Y    (2e)
```

$Al_{13}Fe_4$: A13B4_mC102_12_dg8i5j_4ij - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Al13Fe4'
_chemical_formula_sum 'Al13 Fe4'

loop_
_publ_author_name
 'P. J. Black'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 8
_journal_year 1955
_journal_page_first 43
_journal_page_last 48
_publ_Section_title
;
 The Structure of FeAl$_{3}$. I
;

_aflow_title 'Al$_{13}$Fe$_{4}$ Structure'
_aflow_proto 'A13B4_mC102_12_dg8i5j_4ij'
_aflow_params 'a,b/a,c/a,\beta,y_{2},x_{3},z_{3},x_{4},z_{4},x_{5},z_{5
      ↪ },x_{6},z_{6},x_{7},z_{7},x_{8},z_{8},x_{9},z_{9},x_{10},z_{10
      ↪ },x_{11},z_{11},x_{12},z_{12},x_{13},z_{13},x_{14},z_{14},x_{15
      ↪ },y_{15},z_{15},x_{16},y_{16},z_{16},x_{17},y_{17},z_{17},x_{18
      ↪ },y_{18},z_{18},x_{19},y_{19},z_{19},x_{20},y_{20},z_{20}'
_aflow_params_values '15.489,0.521860675318,0.805474853122,107.71667,
      ↪ 0.244,0.064,0.173,0.322,0.277,0.235,0.539,0.081,0.582,0.231,-
      ↪ 0.028,0.48,0.827,0.31,0.769,0.086,0.781,0.086,0.383,0.401,0.624
      ↪ ,0.09,-0.011,0.4,-0.015,0.188,0.216,0.111,0.373,0.211,0.107,
      ↪ 0.176,0.216,0.334,0.495,0.283,0.329,0.366,0.223,0.479,0.318,
      ↪ 0.285,0.277'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mC102'

_symmetry_space_group_name_H-M "C 1 2/m 1"
_symmetry_Int_Tables_number 12

_cell_length_a     15.48900
_cell_length_b     8.08310
_cell_length_c     12.47600
_cell_angle_alpha  90.00000
_cell_angle_beta   107.71667
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z
3 -x,-y,-z
4 x,-y,z
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Al1 Al 2 d 0.00000 0.50000 0.50000 1.00000
Al2 Al 4 g 0.00000 0.24400 0.00000 1.00000
Al3 Al 4 i 0.06400 0.00000 0.17300 1.00000
Al4 Al 4 i 0.32200 0.00000 0.27700 0.70000
Al5 Al 4 i 0.23500 0.00000 0.53900 1.00000
Al6 Al 4 i 0.08100 0.00000 0.58200 1.00000
Al7 Al 4 i 0.23100 0.00000 -0.02800 1.00000
Al8 Al 4 i 0.48000 0.00000 0.82700 1.00000
Al9 Al 4 i 0.31000 0.00000 0.76900 1.00000
Al10 Al 4 i 0.08600 0.00000 0.78100 1.00000
Fe1 Fe 4 i 0.08600 0.00000 0.38300 1.00000
Fe2 Fe 4 i 0.40100 0.00000 0.62400 1.00000
Fe3 Fe 4 i 0.09000 0.00000 -0.01100 1.00000
Fe4 Fe 4 i 0.40000 0.00000 -0.01500 1.00000
Al11 Al 8 j 0.18800 0.21600 0.11100 1.00000
Al12 Al 8 j 0.37300 0.21100 0.10700 1.00000
Al13 Al 8 j 0.17600 0.21600 0.33400 1.00000
Al14 Al 8 j 0.49500 0.28300 0.32900 1.00000
Al15 Al 8 j 0.36600 0.22300 0.47900 1.00000
Fe5 Fe 8 j 0.31800 0.28500 0.27700 1.00000
```

$Al_{13}Fe_4$: A13B4_mC102_12_dg8i5j_4ij - POSCAR

```
A13B4_mC102_12_dg8i5j_4ij & a,b/a,c/a,beta,y2,x3,z3,x4,z4,x5,z5,x6,z6,x7
      ↪ ,z7,x8,z8,x9,z9,x10,z10,x11,z11,x12,z12,x13,z13,x14,z14,x15,y15
      ↪ ,z15,x16,y16,z16,x17,y17,z17,x18,y18,z18,x19,y19,z19,x20,y20,
      ↪ z20 --params=15.489,0.521860675318,0.805474853122,107.71667,
      ↪ 0.244,0.064,0.173,0.322,0.277,0.235,0.539,0.081,0.582,0.231,-
      ↪ 0.028,0.48,0.827,0.31,0.769,0.086,0.781,0.086,0.383,0.401,0.624
      ↪ ,0.09,-0.011,0.4,-0.015,0.188,0.216,0.111,0.373,0.211,0.107,
      ↪ 0.176,0.216,0.334,0.495,0.283,0.329,0.366,0.223,0.479,0.318,
      ↪ 0.285,0.277 & C2/m C_{2h}^{3} #12 (dgi^12j^6) & mC102 & None &
      ↪ Al13Fe4 & Al13Fe4 & P. J. Black, Acta Cryst. 8, 43-48 (1955)
   1.00000000000000
   7.74450000000000   -4.04155000000000    0.00000000000000
   7.74450000000000    4.04155000000000    0.00000000000000
  -3.79657432304420    0.00000000000000   11.88430054355750
     Al     Fe
     39     12
Direct
   0.50000000000000    0.50000000000000    0.50000000000000    Al   (2d)
  -0.24400000000000    0.24400000000000    0.00000000000000    Al   (4g)
   0.24400000000000   -0.24400000000000    0.00000000000000    Al   (4g)
```

```
  0.06400000000000   0.06400000000000   0.17300000000000   Al   (4i)
 -0.06400000000000  -0.06400000000000  -0.17300000000000   Al   (4i)
  0.32200000000000   0.32200000000000   0.27700000000000   Al   (4i)
 -0.32200000000000  -0.32200000000000  -0.27700000000000   Al   (4i)
  0.23500000000000   0.23500000000000   0.53900000000000   Al   (4i)
 -0.23500000000000  -0.23500000000000  -0.53900000000000   Al   (4i)
  0.08100000000000   0.08100000000000   0.58200000000000   Al   (4i)
 -0.08100000000000  -0.08100000000000  -0.58200000000000   Al   (4i)
  0.23100000000000   0.23100000000000  -0.02800000000000   Al   (4i)
 -0.23100000000000  -0.23100000000000   0.02800000000000   Al   (4i)
  0.48000000000000   0.48000000000000   0.82700000000000   Al   (4i)
 -0.48000000000000  -0.48000000000000  -0.82700000000000   Al   (4i)
  0.31000000000000   0.31000000000000   0.76900000000000   Al   (4i)
 -0.31000000000000  -0.31000000000000  -0.76900000000000   Al   (4i)
  0.08600000000000   0.08600000000000   0.78100000000000   Al   (4i)
 -0.08600000000000  -0.08600000000000  -0.78100000000000   Al   (4i)
 -0.02800000000000   0.40400000000000   0.11100000000000   Al   (8j)
 -0.40400000000000   0.02800000000000  -0.11100000000000   Al   (8j)
  0.02800000000000  -0.40400000000000  -0.11100000000000   Al   (8j)
  0.40400000000000  -0.02800000000000   0.11100000000000   Al   (8j)
  0.16200000000000   0.58400000000000   0.10700000000000   Al   (8j)
 -0.58400000000000  -0.16200000000000  -0.10700000000000   Al   (8j)
 -0.16200000000000  -0.58400000000000  -0.10700000000000   Al   (8j)
  0.58400000000000   0.16200000000000   0.10700000000000   Al   (8j)
 -0.04000000000000   0.39200000000000   0.33400000000000   Al   (8j)
 -0.39200000000000   0.04000000000000  -0.33400000000000   Al   (8j)
  0.04000000000000  -0.39200000000000  -0.33400000000000   Al   (8j)
  0.39200000000000  -0.04000000000000   0.33400000000000   Al   (8j)
  0.21200000000000   0.77800000000000   0.32900000000000   Al   (8j)
 -0.77800000000000  -0.21200000000000  -0.32900000000000   Al   (8j)
 -0.21200000000000  -0.77800000000000  -0.32900000000000   Al   (8j)
  0.77800000000000   0.21200000000000   0.32900000000000   Al   (8j)
  0.14300000000000   0.58900000000000   0.47900000000000   Al   (8j)
 -0.58900000000000  -0.14300000000000  -0.47900000000000   Al   (8j)
 -0.14300000000000  -0.58900000000000  -0.47900000000000   Al   (8j)
  0.58900000000000   0.14300000000000   0.47900000000000   Al   (8j)
  0.08600000000000   0.08600000000000   0.38300000000000   Fe   (4i)
 -0.08600000000000  -0.08600000000000  -0.38300000000000   Fe   (4i)
  0.40100000000000   0.40100000000000   0.62400000000000   Fe   (4i)
 -0.40100000000000  -0.40100000000000  -0.62400000000000   Fe   (4i)
  0.09000000000000   0.09000000000000  -0.01100000000000   Fe   (4i)
 -0.09000000000000  -0.09000000000000   0.01100000000000   Fe   (4i)
  0.40000000000000   0.40000000000000  -0.01500000000000   Fe   (4i)
 -0.40000000000000  -0.40000000000000   0.01500000000000   Fe   (4i)
  0.03300000000000   0.60300000000000   0.27700000000000   Fe   (8j)
 -0.60300000000000  -0.03300000000000  -0.27700000000000   Fe   (8j)
 -0.03300000000000  -0.60300000000000  -0.27700000000000   Fe   (8j)
  0.60300000000000   0.03300000000000   0.27700000000000   Fe   (8j)
```

Os_4Al_{13}: A13B4_mC34_12_b6i_2i - CIF

```
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_audit_creation_method FINDSYM

_chemical_name_mineral 'Al13Os4'
_chemical_formula_sum 'Al13 Os4'

loop_
_publ_author_name
 'L.-E. Edshammar'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 18
_journal_year 1964
_journal_page_first 2294
_journal_page_last 2302
_publ_Section_title
;
 The Crystal Structure of Os$_{4}$Al$_{13}$
;

_aflow_title 'Os$_{4}$Al$_{13}$ Structure'
_aflow_proto 'A13B4_mC34_12_b6i_2i'
_aflow_params 'a,b/a,c/a,\beta,x_{2},z_{2},x_{3},z_{3},x_{4},z_{4},x_{5
      ↪ },z_{5},x_{6},z_{6},x_{7},z_{7},x_{8},z_{8},x_{9},z_{9}'
_aflow_params_values '17.64,0.239682539683,0.440646258503,115.15,0.587,
      ↪ 0.368,0.257,0.613,0.132,0.162,0.79,0.087,-0.086,0.432,0.409,
      ↪ 0.194,0.294,0.2915,-0.0081,0.1947'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mC34'

_symmetry_space_group_name_H-M "C 1 2/m 1"
_symmetry_Int_Tables_number 12

_cell_length_a     17.64000
_cell_length_b     4.22800
_cell_length_c     7.77300
_cell_angle_alpha  90.00000
_cell_angle_beta   115.15000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
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2 -x,y,-z
3 -x,-y,-z
4 x,-y,z
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Al1  Al  2  b  0.00000  0.50000  0.00000  1.00000
Al2  Al  4  i  0.58700  0.00000  0.36800  1.00000
Al3  Al  4  i  0.25700  0.00000  0.61300  1.00000
Al4  Al  4  i  0.13200  0.00000  0.16200  1.00000
Al5  Al  4  i  0.79000  0.00000  0.08700  1.00000
Al6  Al  4  i -0.08600  0.00000  0.43200  1.00000
Al7  Al  4  i  0.40900  0.00000  0.19400  1.00000
Os1  Os  4  i  0.29400  0.00000  0.29150  1.00000
Os2  Os  4  i -0.00810  0.00000  0.19470  1.00000
```

Os_4Al_{13}: A13B4_mC34_12_b6i_2i - POSCAR

```
A13B4_mC34_12_b6i_2i & a,b/a,c/a,beta,x2,z2,x3,z3,x4,z4,x5,z5,x6,z6,x7,
      ↪ z7,x8,z8,x9,z9 --params=17.64,0.239682539683,0.440646258503,
      ↪ 115.15,0.587,0.368,0.257,0.613,0.132,0.162,0.79,0.087,-0.086,
      ↪ 0.432,0.409,0.194,0.294,0.2915,-0.0081,0.1947 & C2/m C_{2h}^{3}
      ↪  #12 (bi^8) & mC34 & None & Al13Os4 & Al13Os4 & L.-E. Edshammar
      ↪ , Acta Chem. Scand. 18, 2294-2302 (1964)
   1.00000000000000
   8.82000000000000  -2.11400000000000   0.00000000000000
   8.82000000000000   2.11400000000000   0.00000000000000
  -3.30344353102559   0.00000000000000   7.03610615591644
     Al      Os
     13       4
Direct
   0.50000000000000   0.50000000000000   0.00000000000000   Al   (2b)
   0.58700000000000   0.58700000000000   0.36800000000000   Al   (4i)
  -0.58700000000000  -0.58700000000000  -0.36800000000000   Al   (4i)
   0.25700000000000   0.25700000000000   0.61300000000000   Al   (4i)
  -0.25700000000000  -0.25700000000000  -0.61300000000000   Al   (4i)
   0.13200000000000   0.13200000000000   0.16200000000000   Al   (4i)
  -0.13200000000000  -0.13200000000000  -0.16200000000000   Al   (4i)
   0.79000000000000   0.79000000000000   0.08700000000000   Al   (4i)
  -0.79000000000000  -0.79000000000000  -0.08700000000000   Al   (4i)
  -0.08600000000000  -0.08600000000000   0.43200000000000   Al   (4i)
   0.08600000000000   0.08600000000000  -0.43200000000000   Al   (4i)
   0.40900000000000   0.40900000000000   0.19400000000000   Al   (4i)
  -0.40900000000000  -0.40900000000000  -0.19400000000000   Al   (4i)
   0.29400000000000   0.29400000000000   0.29150000000000   Os   (4i)
  -0.29400000000000  -0.29400000000000  -0.29150000000000   Os   (4i)
  -0.00810000000000  -0.00810000000000   0.19470000000000   Os   (4i)
   0.00810000000000   0.00810000000000  -0.19470000000000   Os   (4i)
```

Bischofite ($MgCl_2{\cdot}6H_2O$, $J1_7$): A2B12CD6_mC42_12_i_2i2j_a_ij - CIF

```
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_chemical_name_mineral 'Bischofite'
_chemical_formula_sum 'Cl2 H12 Mg O6'

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_publ_author_name
 'P. A. Agron'
 'W. R. Busing'
_journal_name_full_name
;
 Acta Crystallographica Section C: Structural Chemistry
;
_journal_volume 41
_journal_year 1985
_journal_page_first 8
_journal_page_last 10
_publ_Section_title
;
 Magnesium dichloride hexahydrate, MgCl$_{2}$$\cdot$6H$_{2}$O, by
      ↪ neutron diffraction
;

_aflow_title 'Bischofite (MgCl$_{2}$$\cdot$6H$_{2}$O, $J1_{7}$)
      ↪ Structure'
_aflow_proto 'A2B12CD6_mC42_12_i_2i2j_a_ij'
_aflow_params 'a,b/a,c/a,\beta,x_{2},z_{2},x_{3},z_{3},x_{4},z_{4},x_{5
      ↪ },z_{5},x_{6},y_{6},z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},z_{8}'
_aflow_params_values '9.8607,0.720750048171,0.615950186092,93.758,0.3176
      ↪ ,0.6122,0.2372,0.2583,0.2693,0.0083,0.2019,0.1095,0.0209,0.2997
      ↪ ,0.2784,0.8839,0.1984,0.3151,-0.0429,0.2067,0.2233'
_aflow_Strukturbericht '$J1_{7}$'
_aflow_Pearson 'mC42'

_symmetry_space_group_name_H-M "C 1 2/m 1"
_symmetry_Int_Tables_number 12

_cell_length_a     9.86070
_cell_length_b     7.10710
_cell_length_c     6.07370
_cell_angle_alpha  90.00000
_cell_angle_beta   93.75800
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
```

```
2 -x,y,-z
3 -x,-y,-z
4 x,-y,z
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Mg1 Mg 2 a 0.00000 0.00000 0.00000 1.00000
Cl1 Cl 4 i 0.31760 0.00000 0.61220 1.00000
H1 H 4 i 0.23720 0.00000 0.25830 1.00000
H2 H 4 i 0.26930 0.00000 0.00830 1.00000
O1 O 4 i 0.20190 0.00000 0.10950 1.00000
H3 H 8 j 0.02090 0.29970 0.27840 1.00000
H4 H 8 j 0.88390 0.19840 0.31510 1.00000
O2 O 8 j -0.04290 0.20670 0.22330 1.00000
```

Bischofite ($MgCl_2{\cdot}6H_2O$, $J1_7$): A2B12CD6_mC42_12_i_2i2j_a_ij - POSCAR

```
A2B12CD6_mC42_12_i_2i2j_a_ij & a,b/a,c/a,beta,x2,z2,x3,z3,x4,z4,x5,z5,x6
    ↪ ,y6,z6,x7,y7,z7,x8,y8,z8 --params=9.8607,0.720750048171,
    ↪ 0.615950186092,93.758,0.3176,0.6122,0.2372,0.2583,0.2693,0.0083
    ↪ ,0.2019,0.1095,0.0209,0.2997,0.2784,0.8839,0.1984,0.3151,-
    ↪ 0.0429,0.2067,0.2233 & C2/m C_{2h}^{3} #12 (ai^4j^3) & mC42 &
    ↪ $J1_{7}$ & Cl2H12MgO6 & Bischofite & P. A. Agron and W. R.
    ↪ Busing, Acta Crystallogr. C 41, 8-10 (1985)
  1.00000000000000
  4.93035000000000  -3.55355000000000   0.00000000000000
  4.93035000000000   3.55355000000000   0.00000000000000
 -0.39808521489547   0.00000000000000   6.06064021797051
    Cl   H  Mg   O
     2  12   1   6
Direct
  0.31760000000000   0.31760000000000   0.61220000000000  Cl  (4i)
 -0.31760000000000  -0.31760000000000  -0.61220000000000  Cl  (4i)
  0.23720000000000   0.23720000000000   0.25830000000000  H  (4i)
 -0.23720000000000  -0.23720000000000  -0.25830000000000  H  (4i)
  0.26930000000000   0.26930000000000   0.00830000000000  H  (4i)
 -0.26930000000000  -0.26930000000000  -0.00830000000000  H  (4i)
 -0.27880000000000   0.32060000000000   0.27840000000000  H  (8j)
 -0.32060000000000   0.27880000000000  -0.27840000000000  H  (8j)
  0.27880000000000  -0.32060000000000  -0.27840000000000  H  (8j)
  0.32060000000000  -0.27880000000000   0.27840000000000  H  (8j)
  0.68550000000000   1.08230000000000   0.31510000000000  H  (8j)
 -1.08230000000000  -0.68550000000000  -0.31510000000000  H  (8j)
 -0.68550000000000  -1.08230000000000  -0.31510000000000  H  (8j)
  1.08230000000000   0.68550000000000   0.31510000000000  H  (8j)
  0.00000000000000   0.00000000000000   0.00000000000000  Mg  (2a)
  0.20190000000000   0.20190000000000   0.10950000000000  O  (4i)
 -0.20190000000000  -0.20190000000000  -0.10950000000000  O  (4i)
 -0.24960000000000   0.16380000000000   0.22330000000000  O  (8j)
 -0.16380000000000   0.24960000000000  -0.22330000000000  O  (8j)
  0.24960000000000  -0.16380000000000  -0.22330000000000  O  (8j)
  0.16380000000000  -0.24960000000000   0.22330000000000  O  (8j)
```

Tremolite ($Ca_2Mg_5Si_8O_{22}(OH)_2$, $S4_2$): A2B2C5D24E8_mC82_12_h_i_agh_2i5j_2j - CIF

```
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_chemical_formula_sum 'Ca2 H2 Mg5 O24 Si8'

loop_
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 'M. Merli'
 'L. Ungaretti'
 'R. Oberti'
_journal_name_full_name
;
 American Mineralogist
;
_journal_volume 85
_journal_year 2000
_journal_page_first 532
_journal_page_last 542
_publ_Section_title
;
 Leverage analysis and structure refinement of minerals
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'Tremolite (Ca${_}2${}Mg${_}5$Si${_}8${}O${_}22${}(OH)${_}2$,
    ↪ $S4_{2}$) Structure'
_aflow_proto 'A2B2C5D24E8_mC82_12_h_i_agh_2i5j_2j'
_aflow_params 'a,b/a,c/a,\beta,y_{2},y_{3},y_{4},x_{5},z_{5},x_{6},z_{6
    ↪ },x_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{10},y_{10
    ↪ },z_{10},x_{11},y_{11},z_{11},x_{12},y_{12},z_{12},x_{13},y_{13
    ↪ },z_{13},x_{14},y_{14},z_{14}'
_aflow_params_values '9.8359,1.83460588253,0.536321028071,104.75,0.1765,
    ↪ 0.278,0.0878,0.196,0.764,0.1085,0.7155,0.3377,0.2928,0.1119,
    ↪ 0.0857,0.218,0.1187,0.1709,0.7244,0.1351,0.2519,0.2069,0.3466,
    ↪ 0.1344,0.1005,0.344,0.1188,0.5891,0.2806,0.0839,0.2972,0.2884,
    ↪ 0.1711,0.8047'
_aflow_Strukturbericht '$S4_{2}$'
_aflow_Pearson 'mC82'

_symmetry_space_group_name_H-M "C 1 2/m 1"
_symmetry_Int_Tables_number 12

_cell_length_a      9.83590
_cell_length_b      18.04500
_cell_length_c      5.27520
_cell_angle_alpha   90.00000
_cell_angle_beta    104.75000
_cell_angle_gamma   90.00000

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_space_group_symop_id
_space_group_symop_operation_xyz
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2 -x,y,-z
3 -x,-y,-z
4 x,-y,z
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Mg1 Mg 2 a 0.00000 0.00000 0.00000 1.00000
Mg2 Mg 4 g 0.00000 0.17650 0.00000 1.00000
Ca1 Ca 4 h 0.00000 0.27800 0.50000 1.00000
Mg3 Mg 4 h 0.00000 0.08780 0.50000 1.00000
H1 H 4 i 0.19600 0.00000 0.76400 1.00000
O1 O 4 i 0.10850 0.00000 0.71550 1.00000
O2 O 4 i 0.33770 0.00000 0.29280 1.00000
O3 O 8 j 0.11190 0.08570 0.21800 1.00000
O4 O 8 j 0.11870 0.17090 0.72440 1.00000
O5 O 8 j 0.13510 0.25190 0.20690 1.00000
O6 O 8 j 0.34660 0.13440 0.10050 1.00000
O7 O 8 j 0.34400 0.11880 0.58910 1.00000
Si1 Si 8 j 0.28060 0.08390 0.29720 1.00000
Si2 Si 8 j 0.28840 0.17110 0.80470 1.00000
```

Tremolite ($Ca_2Mg_5Si_8O_{22}(OH)_2$, $S4_2$): A2B2C5D24E8_mC82_12_h_i_agh_2i5j_2j - POSCAR

```
A2B2C5D24E8_mC82_12_h_i_agh_2i5j_2j & a,b/a,c/a,beta,y2,y3,y4,x5,z5,x6,
    ↪ z6,x7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10,x11,y11,z11,x12,y12,z12,
    ↪ x13,y13,z13,x14,y14,z14 --params=9.8359,1.83460588253,
    ↪ 0.536321028071,104.75,0.1765,0.278,0.0878,0.196,0.764,0.1085,
    ↪ 0.7155,0.3377,0.2928,0.1119,0.0857,0.218,0.1187,0.1709,0.7244,
    ↪ 0.1351,0.2519,0.2069,0.3466,0.1344,0.1005,0.344,0.1188,0.5891,
    ↪ 0.2806,0.0839,0.2972,0.2884,0.1711,0.8047 & C2/m C_{2h}^{3} #12
    ↪ (agh^2i^3j^7) & mC82 & $S4_{2}$ & Ca2H2Mg5O24Si8 & Tremolite &
    ↪ M. Merli and L. Ungaretti and R. Oberti, Am. Mineral. 85,
    ↪ 532-542 (2000)
  1.00000000000000
  4.91795000000000  -9.02250000000000   0.00000000000000
  4.91795000000000   9.02250000000000   0.00000000000000
 -1.34307619717380   0.00000000000000   5.10136073695883
    Ca   H  Mg   O  Si
     2   2   5  24   8
Direct
 -0.27800000000000   0.27800000000000   0.50000000000000  Ca  (4h)
  0.27800000000000  -0.27800000000000   0.50000000000000  Ca  (4h)
  0.19600000000000   0.19600000000000   0.76400000000000  H  (4i)
 -0.19600000000000  -0.19600000000000  -0.76400000000000  H  (4i)
  0.00000000000000   0.00000000000000   0.00000000000000  Mg  (2a)
 -0.17650000000000   0.17650000000000   0.00000000000000  Mg  (4g)
  0.17650000000000  -0.17650000000000   0.00000000000000  Mg  (4g)
 -0.08780000000000   0.08780000000000   0.50000000000000  Mg  (4h)
  0.08780000000000  -0.08780000000000   0.50000000000000  Mg  (4h)
  0.10850000000000   0.10850000000000   0.71550000000000  O  (4i)
 -0.10850000000000  -0.10850000000000  -0.71550000000000  O  (4i)
  0.33770000000000   0.33770000000000   0.29280000000000  O  (4i)
 -0.33770000000000  -0.33770000000000  -0.29280000000000  O  (4i)
  0.02620000000000   0.19760000000000   0.21800000000000  O  (8j)
 -0.19760000000000  -0.02620000000000  -0.21800000000000  O  (8j)
 -0.02620000000000  -0.19760000000000  -0.21800000000000  O  (8j)
  0.19760000000000   0.02620000000000   0.21800000000000  O  (8j)
 -0.05220000000000   0.28960000000000   0.72440000000000  O  (8j)
 -0.28960000000000   0.05220000000000  -0.72440000000000  O  (8j)
  0.05220000000000  -0.28960000000000  -0.72440000000000  O  (8j)
  0.28960000000000  -0.05220000000000   0.72440000000000  O  (8j)
 -0.11680000000000   0.38700000000000   0.20690000000000  O  (8j)
 -0.38700000000000   0.11680000000000  -0.20690000000000  O  (8j)
  0.11680000000000  -0.38700000000000  -0.20690000000000  O  (8j)
  0.38700000000000  -0.11680000000000   0.20690000000000  O  (8j)
  0.21220000000000   0.48100000000000   0.10050000000000  O  (8j)
 -0.48100000000000  -0.21220000000000  -0.10050000000000  O  (8j)
 -0.21220000000000  -0.48100000000000  -0.10050000000000  O  (8j)
  0.48100000000000   0.21220000000000   0.10050000000000  O  (8j)
  0.22520000000000   0.46280000000000   0.58910000000000  O  (8j)
 -0.46280000000000  -0.22520000000000  -0.58910000000000  O  (8j)
 -0.22520000000000  -0.46280000000000  -0.58910000000000  O  (8j)
  0.46280000000000   0.22520000000000   0.58910000000000  O  (8j)
  0.19670000000000   0.36450000000000   0.29720000000000  Si  (8j)
 -0.36450000000000  -0.19670000000000  -0.29720000000000  Si  (8j)
 -0.19670000000000  -0.36450000000000  -0.29720000000000  Si  (8j)
  0.36450000000000   0.19670000000000   0.29720000000000  Si  (8j)
```

```
   0.11730000000000   0.45950000000000   0.80470000000000   Si   (8j)
  -0.45950000000000  -0.11730000000000  -0.80470000000000   Si   (8j)
  -0.11730000000000  -0.45950000000000  -0.80470000000000   Si   (8j)
   0.45950000000000   0.11730000000000   0.80470000000000   Si   (8j)
```

β-Ga_2O_3: A2B3_mC20_12_2i_3i - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Ga2O3'
_chemical_formula_sum 'Ga2 O3'

loop_
_publ_author_name
 'J. {\AA}hman'
 'G. Svensson'
 'J. Albertsson'
_journal_name_full_name
;
 Acta Crystallographica Section C: Structural Chemistry
;
_journal_volume 52
_journal_year 1996
_journal_page_first 1336
_journal_page_last 1338
_publ_Section_title
;
A Reinvestigation of $\beta$-Gallium Oxide
;

_aflow_title '$\beta$-Ga$_{2}$O$_{3}$ Structure'
_aflow_proto 'A2B3_mC20_12_2i_3i'
_aflow_params 'a,b/a,c/a,\beta,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3},x_{4
    ↪ },z_{4},x_{5},z_{5}'
_aflow_params_values '12.214,0.248657278533,0.474709349926,103.83,0.0905
    ↪ ,0.7946,0.65866,0.31402,0.1645,0.1098,0.1733,0.5632,0.4959,
    ↪ 0.2566'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mC20'

_symmetry_space_group_name_H-M "C 1 2/m 1"
_symmetry_Int_Tables_number 12

_cell_length_a     12.21400
_cell_length_b     3.03710
_cell_length_c     5.79810
_cell_angle_alpha  90.00000
_cell_angle_beta   103.83000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z
3 -x,-y,-z
4 x,-y,z
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ga1 Ga 4 i 0.09050 0.00000 0.79460 1.00000
Ga2 Ga 4 i 0.65866 0.00000 0.31402 1.00000
O1 O 4 i 0.16450 0.00000 0.10980 1.00000
O2 O 4 i 0.17330 0.00000 0.56320 1.00000
O3 O 4 i 0.49590 0.00000 0.25660 1.00000
```

β-Ga_2O_3: A2B3_mC20_12_2i_3i - POSCAR

```
A2B3_mC20_12_2i_3i & a,b/a,c/a,beta,x1,z1,x2,z2,x3,z3,x4,z4,x5,z5 --
    ↪ params=12.214,0.248657278533,0.474709349926,103.83,0.0905,
    ↪ 0.7946,0.65866,0.31402,0.1645,0.1098,0.1733,0.5632,0.4959,
    ↪ 0.2566 & C2/m C_{2h}^{3} #12 (i^5) & mC20 & None & Ga2O3 &
    ↪ Ga2O3 & J. {\AA}hman and G. Svensson and J. Albertsson, Acta
    ↪ Crystallogr. C 52, 1336-1338 (1996)
   1.00000000000000
   6.10700000000000  -1.51855000000000   0.00000000000000
   6.10700000000000   1.51855000000000   0.00000000000000
  -1.38598888592188   0.00000000000000   5.63000873803774
     Ga    O
      4    6
Direct
   0.09050000000000   0.09050000000000   0.79460000000000   Ga   (4i)
  -0.09050000000000  -0.09050000000000  -0.79460000000000   Ga   (4i)
   0.65866000000000   0.65866000000000   0.31402000000000   Ga   (4i)
  -0.65866000000000  -0.65866000000000  -0.31402000000000   Ga   (4i)
   0.16450000000000   0.16450000000000   0.10980000000000   O   (4i)
  -0.16450000000000  -0.16450000000000  -0.10980000000000   O   (4i)
   0.17330000000000   0.17330000000000   0.56320000000000   O   (4i)
  -0.17330000000000  -0.17330000000000  -0.56320000000000   O   (4i)
   0.49590000000000   0.49590000000000   0.25660000000000   O   (4i)
  -0.49590000000000  -0.49590000000000  -0.25660000000000   O   (4i)
```

$K_2Ti_2O_5$: A2B5C2_mC18_12_i_a2i_i - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'K2O5Ti2'
_chemical_formula_sum 'K2 O5 Ti2'

loop_
_publ_author_name
 'S. Andersson'
 'A. D. Wadsley'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 15
_journal_year 1961
_journal_page_first 663
_journal_page_last 669
_publ_Section_title
;
The Crystal Structure of K$_{2}$Ti$_{2}$O$_{5}$
;

_aflow_title 'K$_{2}$Ti$_{2}$O$_{5}$ Structure'
_aflow_proto 'A2B5C2_mC18_12_i_a2i_i'
_aflow_params 'a,b/a,c/a,\beta,x_{2},z_{2},x_{3},z_{3},x_{4},z_{4},x_{5
    ↪ },z_{5}'
_aflow_params_values '11.37,0.334212840809,0.582233948989,100.1,0.4022,
    ↪ 0.6439,0.128,0.664,0.325,0.008,0.1495,-0.0928'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mC18'

_symmetry_space_group_name_H-M "C 1 2/m 1"
_symmetry_Int_Tables_number 12

_cell_length_a     11.37000
_cell_length_b     3.80000
_cell_length_c     6.62000
_cell_angle_alpha  90.00000
_cell_angle_beta   100.10000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z
3 -x,-y,-z
4 x,-y,z
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 2 a 0.00000 0.00000 0.00000 1.00000
K1 K 4 i 0.40220 0.00000 0.64390 1.00000
O2 O 4 i 0.12800 0.00000 0.66400 1.00000
O3 O 4 i 0.32500 0.00000 0.00800 1.00000
Ti1 Ti 4 i 0.14950 0.00000 -0.09280 1.00000
```

$K_2Ti_2O_5$: A2B5C2_mC18_12_i_a2i_i - POSCAR

```
A2B5C2_mC18_12_i_a2i_i & a,b/a,c/a,beta,x2,z2,x3,z3,x4,z4,x5,z5 --params
    ↪ =11.37,0.334212840809,0.582233948989,100.1,0.4022,0.6439,0.128,
    ↪ 0.664,0.325,0.008,0.1495,-0.0928 & C2/m C_{2h}^{3} #12 (ai^4) &
    ↪ mC18 & None & K2O5Ti2 & K2O5Ti2 & S. Andersson and A. D.
    ↪ Wadsley, Acta Chem. Scand. 15, 663-669 (1961)
   1.00000000000000
   5.68500000000000  -1.90000000000000   0.00000000000000
   5.68500000000000   1.90000000000000   0.00000000000000
  -1.16092772672895   0.00000000000000   6.51741105143077
     K     O     Ti
     2     5     2
Direct
   0.40220000000000   0.40220000000000   0.64390000000000   K   (4i)
  -0.40220000000000  -0.40220000000000  -0.64390000000000   K   (4i)
   0.00000000000000   0.00000000000000   0.00000000000000   O   (2a)
   0.12800000000000   0.12800000000000   0.66400000000000   O   (4i)
  -0.12800000000000  -0.12800000000000  -0.66400000000000   O   (4i)
   0.32500000000000   0.32500000000000   0.00800000000000   O   (4i)
  -0.32500000000000  -0.32500000000000  -0.00800000000000   O   (4i)
   0.14950000000000   0.14950000000000  -0.09280000000000   Ti   (4i)
  -0.14950000000000  -0.14950000000000   0.09280000000000   Ti   (4i)
```

CaC_2-III: A2B_mC12_12_2i_i - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'C2Ca'
_chemical_formula_sum 'C2 Ca'
```

```
loop_
_publ_author_name
 'M. Knapp'
 'U. Ruschewitz'
_journal_name_full_name
;
 Chemistry - A European Journal
;
_journal_volume 7
_journal_year 2001
_journal_page_first 874
_journal_page_last 880
_publ_Section_title
;
 Structural Phase Transitions in CaC$_{2}$
;

_aflow_title 'CaC$_{2}$-III Structure'
_aflow_proto 'A2B_mC12_12_2i_i'
_aflow_params 'a,b/a,c/a,\beta,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3}'
_aflow_params_values '7.2286,0.532938604986,1.0204465595,107.338,0.439,
    ↪ 0.065,-0.075,0.447,0.2086,0.2486'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mC12'

_symmetry_space_group_name_H-M "C 1 2/m 1"
_symmetry_Int_Tables_number 12

_cell_length_a      7.22860
_cell_length_b      3.85240
_cell_length_c      7.37640
_cell_angle_alpha  90.00000
_cell_angle_beta   107.33800
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z
3 -x,-y,-z
4 x,-y,z
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
C1 C 4 i 0.43900 0.00000 0.06500 1.00000
C2 C 4 i -0.07500 0.00000 0.44700 1.00000
Ca1 Ca 4 i 0.20860 0.00000 0.24860 1.00000
```

CaC_2-III: A2B_mC12_12_2i_i - POSCAR

```
A2B_mC12_12_2i_i & a,b/a,c/a,beta,x1,z1,x2,z2,x3,z3 --params=7.2286,
    ↪ 0.532938604986,1.0204465595,107.338,0.439,0.065,-0.075,0.447,
    ↪ 0.2086,0.2486 & C2/m C_{2h}^{3} #12 (i^3) & mC12 & None & C2Ca
    ↪ & C2Ca & M. Knapp and U. Ruschewitz, Chem. Eur. J. 7, 874-880 (
    ↪ 2001)
   1.00000000000000
   3.61430000000000  -1.92620000000000   0.00000000000000
   3.61430000000000   1.92620000000000   0.00000000000000
  -2.19822643227335   0.00000000000000   7.04124119118603
     C     Ca
     4      2
Direct
   0.43900000000000   0.43900000000000   0.06500000000000   C    (4i)
  -0.43900000000000  -0.43900000000000  -0.06500000000000   C    (4i)
  -0.07500000000000  -0.07500000000000   0.44700000000000   C    (4i)
   0.07500000000000   0.07500000000000  -0.44700000000000   C    (4i)
   0.20860000000000   0.20860000000000   0.24860000000000   Ca   (4i)
  -0.20860000000000  -0.20860000000000  -0.24860000000000   Ca   (4i)
```

Tolbachite ($CuCl_2$): A2B_mC6_12_i_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Tolbachite'
_chemical_formula_sum 'Cl2 Cu'

loop_
_publ_author_name
 'P. C. Burns'
 'F. C. Hawthorne'
_journal_name_full_name
;
 American Mineralogist
;
_journal_volume 78
_journal_year 1993
_journal_page_first 187
_journal_page_last 189
_publ_Section_title
;
 Tolbachite, CuCl$_{2}$, the first example of Cu$_{2}^{+}$ octahedrally
    ↪ coordinated by Cl$^{-}$
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'Tolbachite (CuCl$_{2}$) Structure'
_aflow_proto 'A2B_mC6_12_i_a'
_aflow_params 'a,b/a,c/a,\beta,x_{2},z_{2}'
_aflow_params_values '6.9038,0.477925200614,0.988441148353,122.197,
    ↪ 0.5048,0.2294'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mC6'

_symmetry_space_group_name_H-M "C 1 2/m 1"
_symmetry_Int_Tables_number 12

_cell_length_a      6.90380
_cell_length_b      3.29950
_cell_length_c      6.82400
_cell_angle_alpha  90.00000
_cell_angle_beta   122.19700
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z
3 -x,-y,-z
4 x,-y,z
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cu1 Cu 2 a 0.00000 0.00000 0.00000 1.00000
Cl1 Cl 4 i 0.50480 0.00000 0.22940 1.00000
```

Tolbachite ($CuCl_2$): A2B_mC6_12_i_a - POSCAR

```
A2B_mC6_12_i_a & a,b/a,c/a,beta,x2,z2 --params=6.9038,0.477925200614,
    ↪ 0.988441148353,122.197,0.5048,0.2294 & C2/m C_{2h}^{3} #12 (ai)
    ↪ & mC6 & None & Cl2Cu & Tolbachite & P. C. Burns and F. C.
    ↪ Hawthorne, Am. Mineral. 78, 187-189 (1993)
   1.00000000000000
   3.45190000000000  -1.64975000000000   0.00000000000000
   3.45190000000000   1.64975000000000   0.00000000000000
  -3.63604535488460   0.00000000000000   5.77461255645971
     Cl     Cu
      2      1
Direct
   0.50480000000000   0.50480000000000   0.22940000000000   Cl   (4i)
  -0.50480000000000  -0.50480000000000  -0.22940000000000   Cl   (4i)
   0.00000000000000   0.00000000000000   0.00000000000000   Cu   (2a)
```

δ-Ni_3Sn_4 ($D7_a$): A3B4_mC14_12_ai_2i - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Ni3Sn4'
_chemical_formula_sum 'Ni3 Sn4'

loop_
_publ_author_name
 'W. Jeitschko'
 'B. Jaberg'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 38
_journal_year 1982
_journal_page_first 598
_journal_page_last 600
_publ_Section_title
;
 Structure refinement of Ni$_{3}$Sn$_{4}$
;

_aflow_title '$\delta$-Ni$_{3}$Sn$_{4}$ ($D7_{a}$) Structure'
_aflow_proto 'A3B4_mC14_12_ai_2i'
_aflow_params 'a,b/a,c/a,\beta,x_{2},z_{2},x_{3},z_{3},x_{4},z_{4}'
_aflow_params_values '12.214,0.332405436384,0.427296544948,105.0,0.2147,
    ↪ 0.3369,0.4286,0.6864,0.1718,0.8123'
_aflow_Strukturbericht '$D7_{a}$'
_aflow_Pearson 'mC14'

_symmetry_space_group_name_H-M "C 1 2/m 1"
_symmetry_Int_Tables_number 12

_cell_length_a      12.21400
_cell_length_b      4.06000
_cell_length_c      5.21900
```

```
_cell_angle_alpha  90.00000
_cell_angle_beta   105.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z
3 -x,-y,-z
4 x,-y,z
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ni1 Ni 2 a 0.00000 0.00000 0.00000 1.00000
Ni2 Ni 4 i 0.21470 0.00000 0.33690 1.00000
Sn1 Sn 4 i 0.42860 0.00000 0.68640 1.00000
Sn2 Sn 4 i 0.17180 0.00000 0.81230 1.00000
```

δ-Ni_3Sn_4 ($D7_a$): A3B4_mC14_12_ai_2i - POSCAR

```
A3B4_mC14_12_ai_2i & a,b/a,c/a,beta,x2,z2,x3,z3,x4,z4 --params=12.214,
      ↪ 0.332405436384,0.427296544948,105.0,0.2147,0.3369,0.4286,0.6864
      ↪ ,0.1718,0.8123 & C2/m C_{2h}^{3} #12 (ai^3) & mC14 & $D7_{a}$ &
      ↪  Ni3Sn4 & Ni3Sn4 & W. Jeitschko and B. Jaberg, Acta
      ↪ Crystallogr. Sect. B Struct. Sci. 38, 598-600 (1982)
   1.00000000000000
   6.10700000000000  -2.03000000000000   0.00000000000000
   6.10700000000000   2.03000000000000   0.00000000000000
  -1.35077659639006   0.00000000000000   5.04116688740265
     Ni      Sn
      3       4
Direct
   0.00000000000000   0.00000000000000   0.00000000000000   Ni   (2a)
   0.21470000000000   0.21470000000000   0.33690000000000   Ni   (4i)
  -0.21470000000000  -0.21470000000000  -0.33690000000000   Ni   (4i)
   0.42860000000000   0.42860000000000   0.68640000000000   Sn   (4i)
  -0.42860000000000  -0.42860000000000  -0.68640000000000   Sn   (4i)
   0.17180000000000   0.17180000000000   0.81230000000000   Sn   (4i)
  -0.17180000000000  -0.17180000000000  -0.81230000000000   Sn   (4i)
```

$LiOH{\cdot}H_2O$ ($B36$): A3BC2_mC24_12_ij_h_gi - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'H3LiO2'
_chemical_formula_sum 'H3 Li O2'

loop_
_publ_author_name
 'N. W. Alcock'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 27
_journal_year 1971
_journal_page_first 1682
_journal_page_last 1683
_publ_Section_title
;
 Refinement of the crystal structure of lithium hydroxide monohydrate
;

_aflow_title 'LiOH$\cdot$H$_{2}$O ($B36$) Structure'
_aflow_proto 'A3BC2_mC24_12_ij_h_gi'
_aflow_params 'a,b/a,c/a,\beta,y_{1},y_{2},x_{3},z_{3},x_{4},z_{4},x_{5
      ↪ },y_{5},z_{5}'
_aflow_params_values '7.37,1.12075983718,0.432835820896,110.3,0.2066,
      ↪ 0.3474,0.237,0.631,0.2857,0.3952,0.107,0.118,0.004'
_aflow_Strukturbericht '$B36$'
_aflow_Pearson 'mC24'

_symmetry_space_group_name_H-M "C 1 2/m 1"
_symmetry_Int_Tables_number 12

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_cell_length_b     8.26000
_cell_length_c     3.19000
_cell_angle_alpha  90.00000
_cell_angle_beta   110.30000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z
3 -x,-y,-z
4 x,-y,z
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 4 g 0.00000 0.20660 0.00000 1.00000
Li1 Li 4 h 0.00000 0.34740 0.50000 1.00000
H1 H 4 i 0.23700 0.00000 0.63100 1.00000
O2 O 4 i 0.28570 0.00000 0.39520 1.00000
H2 H 8 j 0.10700 0.11800 0.00400 1.00000
```

$LiOH{\cdot}H_2O$ ($B36$): A3BC2_mC24_12_ij_h_gi - POSCAR

```
A3BC2_mC24_12_ij_h_gi & a,b/a,c/a,beta,y1,y2,x3,z3,x4,z4,x5,y5,z5 --
      ↪ params=7.37,1.12075983718,0.432835820896,110.3,0.2066,0.3474,
      ↪ 0.237,0.631,0.2857,0.3952,0.107,0.118,0.004 & C2/m C_{2h}^{3} #
      ↪ 12 (ghi^2j) & mC24 & $B36$ & H3LiO2 & H3LiO2 & N. W. Alcock,
      ↪ Acta Crystallogr. Sect. B Struct. Sci. 27, 1682-1683 (1971)
   1.00000000000000
   3.68500000000000  -4.13000000000000   0.00000000000000
   3.68500000000000   4.13000000000000   0.00000000000000
  -1.10672472851869   0.00000000000000   2.99186570141195
     H     Li     O
     6      2      4
Direct
   0.23700000000000   0.23700000000000   0.63100000000000   H   (4i)
  -0.23700000000000  -0.23700000000000  -0.63100000000000   H   (4i)
  -0.01100000000000   0.22500000000000   0.00400000000000   H   (8j)
  -0.22500000000000   0.01100000000000  -0.00400000000000   H   (8j)
   0.01100000000000  -0.22500000000000  -0.00400000000000   H   (8j)
   0.22500000000000  -0.01100000000000   0.00400000000000   H   (8j)
  -0.34740000000000   0.34740000000000   0.50000000000000   Li   (4h)
   0.34740000000000  -0.34740000000000   0.50000000000000   Li   (4h)
  -0.20660000000000   0.20660000000000   0.00000000000000   O   (4g)
   0.20660000000000  -0.20660000000000   0.00000000000000   O   (4g)
   0.28570000000000   0.28570000000000   0.39520000000000   O   (4i)
  -0.28570000000000  -0.28570000000000  -0.39520000000000   O   (4i)
```

Staurolite ($Al_5Fe_2O_{10}(OH)_2Si_2$): A5B2C10D2E2_mC84_12_acghj_bdi_5j_2i_j - CIF

```
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_audit_creation_method FINDSYM

_chemical_name_mineral 'Staurolite'
_chemical_formula_sum 'Al5 Fe2 O10 (OH)2 Si2'

loop_
_publ_author_name
 'J. V. Smith'
_journal_name_full_name
;
 American Mineralogist
;
_journal_volume 53
_journal_year 1968
_journal_page_first 1139
_journal_page_last 1155
_publ_Section_title
;
 The crystal structure of staurolite
;

_aflow_title 'Staurolite (Al$_{5}$Fe$_{2}$O$_{10}$(OH)$_{2}$Si$_{2}$)
      ↪ Structure'
_aflow_proto 'A5B2C10D2E2_mC84_12_acghj_bdi_5j_2i_j'
_aflow_params 'a,b/a,c/a,\beta,y_{5},y_{6},x_{7},z_{7},x_{8},z_{8},x_{9
      ↪ },z_{9},x_{10},y_{10},z_{10},x_{11},y_{11},z_{11},x_{12},y_{12
      ↪ },z_{12},x_{13},y_{13},z_{13},x_{14},y_{14},z_{14},x_{15},y_{15
      ↪ },z_{15},x_{16},y_{16},z_{16}'
_aflow_params_values '7.8713,2.11151906292,0.718559831286,90.0,0.67511,
      ↪ 0.67477,0.39281,0.24815,0.23274,-0.03687,0.23438,0.53428,
      ↪ 0.26356,0.41042,0.25122,0.25569,0.16153,0.01527,0.25519,0.16127
      ↪ ,0.48391,0.00143,0.08917,0.24702,0.02156,0.24925,0.25002,
      ↪ 0.52741,0.10004,0.24944,0.13414,0.16612,0.24902'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mC84'

_symmetry_space_group_name_H-M "C 1 2/m 1"
_symmetry_Int_Tables_number 12

_cell_length_a     7.87130
_cell_length_b     16.62040
_cell_length_c     5.65600
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z
3 -x,-y,-z
4 x,-y,z
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z
7 -x+1/2,-y+1/2,-z
```

```
8 x+1/2,-y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Al1 Al 2 a 0.00000 0.00000 0.00000 1.00000
Fe1 Fe 2 b 0.00000 0.50000 0.00000 1.00000
Al2 Al 2 c 0.00000 0.00000 0.50000 1.00000
Fe2 Fe 2 d 0.00000 0.50000 0.50000 1.00000
Al3 Al 4 g 0.00000 0.67511 0.00000 1.00000
Al4 Al 4 h 0.00000 0.67477 0.50000 1.00000
Fe3 Fe 4 i 0.39281 0.00000 0.24815 1.00000
OH1 OH 4 i 0.23274 0.00000 -0.03687 1.00000
OH2 OH 4 i 0.23438 0.00000 0.53428 1.00000
Al5 Al 8 j 0.26356 0.41042 0.25122 1.00000
O1 O 8 j 0.25569 0.16153 0.01527 1.00000
O2 O 8 j 0.25519 0.16127 0.48391 1.00000
O3 O 8 j 0.00143 0.08917 0.24702 1.00000
O4 O 8 j 0.02156 0.24925 0.25002 1.00000
O5 O 8 j 0.52741 0.10004 0.24944 1.00000
Si1 Si 8 j 0.13414 0.16612 0.24902 1.00000
```

Staurolite ($Al_5Fe_2O_{10}(OH)_2Si_2$): A5B2C10D2E2_mC84_12_acghj_bdi_5j_2i_j - POSCAR

```
A5B2C10D2E2_mC84_12_acghj_bdi_5j_2i_j & a,b/a,c/a,beta,y5,y6,x7,z7,x8,z8
    ↪ ,x9,z9,x10,y10,z10,x11,y11,z11,x12,y12,z12,x13,y13,z13,x14,y14,
    ↪ z14,x15,y15,z15,x16,y16,z16 --params=7.8713,2.11151906292,
    ↪ 0.718559831286,90.0,0.67511,0.67477,0.39281,0.24815,0.23274,-
    ↪ 0.03687,0.23438,0.53428,0.26356,0.41042,0.25122,0.25569,0.16153
    ↪ ,0.01527,0.25519,0.16127,0.48391,0.00143,0.08917,0.24702,
    ↪ 0.02156,0.24925,0.25002,0.52741,0.10004,0.24944,0.13414,0.16612
    ↪ ,0.24902 & C2/m C_{2h}^{3} #12 (abcdghi^3j^7) & mC84 & None &
    ↪ Al5Fe2O10(OH)2Si2 & Staurolite & J. V. Smith, Am. Mineral. 53,
    ↪ 1139-1155 (1968)
   1.00000000000000
   3.93565000000000  -8.31020000000000   0.00000000000000
   3.93565000000000   8.31020000000000   0.00000000000000
   0.00000000000000   0.00000000000000   5.65600000000000
    Al    Fe    O    OH    Si
    10    4    20    4    4
Direct
   0.00000000000000   0.00000000000000   0.00000000000000  Al  (2a)
   0.00000000000000   0.00000000000000   0.50000000000000  Al  (2c)
  -0.67511000000000   0.67511000000000   0.00000000000000  Al  (4g)
   0.67511000000000  -0.67511000000000   0.00000000000000  Al  (4g)
  -0.67477000000000   0.67477000000000   0.50000000000000  Al  (4h)
   0.67477000000000  -0.67477000000000   0.50000000000000  Al  (4h)
  -0.14686000000000   0.67398000000000   0.25122000000000  Al  (8j)
  -0.67398000000000   0.14686000000000  -0.25122000000000  Al  (8j)
   0.14686000000000  -0.67398000000000  -0.25122000000000  Al  (8j)
   0.67398000000000  -0.14686000000000   0.25122000000000  Al  (8j)
   0.50000000000000   0.50000000000000   0.00000000000000  Fe  (2b)
   0.50000000000000   0.50000000000000   0.50000000000000  Fe  (2d)
   0.39281000000000   0.39281000000000   0.24815000000000  Fe  (4i)
  -0.39281000000000  -0.39281000000000  -0.24815000000000  Fe  (4i)
   0.09416000000000   0.41722000000000   0.01527000000000  O  (8j)
  -0.41722000000000  -0.09416000000000  -0.01527000000000  O  (8j)
  -0.09416000000000  -0.41722000000000  -0.01527000000000  O  (8j)
   0.41722000000000   0.09416000000000   0.01527000000000  O  (8j)
   0.09392000000000   0.41646000000000   0.48391000000000  O  (8j)
  -0.41646000000000  -0.09392000000000  -0.48391000000000  O  (8j)
  -0.09392000000000  -0.41646000000000  -0.48391000000000  O  (8j)
   0.41646000000000   0.09392000000000   0.48391000000000  O  (8j)
  -0.08774000000000   0.09060000000000   0.24702000000000  O  (8j)
  -0.09060000000000   0.08774000000000  -0.24702000000000  O  (8j)
   0.08774000000000  -0.09060000000000  -0.24702000000000  O  (8j)
   0.09060000000000  -0.08774000000000   0.24702000000000  O  (8j)
  -0.22769000000000   0.27081000000000   0.25002000000000  O  (8j)
  -0.27081000000000   0.22769000000000  -0.25002000000000  O  (8j)
   0.22769000000000  -0.27081000000000  -0.25002000000000  O  (8j)
   0.27081000000000  -0.22769000000000   0.25002000000000  O  (8j)
   0.42737000000000   0.62745000000000   0.24944000000000  O  (8j)
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  -0.42737000000000  -0.62745000000000  -0.24944000000000  O  (8j)
   0.62745000000000   0.42737000000000   0.24944000000000  O  (8j)
   0.23274000000000   0.23274000000000  -0.03687000000000  OH  (4i)
  -0.23274000000000  -0.23274000000000   0.03687000000000  OH  (4i)
   0.23438000000000   0.23438000000000   0.53428000000000  OH  (4i)
  -0.23438000000000  -0.23438000000000  -0.53428000000000  OH  (4i)
  -0.03198000000000   0.30026000000000   0.24902000000000  Si  (8j)
  -0.30026000000000   0.03198000000000  -0.24902000000000  Si  (8j)
   0.03198000000000  -0.30026000000000  -0.24902000000000  Si  (8j)
   0.30026000000000  -0.03198000000000   0.24902000000000  Si  (8j)
```

Manganese-leonite [$K_2Mn(SO_4)_2{\cdot}4H_2O$, $H4_{23}$]: A8B2CD15E2_mC112_12_2i3j_j_ad_g4i5j_2i - CIF

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_audit_creation_method FINDSYM

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_chemical_formula_sum 'H8 K2 Mn O15 S2'

loop_
_publ_author_name
 'B. Hertweck'
 'G. Giester'
 'E. Libowitzky'
_journal_name_full_name
;
 American Mineralogist
;
_journal_volume 86
_journal_year 2001
_journal_page_first 1282
_journal_page_last 1292
_publ_Section_title
;
 The crystal structures of the low-temperature phases of leonite-type
    ↪ compounds, K$_{2}$$Me$(SO$_{4}$)$_{2}$$\cdot$4H$_{2}$O ($Me^{2
    ↪ +}$ = Mg, Mn, Fe)
;

_aflow_title 'Manganese-leonite [K$_{2}$Mn(SO$_{4}$)$_{2}$$\cdot$4H$_{2}
    ↪ $O, $H4_{23}$] Structure'
_aflow_proto 'A8B2CD15E2_mC112_12_2i3j_j_ad_g4i5j_2i'
_aflow_params 'a,b/a,c/a,\beta,y_{3},x_{4},z_{4},x_{5},z_{5},x_{6},z_{6
    ↪ },x_{7},z_{7},x_{8},z_{8},x_{9},z_{9},x_{10},z_{10},x_{11},z_{
    ↪ 11},x_{12},y_{12},z_{12},x_{13},y_{13},z_{13},x_{14},y_{14},z_{
    ↪ 14},x_{15},y_{15},z_{15},x_{16},y_{16},z_{16},x_{17},y_{17},z_{
    ↪ 17},x_{18},y_{18},z_{18},x_{19},y_{19},z_{19},x_{20},y_{20},z_{
    ↪ 20}'
_aflow_params_values '12.017,0.798368977282,0.827910460181,95.03,0.2267,
    ↪ 0.061,0.743,-0.041,0.728,0.1728,0.0594,0.6841,0.0965,0.6686,
    ↪ 0.4354,0.0104,0.7806,0.70695,-0.04418,0.21077,0.52021,0.072,
    ↪ 0.276,0.621,0.079,0.349,0.726,0.047,0.271,-0.027,0.17022,
    ↪ 0.26326,0.24936,0.1588,0.3743,0.8871,0.1519,0.0973,0.5987,
    ↪ 0.1723,0.1457,0.548,0.1918,0.0236,0.3754,0.0387,0.3351,0.6481'
_aflow_Strukturbericht '$H4_{23}$'
_aflow_Pearson 'mC112'

_symmetry_space_group_name_H-M "C 1 2/m 1"
_symmetry_Int_Tables_number 12

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_cell_length_b      9.59400
_cell_length_c      9.94900
_cell_angle_alpha   90.00000
_cell_angle_beta    95.03000
_cell_angle_gamma   90.00000

loop_
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_space_group_symop_operation_xyz
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2 -x,y,-z
3 -x,-y,-z
4 x,-y,z
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
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Mn2 Mn 2 d 0.00000 0.50000 0.50000 1.00000
O1 O 4 g 0.00000 0.22670 0.00000 1.00000
H1 H 4 i 0.06100 0.00000 0.74300 1.00000
H2 H 4 i -0.04100 0.00000 0.72800 1.00000
O2 O 4 i 0.17280 0.00000 0.05940 1.00000
O3 O 4 i 0.68410 0.00000 0.09650 1.00000
O4 O 4 i 0.66860 0.00000 0.43540 1.00000
O5 O 4 i 0.01040 0.00000 0.78060 1.00000
S1 S 4 i 0.70695 0.00000 -0.04418 1.00000
S2 S 4 i 0.21077 0.00000 0.52021 1.00000
H3 H 8 j 0.07200 0.27600 0.62100 1.00000
H4 H 8 j 0.07900 0.34900 0.72600 1.00000
H5 H 8 j 0.04700 0.27100 -0.02700 1.00000
K1 K 8 j 0.17022 0.26326 0.24936 1.00000
O6 O 8 j 0.15880 0.37430 0.88710 1.00000
O7 O 8 j 0.15190 0.09730 0.59870 0.50000
O8 O 8 j 0.17230 0.14570 0.54800 0.50000
O9 O 8 j 0.19180 0.02360 0.37540 0.50000
O10 O 8 j 0.03870 0.33510 0.64810 1.00000
```

Manganese-leonite [$K_2Mn(SO_4)_2{\cdot}4H_2O$, $H4_{23}$]: A8B2CD15E2_mC112_12_2i3j_j_ad_g4i5j_2i - POSCAR

```
A8B2CD15E2_mC112_12_2i3j_j_ad_g4i5j_2i & a,b/a,c/a,beta,y3,x4,z4,x5,z5,
    ↪ x6,z6,x7,z7,x8,z8,x9,z9,x10,z10,x11,z11,x12,y12,z12,x13,y13,z13
    ↪ ,x14,y14,z14,x15,y15,z15,x16,y16,z16,x17,y17,z17,x18,y18,z18,
    ↪ x19,y19,z19,x20,y20,z20 --params=12.017,0.798368977282,
    ↪ 0.827910460181,95.03,0.2267,0.061,0.743,-0.041,0.728,0.1728,
    ↪ 0.0594,0.6841,0.0965,0.6686,0.4354,0.0104,0.7806,0.70695,-
    ↪ 0.04418,0.21077,0.52021,0.072,0.276,0.621,0.079,0.349,0.726,
    ↪ 0.047,0.271,-0.027,0.17022,0.26326,0.24936,0.1588,0.3743,0.8871
    ↪ ,0.1519,0.0973,0.5987,0.1723,0.1457,0.548,0.1918,0.0236,0.3754,
    ↪ 0.0387,0.3351,0.6481 & C2/m C_{2h}^{3} #12 (adgi^8j^9) & mC112
    ↪ & $H4_{23}$ & H8K2MnO12S2 & Manganese-leonite & B. Hertweck and
    ↪  G. Giester and E. Libowitzky, Am. Mineral. 86, 1282-1292 (2001
    ↪ )
   1.00000000000000
   6.00850000000000  -4.79700000000000   0.00000000000000
   6.00850000000000   4.79700000000000   0.00000000000000
  -0.87230182681668   0.00000000000000   9.91068567370252
    H    K    Mn    O    S
```

```
    16      4      2     30      4
Direct
   0.06100000000000   0.06100000000000   0.74300000000000   H   (4i)
  -0.06100000000000  -0.06100000000000  -0.74300000000000   H   (4i)
  -0.04100000000000  -0.04100000000000   0.72800000000000   H   (4i)
   0.04100000000000   0.04100000000000  -0.72800000000000   H   (4i)
  -0.20400000000000   0.34800000000000   0.62100000000000   H   (8j)
  -0.34800000000000   0.20400000000000  -0.62100000000000   H   (8j)
   0.20400000000000  -0.34800000000000  -0.62100000000000   H   (8j)
   0.34800000000000  -0.20400000000000   0.62100000000000   H   (8j)
  -0.27000000000000   0.42800000000000   0.72600000000000   H   (8j)
  -0.42800000000000   0.27000000000000  -0.72600000000000   H   (8j)
   0.27000000000000  -0.42800000000000  -0.72600000000000   H   (8j)
   0.42800000000000  -0.27000000000000   0.72600000000000   H   (8j)
  -0.22400000000000   0.31800000000000  -0.02700000000000   H   (8j)
  -0.31800000000000   0.22400000000000   0.02700000000000   H   (8j)
   0.22400000000000  -0.31800000000000   0.02700000000000   H   (8j)
   0.31800000000000  -0.22400000000000  -0.02700000000000   H   (8j)
  -0.09304000000000   0.43348000000000   0.24936000000000   K   (8j)
  -0.43348000000000   0.09304000000000  -0.24936000000000   K   (8j)
   0.09304000000000  -0.43348000000000  -0.24936000000000   K   (8j)
   0.43348000000000  -0.09304000000000   0.24936000000000   K   (8j)
   0.00000000000000   0.00000000000000   0.00000000000000   Mn  (2a)
   0.50000000000000   0.50000000000000   0.50000000000000   Mn  (2d)
  -0.22670000000000   0.22670000000000   0.00000000000000   O   (4g)
   0.22670000000000  -0.22670000000000   0.00000000000000   O   (4g)
   0.17280000000000   0.17280000000000   0.05940000000000   O   (4i)
  -0.17280000000000  -0.17280000000000  -0.05940000000000   O   (4i)
   0.68410000000000   0.68410000000000   0.09650000000000   O   (4i)
  -0.68410000000000  -0.68410000000000  -0.09650000000000   O   (4i)
   0.66860000000000   0.66860000000000   0.43540000000000   O   (4i)
  -0.66860000000000  -0.66860000000000  -0.43540000000000   O   (4i)
   0.01040000000000   0.01040000000000   0.78060000000000   O   (4i)
  -0.01040000000000  -0.01040000000000  -0.78060000000000   O   (4i)
  -0.21550000000000   0.53310000000000   0.88710000000000   O   (8j)
  -0.53310000000000   0.21550000000000  -0.88710000000000   O   (8j)
   0.21550000000000  -0.53310000000000  -0.88710000000000   O   (8j)
   0.53310000000000  -0.21550000000000   0.88710000000000   O   (8j)
   0.05460000000000   0.24920000000000   0.59870000000000   O   (8j)
  -0.24920000000000  -0.05460000000000  -0.59870000000000   O   (8j)
  -0.05460000000000  -0.24920000000000  -0.59870000000000   O   (8j)
   0.24920000000000   0.05460000000000   0.59870000000000   O   (8j)
   0.02660000000000   0.31800000000000   0.54800000000000   O   (8j)
  -0.31800000000000  -0.02660000000000  -0.54800000000000   O   (8j)
  -0.02660000000000  -0.31800000000000  -0.54800000000000   O   (8j)
   0.31800000000000   0.02660000000000   0.54800000000000   O   (8j)
   0.16820000000000   0.21540000000000   0.37540000000000   O   (8j)
  -0.21540000000000  -0.16820000000000  -0.37540000000000   O   (8j)
  -0.16820000000000  -0.21540000000000  -0.37540000000000   O   (8j)
   0.21540000000000   0.16820000000000   0.37540000000000   O   (8j)
  -0.29640000000000   0.37380000000000   0.64810000000000   O   (8j)
  -0.37380000000000   0.29640000000000  -0.64810000000000   O   (8j)
   0.29640000000000  -0.37380000000000  -0.64810000000000   O   (8j)
   0.37380000000000  -0.29640000000000   0.64810000000000   O   (8j)
   0.70695000000000   0.70695000000000  -0.04418000000000   S   (4i)
  -0.70695000000000  -0.70695000000000   0.04418000000000   S   (4i)
   0.21077000000000   0.21077000000000   0.52021000000000   S   (4i)
  -0.21077000000000  -0.21077000000000  -0.52021000000000   S   (4i)
```

Monoclinic $FeTlSe_2$: AB2C_mC16_12_g_2i_i - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'FeSe2Tl'
_chemical_formula_sum 'Fe Se2 Tl'

loop_
_publ_author_name
 'K. Klepp'
 'H. Boller'
_journal_name_full_name
;
 Monatshefte f{\"u}r Chemie - Chemical Monthly
;
_journal_volume 110
_journal_year 1979
_journal_page_first 1045
_journal_page_last 1055
_publ_Section_title
;
 Die Kristallstruktur von TlFeSe$_{2}$ und TlFeS$_{2}$
;

_aflow_title 'Monoclinic FeTlSe$_{2}$ Structure'
_aflow_proto 'AB2C_mC16_12_g_2i_i'
_aflow_params 'a,b/a,c/a,\beta,y_{1},x_{2},z_{2},x_{3},z_{3},x_{4},z_{4
      ↪ }'
_aflow_params_values '11.973,0.458531696317,0.593836131295,118.2,0.2493,
      ↪ 0.5357,0.2899,0.1783,0.0907,0.1752,0.6314'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mC16'

_symmetry_space_group_name_H-M "C 1 2/m 1"
_symmetry_Int_Tables_number 12

_cell_length_a     11.97300
_cell_length_b     5.49000
_cell_length_c     7.11000
_cell_angle_alpha  90.00000
_cell_angle_beta   118.20000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z
3 -x,-y,-z
4 x,-y,z
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Fe1 Fe 4 g 0.00000 0.24930 0.00000 1.00000
Se1 Se 4 i 0.53570 0.00000 0.28990 1.00000
Se2 Se 4 i 0.17830 0.00000 0.09070 1.00000
Tl1 Tl 4 i 0.17520 0.00000 0.63140 1.00000
```

Monoclinic $FeTlSe_2$: AB2C_mC16_12_g_2i_i - POSCAR

```
AB2C_mC16_12_g_2i_i & a,b/a,c/a,beta,y1,x2,z2,x3,z3,x4,z4 --params=
      ↪ 11.973,0.458531696317,0.593836131295,118.2,0.2493,0.5357,0.2899
      ↪ ,0.1783,0.0907,0.1752,0.6314 & C2/m C_{2h}^{3} #12 (gi^3) &
      ↪ mC16 & None & FeSe2Tl & FeSe2Tl & K. Klepp and H. Boller,
      ↪ Monatshefte f{"u}r Chemie - Chemical Monthly 110, 1045-1055 (
      ↪ 1979)
   1.00000000000000
   5.98650000000000  -2.74500000000000   0.00000000000000
   5.98650000000000   2.74500000000000   0.00000000000000
  -3.35983593821897   0.00000000000000   6.26606754418210
     Fe     Se     Tl
      2      4      2
Direct
  -0.24930000000000   0.24930000000000   0.00000000000000   Fe   (4g)
   0.24930000000000  -0.24930000000000   0.00000000000000   Fe   (4g)
   0.53570000000000   0.53570000000000   0.28990000000000   Se   (4i)
  -0.53570000000000  -0.53570000000000  -0.28990000000000   Se   (4i)
   0.17830000000000   0.17830000000000   0.09070000000000   Se   (4i)
  -0.17830000000000  -0.17830000000000  -0.09070000000000   Se   (4i)
   0.17520000000000   0.17520000000000   0.63140000000000   Tl   (4i)
  -0.17520000000000  -0.17520000000000  -0.63140000000000   Tl   (4i)
```

$NbTe_2$: AB2_mC18_12_ai_3i - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'NbTe2'
_chemical_formula_sum 'Nb Te2'

loop_
_publ_author_name
 'B. E. Brown'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 20
_journal_year 1966
_journal_page_first 264
_journal_page_last 267
_publ_Section_title
;
 The Crystal Structures of NbTe$_{2}$ and TaTe$_{2}$
;

# Found in Pearson's Handbook of Crystallographic Data for Intermetallic
      ↪  Phases, 1991

_aflow_title 'NbTe$_{2}$ Structure'
_aflow_proto 'AB2_mC18_12_ai_3i'
_aflow_params 'a,b/a,c/a,\beta,x_{2},z_{2},x_{3},z_{3},x_{4},z_{4},x_{5
      ↪ },z_{5}'
_aflow_params_values '14.73541,0.247159732915,0.636222541483,110.41211,
      ↪ 0.36025,0.7087,0.3503,-0.0096,0.703,0.6208,0.0039,0.3098'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mC18'

_symmetry_space_group_name_H-M "C 1 2/m 1"
_symmetry_Int_Tables_number 12

_cell_length_a     14.73541
_cell_length_b     3.64200
_cell_length_c     9.37500
_cell_angle_alpha  90.00000
_cell_angle_beta   110.41211
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z
3 -x,-y,-z
4 x,-y,z
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z
```

```
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Nb1 Nb 2 a 0.00000 0.00000 0.00000 1.00000
Nb2 Nb 4 i 0.36025 0.00000 0.70870 1.00000
Te1 Te 4 i 0.35030 0.00000 -0.00960 1.00000
Te2 Te 4 i 0.70300 0.00000 0.62080 1.00000
Te3 Te 4 i 0.00390 0.00000 0.30980 1.00000
```

$NbTe_2$: AB2_mC18_12_ai_3i - POSCAR

```
AB2_mC18_12_ai_3i & a,b/a,c/a,beta,x2,z2,x3,z3,x4,z4,x5,z5 --params=
    ↪ 14.73541,0.247159732915,0.636222541483,110.41211,0.36025,0.7087
    ↪ ,0.3503,-0.0096,0.703,0.6208,0.0039,0.3098 & C2/m C_{2h}^{3} #
    ↪ 12 (ai^4) & mC18 & None & NbTe2 & NbTe2 & B. E. Brown, Acta
    ↪ Cryst. 20, 264-267 (1966)
   1.00000000000000
   7.36770500000000   -1.82100000000000    0.00000000000000
   7.36770500000000    1.82100000000000    0.00000000000000
  -3.26972008938351    0.00000000000000    8.78632776176042
     Nb      Te
      3       6
Direct
    0.00000000000000    0.00000000000000    0.00000000000000    Nb    (2a)
    0.36025000000000    0.36025000000000    0.70870000000000    Nb    (4i)
   -0.36025000000000   -0.36025000000000   -0.70870000000000    Nb    (4i)
    0.35030000000000    0.35030000000000   -0.00960000000000    Te    (4i)
   -0.35030000000000   -0.35030000000000    0.00960000000000    Te    (4i)
    0.70300000000000    0.70300000000000    0.62080000000000    Te    (4i)
   -0.70300000000000   -0.70300000000000   -0.62080000000000    Te    (4i)
    0.00390000000000    0.00390000000000    0.30980000000000    Te    (4i)
   -0.00390000000000   -0.00390000000000   -0.30980000000000    Te    (4i)
```

$D2_2$ ($MgZn_5$?) (*problematic*): AB5_mC48_12_2i_ac5i2j - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'MgZn5'
_chemical_formula_sum 'Mg Zn5'

loop_
_publ_author_name
 'C. Hermann'
 'O. Lohrmann'
 'H. Philipp'
_journal_year 1937
_publ_Section_title
;
 Strukturbericht Band II 1928-1932
;

_aflow_title '$D2_{2}$ (MgZn$_{5}$?) ({\em{problematic}}) Structure'
_aflow_proto 'AB5_mC48_12_2i_ac5i2j'
_aflow_params 'a,b/a,c/a,\beta,x_{3},z_{3},x_{4},z_{4},x_{5},z_{5},x_{6
    ↪ },z_{6},x_{7},z_{7},x_{8},z_{8},x_{9},z_{9},x_{10},y_{10},z_{10
    ↪ },x_{11},y_{11},z_{11}'
_aflow_params_values '9.92,1.66129032258,1.0,120.0,0.66667,0.19792,
    ↪ 0.66667,0.38542,0.66667,0.66667,0.0,0.25,0.33333,0.08333,
    ↪ 0.16667,0.79167,0.16667,0.29167,0.41667,0.25,0.04167,-0.08333,
    ↪ 0.25,0.54167'
_aflow_Strukturbericht '$D2_{2}$'
_aflow_Pearson 'mC48'

_symmetry_space_group_name_H-M "C 1 2/m 1"
_symmetry_Int_Tables_number 12

_cell_length_a     9.92000
_cell_length_b     16.48000
_cell_length_c     9.92000
_cell_angle_alpha  90.00000
_cell_angle_beta   120.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z
3 -x,-y,-z
4 x,-y,z
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Zn1 Zn 2 a 0.00000 0.00000 0.00000 1.00000
Zn2 Zn 2 c 0.00000 0.00000 0.50000 1.00000
Mg1 Mg 4 i 0.66667 0.00000 0.19792 1.00000
Mg2 Mg 4 i 0.66667 0.00000 0.38542 1.00000
Zn3 Zn 4 i 0.66667 0.00000 0.66667 1.00000
Zn4 Zn 4 i 0.00000 0.00000 0.25000 1.00000
Zn5 Zn 4 i 0.33333 0.00000 0.08333 1.00000
Zn6 Zn 4 i 0.16667 0.00000 0.79167 1.00000
Zn7 Zn 4 i 0.16667 0.00000 0.29167 1.00000
Zn8 Zn 8 j 0.41667 0.25000 0.04167 1.00000
Zn9 Zn 8 j -0.08333 0.25000 0.54167 1.00000
```

$D2_2$ ($MgZn_5$?) (*problematic*): AB5_mC48_12_2i_ac5i2j - POSCAR

```
AB5_mC48_12_2i_ac5i2j & a,b/a,c/a,beta,x3,z3,x4,z4,x5,z5,x6,z6,x7,z7,x8,
    ↪ z8,x9,z9,x10,y10,z10,x11,y11,z11 --params=9.92,1.66129032258,
    ↪ 1.0,120.0,0.66667,0.19792,0.66667,0.38542,0.66667,0.66667,0.0,
    ↪ 0.25,0.33333,0.08333,0.16667,0.79167,0.16667,0.29167,0.41667,
    ↪ 0.25,0.04167,-0.08333,0.25,0.54167 & C2/m C_{2h}^{3} #12 (aci^
    ↪ 7j^2) & mC48 & $D2_{2}$ & MgZn5 & MgZn5 & C. Hermann and O.
    ↪ Lohrmann and H. Philipp, (1937)
   1.00000000000000
   4.96000000000000   -8.24000000000000    0.00000000000000
   4.96000000000000    8.24000000000000    0.00000000000000
  -4.96000000000000    0.00000000000000    8.59097200554163
     Mg      Zn
      4      20
Direct
    0.66667000000000    0.66667000000000    0.19792000000000    Mg    (4i)
   -0.66667000000000   -0.66667000000000   -0.19792000000000    Mg    (4i)
    0.66667000000000    0.66667000000000    0.38542000000000    Mg    (4i)
   -0.66667000000000   -0.66667000000000   -0.38542000000000    Mg    (4i)
    0.00000000000000    0.00000000000000    0.00000000000000    Zn    (2a)
    0.00000000000000    0.00000000000000    0.50000000000000    Zn    (2c)
    0.66667000000000    0.66667000000000    0.66667000000000    Zn    (4i)
   -0.66667000000000   -0.66667000000000   -0.66667000000000    Zn    (4i)
    0.00000000000000    0.00000000000000    0.25000000000000    Zn    (4i)
    0.00000000000000    0.00000000000000   -0.25000000000000    Zn    (4i)
    0.33333000000000    0.33333000000000    0.08333000000000    Zn    (4i)
   -0.33333000000000   -0.33333000000000   -0.08333000000000    Zn    (4i)
    0.16667000000000    0.16667000000000    0.79167000000000    Zn    (4i)
   -0.16667000000000   -0.16667000000000   -0.79167000000000    Zn    (4i)
    0.16667000000000    0.16667000000000    0.29167000000000    Zn    (4i)
   -0.16667000000000   -0.16667000000000   -0.29167000000000    Zn    (4i)
    0.16667000000000    0.66667000000000    0.04167000000000    Zn    (8j)
   -0.66667000000000   -0.16667000000000   -0.04167000000000    Zn    (8j)
   -0.16667000000000   -0.66667000000000   -0.04167000000000    Zn    (8j)
    0.66667000000000    0.16667000000000    0.04167000000000    Zn    (8j)
   -0.33333000000000    0.16667000000000    0.54167000000000    Zn    (8j)
   -0.16667000000000    0.33333000000000   -0.54167000000000    Zn    (8j)
    0.33333000000000   -0.16667000000000   -0.54167000000000    Zn    (8j)
    0.16667000000000   -0.33333000000000    0.54167000000000    Zn    (8j)
```

Chrysotile ($H_4Mg_3Si_2O_9$, $S4_5$): AB6C11D6E4_mC112_12_e_gi2j_i5j_2i2j_2j - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Chrysotile'
_chemical_formula_sum '(H2O) Mg6 O11 (OH)6 Si4'

loop_
_publ_author_name
 'B. E. Warren'
 'W. L. Bragg'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 76
_journal_year 1931
_journal_page_first 201
_journal_page_last 210
_publ_Section_title
;
 The Structure of Chrysotile H$_{4}$Mg$_{3}$Si$_{2}$O$_{9}$
;

# Found in Strukturbericht Band II 1928-1932, 1937

_aflow_title 'Chrysotile (H$_{4}$Mg$_{3}$Si$_{2}$O$_{9}$, $S4_{5}$)
    ↪ Structure'
_aflow_proto 'AB6C11D6E4_mC112_12_e_gi2j_i5j_2i2j_2j'
_aflow_params 'a,b/a,c/a,\beta,y_{2},x_{3},z_{3},x_{4},z_{4},x_{5},z_{5
    ↪ },x_{6},z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},
    ↪ z_{9},x_{10},y_{10},z_{10},x_{11},y_{11},z_{11},x_{12},y_{12},
    ↪ z_{12},x_{13},y_{13},z_{13},x_{14},y_{14},z_{14},x_{15},y_{15},
    ↪ z_{15},x_{16},y_{16},z_{16},x_{17},y_{17},z_{17}'
_aflow_params_values '15.31076,1.20830056771,0.348121190588,107.07121,
    ↪ 0.75,0.32,0.72,0.9,0.9,0.75,0.45,0.4,0.4,0.32,0.17,0.22,0.32,
    ↪ 0.08,0.72,0.4,0.37,0.25,0.4,0.37,0.75,0.4,0.25,0.4,0.25,0.42,-
    ↪ 0.05,0.25,0.33,0.45,0.4,0.17,0.9,0.4,0.08,0.4,0.37,0.42,-0.03,
    ↪ 0.37,0.33,0.47'
_aflow_Strukturbericht '$S4_{5}$'
_aflow_Pearson 'mC112'

_symmetry_space_group_name_H-M "C 1 2/m 1"
_symmetry_Int_Tables_number 12

_cell_length_a     15.31076
_cell_length_b     18.50000
_cell_length_c     5.33000
_cell_angle_alpha  90.00000
_cell_angle_beta   107.07121
```

```
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z
3 -x,-y,-z
4 x,-y,z
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
H2O1 H2O 4 e 0.25000 0.25000 0.00000 1.00000
Mg1 Mg 4 g 0.00000 0.75000 0.00000 1.00000
Mg2 Mg 4 i 0.32000 0.00000 0.72000 1.00000
O1 O 4 i 0.90000 0.00000 0.90000 1.00000
OH1 OH 4 i 0.75000 0.00000 0.45000 1.00000
OH2 OH 4 i 0.40000 0.00000 0.40000 1.00000
Mg3 Mg 8 j 0.32000 0.17000 0.22000 1.00000
Mg4 Mg 8 j 0.32000 0.08000 0.72000 1.00000
O2 O 8 j 0.40000 0.37000 0.25000 1.00000
O3 O 8 j 0.40000 0.37000 0.75000 1.00000
O4 O 8 j 0.40000 0.25000 0.40000 1.00000
O5 O 8 j 0.25000 0.42000 -0.05000 1.00000
O6 O 8 j 0.25000 0.33000 0.45000 1.00000
OH3 OH 8 j 0.40000 0.17000 0.90000 1.00000
OH4 OH 8 j 0.40000 0.08000 0.40000 1.00000
Si1 Si 8 j 0.37000 0.42000 -0.03000 1.00000
Si2 Si 8 j 0.37000 0.33000 0.47000 1.00000
```

Chrysotile ($H_4Mg_3Si_2O_9$, $S4_5$): AB6C11D6E4_mC112_12_e_gi2j_i5j_2i2j_2j - POSCAR

```
AB6C11D6E4_mC112_12_e_gi2j_i5j_2i2j_2j & a,b/a,c/a,beta,y2,x3,z3,x4,z4,
    ↪ x5,z5,x6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10,x11,y11,z11,
    ↪ x12,y12,z12,x13,y13,z13,x14,y14,z14,x15,y15,z15,x16,y16,z16,x17
    ↪ ,y17,z17 --params=15.31076,1.20830056771,0.348121190588,
    ↪ 107.07121,0.75,0.32,0.72,0.9,0.9,0.75,0.45,0.4,0.4,0.32,0.17,
    ↪ 0.22,0.32,0.08,0.72,0.4,0.37,0.25,0.4,0.37,0.75,0.4,0.25,0.4,
    ↪ 0.25,0.42,-0.05,0.25,0.33,0.45,0.4,0.17,0.9,0.4,0.08,0.4,0.37,
    ↪ 0.42,-0.03,0.37,0.33,0.47 & C2/m C_{2h}^{3} #12 (egi^4j^11) &
    ↪ mC112 & $S4_{5}$ & (H2O)(OH)6O11Mg6Si4 & Chrysotile & B. E.
    ↪ Warren and W. L. Bragg, Zeitschrift f{"u}r Kristallographie -
    ↪ Crystalline Materials 76, 201-210 (1931)
   1.00000000000000
   7.65538000000000  -9.25000000000000   0.00000000000000
   7.65538000000000   9.25000000000000   0.00000000000000
  -1.56467491181852   0.00000000000000   5.09516363037790
   H2O  Mg  O  OH  Si
     2  12  22  12   8
Direct
   0.00000000000000   0.50000000000000   0.00000000000000  H2O  (4e)
   0.50000000000000   0.00000000000000   0.00000000000000  H2O  (4e)
  -0.75000000000000   0.75000000000000   0.00000000000000  Mg   (4g)
   0.75000000000000  -0.75000000000000   0.00000000000000  Mg   (4g)
   0.32000000000000   0.32000000000000   0.72000000000000  Mg   (4i)
  -0.32000000000000  -0.32000000000000  -0.72000000000000  Mg   (4i)
   0.15000000000000   0.49000000000000   0.22000000000000  Mg   (8j)
  -0.49000000000000  -0.15000000000000  -0.22000000000000  Mg   (8j)
  -0.15000000000000  -0.49000000000000  -0.22000000000000  Mg   (8j)
   0.49000000000000   0.15000000000000   0.22000000000000  Mg   (8j)
   0.24000000000000   0.40000000000000   0.72000000000000  Mg   (8j)
  -0.40000000000000  -0.24000000000000  -0.72000000000000  Mg   (8j)
  -0.24000000000000  -0.40000000000000  -0.72000000000000  Mg   (8j)
   0.40000000000000   0.24000000000000   0.72000000000000  Mg   (8j)
   0.90000000000000   0.90000000000000   0.90000000000000  O    (4i)
  -0.90000000000000  -0.90000000000000  -0.90000000000000  O    (4i)
   0.03000000000000   0.77000000000000   0.25000000000000  O    (8j)
  -0.77000000000000  -0.03000000000000  -0.25000000000000  O    (8j)
  -0.03000000000000  -0.77000000000000  -0.25000000000000  O    (8j)
   0.77000000000000   0.03000000000000   0.25000000000000  O    (8j)
   0.03000000000000   0.77000000000000   0.75000000000000  O    (8j)
  -0.77000000000000  -0.03000000000000  -0.75000000000000  O    (8j)
  -0.03000000000000  -0.77000000000000  -0.75000000000000  O    (8j)
   0.77000000000000   0.03000000000000   0.75000000000000  O    (8j)
   0.15000000000000   0.65000000000000   0.40000000000000  O    (8j)
  -0.65000000000000  -0.15000000000000  -0.40000000000000  O    (8j)
  -0.15000000000000  -0.65000000000000  -0.40000000000000  O    (8j)
   0.65000000000000   0.15000000000000   0.40000000000000  O    (8j)
  -0.17000000000000   0.67000000000000  -0.05000000000000  O    (8j)
  -0.67000000000000   0.17000000000000   0.05000000000000  O    (8j)
   0.17000000000000  -0.67000000000000   0.05000000000000  O    (8j)
   0.67000000000000  -0.17000000000000  -0.05000000000000  O    (8j)
  -0.08000000000000   0.58000000000000   0.45000000000000  O    (8j)
  -0.58000000000000   0.08000000000000  -0.45000000000000  O    (8j)
   0.08000000000000  -0.58000000000000  -0.45000000000000  O    (8j)
   0.58000000000000  -0.08000000000000   0.45000000000000  O    (8j)
   0.75000000000000   0.75000000000000   0.45000000000000  OH   (4i)
  -0.75000000000000  -0.75000000000000  -0.45000000000000  OH   (4i)
   0.40000000000000   0.40000000000000   0.40000000000000  OH   (4i)
  -0.40000000000000  -0.40000000000000  -0.40000000000000  OH   (4i)
   0.23000000000000   0.57000000000000   0.90000000000000  OH   (8j)
  -0.57000000000000  -0.23000000000000  -0.90000000000000  OH   (8j)
  -0.23000000000000  -0.57000000000000  -0.90000000000000  OH   (8j)
   0.57000000000000   0.23000000000000   0.90000000000000  OH   (8j)
   0.32000000000000   0.48000000000000   0.40000000000000  OH   (8j)
  -0.48000000000000  -0.32000000000000  -0.40000000000000  OH   (8j)
  -0.32000000000000  -0.48000000000000  -0.40000000000000  OH   (8j)
   0.48000000000000   0.32000000000000   0.40000000000000  OH   (8j)
  -0.05000000000000   0.79000000000000  -0.03000000000000  Si   (8j)
  -0.79000000000000   0.05000000000000   0.03000000000000  Si   (8j)
   0.05000000000000  -0.79000000000000   0.03000000000000  Si   (8j)
   0.79000000000000  -0.05000000000000  -0.03000000000000  Si   (8j)
   0.04000000000000   0.70000000000000   0.47000000000000  Si   (8j)
  -0.70000000000000  -0.04000000000000  -0.47000000000000  Si   (8j)
  -0.04000000000000  -0.70000000000000  -0.47000000000000  Si   (8j)
   0.70000000000000   0.04000000000000   0.47000000000000  Si   (8j)
```

Sr_2NiTeO_6: AB6C2D_mC40_12_ad_gh4i_j_bc - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'NiO6Sr2Te'
_chemical_formula_sum 'Ni O6 Sr2 Te'

loop_
_publ_author_name
 'D. Iwanaga'
 'Y. Inaguma'
 'M. Itoh'
_journal_name_full_name
;
 Materials Research Bulletin
;
_journal_volume 35
_journal_year 2000
_journal_page_first 449
_journal_page_last 457
_publ_Section_title
;
 Structure and Magnetic Properties of Sr$_{2}$Ni$A$O$_{6}$ ($A$ = W, Te)
;

_aflow_title 'Sr$_{2}$NiTeO$_{6}$ Structure'
_aflow_proto 'AB6C2D_mC40_12_ad_gh4i_j_bc'
_aflow_params 'a,b/a,c/a,\beta,y_{5},y_{6},x_{7},z_{7},x_{8},z_{8},x_{9
    ↪ },z_{9},x_{10},z_{10},x_{11},y_{11},z_{11}'
_aflow_params_values '7.9174,0.99483416273,0.999823174274,90.378,0.26,
    ↪ 0.759,0.266,0.028,-0.031,0.255,0.24,0.454,0.535,0.247,0.248,
    ↪ 0.251,0.249'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mC40'

_symmetry_space_group_name_H-M "C 1 2/m 1"
_symmetry_Int_Tables_number 12

_cell_length_a     7.91740
_cell_length_b     7.87650
_cell_length_c     7.91600
_cell_angle_alpha  90.00000
_cell_angle_beta   90.37800
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z
3 -x,-y,-z
4 x,-y,z
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ni1 Ni 2 a 0.00000 0.00000 0.00000 1.00000
Te1 Te 2 b 0.00000 0.50000 0.00000 1.00000
Te2 Te 2 c 0.00000 0.00000 0.50000 1.00000
Ni2 Ni 2 d 0.00000 0.50000 0.50000 1.00000
O1 O 4 g 0.00000 0.26000 0.00000 1.00000
O2 O 4 h 0.00000 0.75900 0.50000 1.00000
O3 O 4 i 0.26600 0.00000 0.02800 1.00000
O4 O 4 i -0.03100 0.00000 0.25500 1.00000
O5 O 4 i 0.24000 0.00000 0.45400 1.00000
O6 O 4 i 0.53500 0.00000 0.24700 1.00000
Sr1 Sr 8 j 0.24800 0.25100 0.24900 1.00000
```

Sr_2NiTeO_6: AB6C2D_mC40_12_ad_gh4i_j_bc - POSCAR

```
AB6C2D_mC40_12_ad_gh4i_j_bc & a,b/a,c/a,beta,y5,y6,x7,z7,x8,z8,x9,z9,x10
    ↪ ,z10,x11,y11,z11 --params=7.9174,0.99483416273,0.999823174274,
    ↪ 90.378,0.26,0.759,0.266,0.028,-0.031,0.255,0.24,0.454,0.535,
    ↪ 0.247,0.248,0.251,0.249 & C2/m C_{2h}^{3} #12 (abcdghi^4j) &
    ↪ mC40 & None & NiO6Sr2Te & NiO6Sr2Te & D. Iwanaga and Y. Inaguma
    ↪ and M. Itoh, Mater. Res. Bull. 35, 449-457 (2000)
   1.00000000000000
   3.95870000000000  -3.93825000000000   0.00000000000000
   3.95870000000000   3.93825000000000   0.00000000000000
  -0.05222420079162   0.00000000000000   7.91582772885134
```

```
    Ni     O     Sr     Te
     2    12      4      2
Direct
   0.00000000000000   0.00000000000000   0.00000000000000   Ni   (2a)
   0.50000000000000   0.50000000000000   0.50000000000000   Ni   (2d)
  -0.26000000000000   0.26000000000000   0.00000000000000   O   (4g)
   0.26000000000000  -0.26000000000000   0.00000000000000   O   (4g)
  -0.75900000000000   0.75900000000000   0.50000000000000   O   (4h)
   0.75900000000000  -0.75900000000000   0.50000000000000   O   (4h)
   0.26600000000000   0.26600000000000   0.02800000000000   O   (4i)
  -0.26600000000000  -0.26600000000000  -0.02800000000000   O   (4i)
  -0.03100000000000  -0.03100000000000   0.25500000000000   O   (4i)
   0.03100000000000   0.03100000000000  -0.25500000000000   O   (4i)
   0.24000000000000   0.24000000000000   0.45400000000000   O   (4i)
  -0.24000000000000  -0.24000000000000  -0.45400000000000   O   (4i)
   0.53500000000000   0.53500000000000   0.24700000000000   O   (4i)
  -0.53500000000000  -0.53500000000000  -0.24700000000000   O   (4i)
  -0.00300000000000   0.49900000000000   0.24900000000000   Sr   (8j)
  -0.49900000000000   0.00300000000000  -0.24900000000000   Sr   (8j)
   0.00300000000000  -0.49900000000000  -0.24900000000000   Sr   (8j)
   0.49900000000000  -0.00300000000000   0.24900000000000   Sr   (8j)
   0.50000000000000   0.50000000000000   0.00000000000000   Te   (2b)
   0.00000000000000   0.00000000000000   0.50000000000000   Te   (2c)
```

Ta_2PdSe_6: AB6C2_mC18_12_a_3i_i - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'PdSe6Ta2'
_chemical_formula_sum 'Pd Se6 Ta2'

loop_
_publ_author_name
 'D. A. Keszler'
 'P. J. Squattrito'
 'N. E. Brese'
 'J. A. Ibers'
 'M. Shang'
 'J. Lu'
_journal_year 1985
_publ_Section_title
;
New layered ternary chalcogenides: tantalum palladium sulfide (Ta$_{2}
    ↪ $PdS$_{6}$), tantalum palladium selenide (Ta$_{2}$PdSe$_{6}$),
    ↪  niobium palladium sulfide (Nb$_{2}$PdS$_{6}$), niobium
    ↪ palladium selenide (Nb$_{2}$PdSe$_{6}$)
;

_aflow_title 'Ta$_{2}$PdSe$_{6}$ Structure'
_aflow_proto 'AB6C2_mC18_12_a_3i_i'
_aflow_params 'a,b/a,c/a,\beta,x_{2},z_{2},x_{3},z_{3},x_{4},z_{4},x_{5
    ↪ },z_{5}'
_aflow_params_values '12.46083,0.270848731585,0.836461134611,116.31902,-
    ↪ 0.00589,0.76131,0.65848,0.53954,0.21851,0.12521,0.67999,0.28976
    ↪ '
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mC18'

_symmetry_space_group_name_H-M "C 1 2/m 1"
_symmetry_Int_Tables_number 12

_cell_length_a    12.46083
_cell_length_b    3.37500
_cell_length_c    10.42300
_cell_angle_alpha  90.00000
_cell_angle_beta  116.31902
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z
3 -x,-y,-z
4 x,-y,z
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Pd1 Pd 2 a 0.00000 0.00000 0.00000 1.00000
Se1 Se 4 i -0.00589 0.00000 0.76131 1.00000
Se2 Se 4 i 0.65848 0.00000 0.53954 1.00000
Se3 Se 4 i 0.21851 0.00000 0.12521 1.00000
Ta1 Ta 4 i 0.67999 0.00000 0.28976 1.00000
```

Ta_2PdSe_6: AB6C2_mC18_12_a_3i_i - POSCAR

```
AB6C2_mC18_12_a_3i_i & a,b/a,c/a,beta,x2,z2,x3,z3,x4,z4,x5,z5 --params=
    ↪ 12.46083,0.270848731585,0.836461134611,116.31902,-0.00589,
    ↪ 0.76131,0.65848,0.53954,0.21851,0.12521,0.67999,0.28976 & C2/m
    ↪ C_{2h}^{3} #12 (ai^4) & mC18 & None & PdSe6Ta2 & PdSe6Ta2 & D.
    ↪ A. Keszler et al., (1985)
   1.00000000000000
   6.23041500000000  -1.68750000000000   0.00000000000000
   6.23041500000000   1.68750000000000   0.00000000000000
  -4.62123264277420   0.00000000000000   9.34254450679033
    Pd    Se    Ta
     1     6     2
Direct
   0.00000000000000   0.00000000000000   0.00000000000000   Pd   (2a)
  -0.00589000000000  -0.00589000000000   0.76131000000000   Se   (4i)
   0.00589000000000   0.00589000000000  -0.76131000000000   Se   (4i)
   0.65848000000000   0.65848000000000   0.53954000000000   Se   (4i)
  -0.65848000000000  -0.65848000000000  -0.53954000000000   Se   (4i)
   0.21851000000000   0.21851000000000   0.12521000000000   Se   (4i)
  -0.21851000000000  -0.21851000000000  -0.12521000000000   Se   (4i)
   0.67999000000000   0.67999000000000   0.28976000000000   Ta   (4i)
  -0.67999000000000  -0.67999000000000  -0.28976000000000   Ta   (4i)
```

Sanidine ($KAlSi_3O_8$, $S6_7$): AB8C4_mC52_12_i_gi3j_2j - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Sanidine'
_chemical_formula_sum 'K O8 Si4'

loop_
_publ_author_name
 'T. A. Scambos'
 'J. R. Smyth'
 'T. C. {McCormick}'
_journal_name_full_name
;
 American Mineralogist
;
_journal_volume 72
_journal_year 1987
_journal_page_first 973
_journal_page_last 978
_publ_Section_title
;
 Crystal-structure refinement of high sanidine from the upper mantle
;

_aflow_title 'Sanidine ($KAlSi$_{3}$O$_{8}$, $S6_{7}$) Structure'
_aflow_proto 'AB8C4_mC52_12_i_gi3j_2j'
_aflow_params 'a,b/a,c/a,\beta,y_{1},x_{2},z_{2},x_{3},z_{3},x_{4},y_{4
    ↪ },z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{6},x_{7},y_{7},z_{7},
    ↪ x_{8},y_{8},z_{8}'
_aflow_params_values '8.595,1.51576497964,0.834787667248,115.94,0.147,
    ↪ 0.2866,0.138,0.64,0.2849,0.8302,0.1476,0.2269,0.0351,0.3105,
    ↪ 0.2569,0.1789,0.1265,0.4083,0.00991,0.1856,0.22381,0.71075,
    ↪ 0.11813,0.34438'
_aflow_Strukturbericht '$S6_{7}$'
_aflow_Pearson 'mC52'

_symmetry_space_group_name_H-M "C 1 2/m 1"
_symmetry_Int_Tables_number 12

_cell_length_a    8.59500
_cell_length_b    13.02800
_cell_length_c    7.17500
_cell_angle_alpha  90.00000
_cell_angle_beta  115.94000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z
3 -x,-y,-z
4 x,-y,z
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 4 g 0.00000 0.14700 0.00000 1.00000
K1 K 4 i 0.28660 0.00000 0.13800 1.00000
O2 O 4 i 0.64000 0.00000 0.28490 1.00000
O3 O 8 j 0.83020 0.14760 0.22690 1.00000
O4 O 8 j 0.03510 0.31050 0.25690 1.00000
O5 O 8 j 0.17890 0.12650 0.40830 1.00000
Si1 Si 8 j 0.00991 0.18560 0.22381 1.00000
Si2 Si 8 j 0.71075 0.11813 0.34438 1.00000
```

Sanidine ($KAlSi_3O_8$, $S6_7$): AB8C4_mC52_12_i_gi3j_2j - POSCAR

```
AB8C4_mC52_12_i_gi3j_2j & a,b/a,c/a,beta,y1,x2,z2,x3,z3,x4,y4,z4,x5,y5,
    ↪ z5,x6,y6,z6,x7,y7,z7,x8,y8,z8 --params=8.595,1.51576497964,
    ↪ 0.834787667248,115.94,0.147,0.2866,0.138,0.64,0.2849,0.8302,
    ↪ 0.1476,0.2269,0.0351,0.3105,0.2569,0.1789,0.1265,0.4083,0.00991
    ↪ ,0.1856,0.22381,0.71075,0.11813,0.34438 & C2/m C_{2h}^{3} #12 (
    ↪ gi^2j^5) & mC52 & $S6_{7}$ & AlKO8Si3 & Sanidine & T. A.
    ↪ Scambos and J. R. Smyth and T. C. {McCormick}, Am. Mineral. 72,
    ↪  973-978 (1987)
```

```
   1.00000000000000
   4.29750000000000   -6.51400000000000    0.00000000000000
   4.29750000000000    6.51400000000000    0.00000000000000
  -3.13855803775329    0.00000000000000    6.45213750966718
     K     O     Si
     2    16     8
Direct
   0.28660000000000   0.28660000000000   0.13800000000000   K   (4i)
  -0.28660000000000  -0.28660000000000  -0.13800000000000   K   (4i)
  -0.14700000000000   0.14700000000000   0.00000000000000   O   (4g)
   0.14700000000000  -0.14700000000000   0.00000000000000   O   (4g)
   0.64000000000000   0.64000000000000   0.28490000000000   O   (4i)
  -0.64000000000000  -0.64000000000000  -0.28490000000000   O   (4i)
   0.68260000000000   0.97780000000000   0.22690000000000   O   (8j)
  -0.97780000000000  -0.68260000000000  -0.22690000000000   O   (8j)
  -0.68260000000000  -0.97780000000000  -0.22690000000000   O   (8j)
   0.97780000000000   0.68260000000000   0.22690000000000   O   (8j)
  -0.27540000000000   0.34560000000000   0.25690000000000   O   (8j)
  -0.34560000000000   0.27540000000000  -0.25690000000000   O   (8j)
   0.27540000000000  -0.34560000000000  -0.25690000000000   O   (8j)
   0.34560000000000  -0.27540000000000   0.25690000000000   O   (8j)
   0.05240000000000   0.30540000000000   0.40830000000000   O   (8j)
  -0.30540000000000  -0.05240000000000  -0.40830000000000   O   (8j)
  -0.05240000000000  -0.30540000000000  -0.40830000000000   O   (8j)
   0.30540000000000   0.05240000000000   0.40830000000000   O   (8j)
  -0.17569000000000   0.19551000000000   0.22381000000000   Si  (8j)
  -0.19551000000000   0.17569000000000  -0.22381000000000   Si  (8j)
   0.17569000000000  -0.19551000000000  -0.22381000000000   Si  (8j)
   0.19551000000000  -0.17569000000000   0.22381000000000   Si  (8j)
   0.59262000000000   0.82888000000000   0.34438000000000   Si  (8j)
  -0.82888000000000  -0.59262000000000  -0.34438000000000   Si  (8j)
  -0.59262000000000  -0.82888000000000  -0.34438000000000   Si  (8j)
   0.82888000000000   0.59262000000000   0.34438000000000   Si  (8j)
```

MnPS$_3$: ABC3_mC20_12_g_i_ij - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'MnPS3'
_chemical_formula_sum 'Mn P S3'

loop_
_publ_author_name
 'G. Ouvrard'
 'R. Brec'
 'J. Rouxel'
_journal_name_full_name
;
 Materials Research Bulletin
;
_journal_volume 20
_journal_year 1985
_journal_page_first 1181
_journal_page_last 1189
_publ_Section_title
;
 Structural determination of some $M$PS$_{3}$ layered phases ($M$ = Mn,
      ↪ Fe, Co, Ni and Cd)
;

# Found in Electronic band structure of the magnetic layered
      ↪ semiconductors $M$PS$_{3}$ ($M$ = Mn, Fe and Ni), 1996

_aflow_title 'MnPS$_{3}$ Structure'
_aflow_proto 'ABC3_mC20_12_g_i_ij'
_aflow_params 'a,b/a,c/a,\beta,y_{1},x_{2},z_{2},x_{3},z_{3},x_{4},y_{4
      ↪ },z_{4}'
_aflow_params_values '6.077,1.73177554714,1.11831495804,107.35,0.33258,
      ↪ 0.0556,0.1686,0.7593,0.2497,0.2438,0.1612,0.2516'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mC20'

_symmetry_space_group_name_H-M "C 1 2/m 1"
_symmetry_Int_Tables_number 12

_cell_length_a      6.07700
_cell_length_b      10.52400
_cell_length_c      6.79600
_cell_angle_alpha   90.00000
_cell_angle_beta    107.35000
_cell_angle_gamma   90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z
3 -x,-y,-z
4 x,-y,z
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
```

```
Mn1 Mn 4 g 0.00000 0.33258 0.00000 1.00000
P1 P 4 i 0.05560 0.00000 0.16860 1.00000
S1 S 4 i 0.75930 0.00000 0.24970 1.00000
S2 S 8 j 0.24380 0.16120 0.25160 1.00000
```

MnPS$_3$: ABC3_mC20_12_g_i_ij - POSCAR

```
ABC3_mC20_12_g_i_ij & a,b/a,c/a,beta,y1,x2,z2,x3,z3,x4,y4,z4 --params=
      ↪ 6.077,1.73177554714,1.11831495804,107.35,0.33258,0.0556,0.1686,
      ↪ 0.7593,0.2497,0.2438,0.1612,0.2516 & C2/m C_{2h}^{3} #12 (gi^2j
      ↪ ) & mC20 & None & MnPS3 & MnPS3 & G. Ouvrard and R. Brec and J.
      ↪  Rouxel, Mater. Res. Bull. 20, 1181-1189 (1985)
   1.00000000000000
   3.03850000000000   -5.26200000000000    0.00000000000000
   3.03850000000000    5.26200000000000    0.00000000000000
  -2.02662120588260    0.00000000000000    6.48678830299455
    Mn     P     S
     2     2     6
Direct
  -0.33258000000000    0.33258000000000    0.00000000000000   Mn   (4g)
   0.33258000000000   -0.33258000000000    0.00000000000000   Mn   (4g)
   0.05560000000000    0.05560000000000    0.16860000000000   P    (4i)
  -0.05560000000000   -0.05560000000000   -0.16860000000000   P    (4i)
   0.75930000000000    0.75930000000000    0.24970000000000   S    (4i)
  -0.75930000000000   -0.75930000000000   -0.24970000000000   S    (4i)
   0.08260000000000    0.40500000000000    0.25160000000000   S    (8j)
  -0.40500000000000   -0.08260000000000   -0.25160000000000   S    (8j)
  -0.08260000000000   -0.40500000000000   -0.25160000000000   S    (8j)
   0.40500000000000    0.08260000000000    0.25160000000000   S    (8j)
```

AlNbO$_4$: ABC4_mC24_12_i_i_4i - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'AlNbO4'
_chemical_formula_sum 'Al Nb O4'

loop_
_publ_author_name
 'M. Ardit'
 'M. Dondi'
 'G. Cruciani'
_journal_name_full_name
;
 American Mineralogist
;
_journal_volume 97
_journal_year 2012
_journal_page_first 910
_journal_page_last 917
_publ_Section_title
;
 Structural stability, cation ordering, and local relaxation along the
      ↪ AlNbO$_{4}$-Al$_{0.5}$Cr$_{0.5}$NbO$_{4}$ join
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'AlNbO$_{4}$ Structure'
_aflow_proto 'ABC4_mC24_12_i_i_4i'
_aflow_params 'a,b/a,c/a,\beta,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3},x_{4
      ↪ },z_{4},x_{5},z_{5},x_{6},z_{6}'
_aflow_params_values '12.15449,0.307293847788,0.533825771382,107.6206,
      ↪ 0.19362,0.30116,0.1025,0.73176,0.1358,0.0099,0.0564,0.3636,
      ↪ 0.3615,0.2985,0.2622,0.6432'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mC24'

_symmetry_space_group_name_H-M "C 1 2/m 1"
_symmetry_Int_Tables_number 12

_cell_length_a      12.15449
_cell_length_b      3.73500
_cell_length_c      6.48838
_cell_angle_alpha   90.00000
_cell_angle_beta    107.62060
_cell_angle_gamma   90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z
3 -x,-y,-z
4 x,-y,z
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Al1 Al 4 i 0.19362 0.00000 0.30116 1.00000
Nb1 Nb 4 i 0.10250 0.00000 0.73176 1.00000
O1 O 4 i 0.13580 0.00000 0.00990 1.00000
O2 O 4 i 0.05640 0.00000 0.36360 1.00000
```

```
O3 O 4 i 0.36150 0.00000 0.29850 1.00000
O4 O 4 i 0.26220 0.00000 0.64320 1.00000
```

$AlNbO_4$: ABC4_mC24_12_i_i_4i - POSCAR

```
ABC4_mC24_12_i_i_4i & a,b/a,c/a,beta,x1,z1,x2,z2,x3,z3,x4,z4,x5,z5,x6,z6
    ↪ --params=12.15449,0.307293847788,0.533825771382,107.6206,
    ↪ 0.19362,0.30116,0.1025,0.73176,0.1358,0.0099,0.0564,0.3636,
    ↪ 0.3615,0.2985,0.2622,0.6432 & C2/m C_{2h}^{3} #12 (i^6) & mC24
    ↪ & None & AlNbO4 & AlNbO4 & M. Ardit and M. Dondi and G.
    ↪ Cruciani, Am. Mineral. 97, 910-917 (2012)
  1.00000000000000
  6.07724500000000  -1.86750000000000   0.00000000000000
  6.07724500000000   1.86750000000000   0.00000000000000
 -1.96411424527948   0.00000000000000   6.18395749143623
   Al   Nb   O
    2    2   8
Direct
  0.19362000000000   0.19362000000000   0.30116000000000  Al  (4i)
 -0.19362000000000  -0.19362000000000  -0.30116000000000  Al  (4i)
  0.10250000000000   0.10250000000000   0.73176000000000  Nb  (4i)
 -0.10250000000000  -0.10250000000000  -0.73176000000000  Nb  (4i)
  0.13580000000000   0.13580000000000   0.00990000000000  O  (4i)
 -0.13580000000000  -0.13580000000000  -0.00990000000000  O  (4i)
  0.05640000000000   0.05640000000000   0.36360000000000  O  (4i)
 -0.05640000000000  -0.05640000000000  -0.36360000000000  O  (4i)
  0.36150000000000   0.36150000000000   0.29850000000000  O  (4i)
 -0.36150000000000  -0.36150000000000  -0.29850000000000  O  (4i)
  0.26220000000000   0.26220000000000   0.64320000000000  O  (4i)
 -0.26220000000000  -0.26220000000000  -0.64320000000000  O  (4i)
```

SiAs: AB_mC24_12_3i_3i - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'AsSi'
_chemical_formula_sum 'As Si'

loop_
_publ_author_name
 'T. Wadsten'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 19
_journal_year 1965
_journal_page_first 1232
_journal_page_last 1238
_publ_Section_title
;
 The Crystal Structure of SiAs
;

_aflow_title 'SiAs Structure'
_aflow_proto 'AB_mC24_12_3i_3i'
_aflow_params 'a,b/a,c/a,\beta,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3},x_{4
    ↪ },z_{4},x_{5},z_{5},x_{6},z_{6}'
_aflow_params_values '15.98,0.229536921151,0.596307884856,106.0,0.6632,
    ↪ 0.1738,-0.0369,0.1761,0.3479,0.4543,0.2613,0.2076,0.3697,0.0838
    ↪ ,-0.0661,0.4116'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mC24'

_symmetry_space_group_name_H-M "C 1 2/m 1"
_symmetry_Int_Tables_number 12

_cell_length_a    15.98000
_cell_length_b    3.66800
_cell_length_c    9.52900
_cell_angle_alpha 90.00000
_cell_angle_beta  106.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z
3 -x,-y,-z
4 x,-y,z
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
As1 As 4 i 0.66320 0.00000 0.17380 1.00000
As2 As 4 i -0.03690 0.00000 0.17610 1.00000
As3 As 4 i 0.34790 0.00000 0.45430 1.00000
Si1 Si 4 i 0.26130 0.00000 0.20760 1.00000
Si2 Si 4 i 0.36970 0.00000 0.08380 1.00000
Si3 Si 4 i -0.06610 0.00000 0.41160 1.00000
```

SiAs: AB_mC24_12_3i_3i - POSCAR

```
AB_mC24_12_3i_3i & a,b/a,c/a,beta,x1,z1,x2,z2,x3,z3,x4,z4,x5,z5,x6,z6 --
    ↪ params=15.98,0.229536921151,0.596307884856,106.0,0.6632,0.1738
    ↪ ,-0.0369,0.1761,0.3479,0.4543,0.2613,0.2076,0.3697,0.0838,-
    ↪ 0.0661,0.4116 & C2/m C_{2h}^{3} #12 (i^6) & mC24 & None & AsSi
    ↪ & AsSi & T. Wadsten, Acta Chem. Scand. 19, 1232-1238 (1965)
  1.00000000000000
  7.99000000000000  -1.83400000000000   0.00000000000000
  7.99000000000000   1.83400000000000   0.00000000000000
 -2.62654836358018   0.00000000000000   9.15986270059624
   As   Si
    6    6
Direct
  0.66320000000000   0.66320000000000   0.17380000000000  As  (4i)
 -0.66320000000000  -0.66320000000000  -0.17380000000000  As  (4i)
 -0.03690000000000  -0.03690000000000   0.17610000000000  As  (4i)
  0.03690000000000   0.03690000000000  -0.17610000000000  As  (4i)
  0.34790000000000   0.34790000000000   0.45430000000000  As  (4i)
 -0.34790000000000  -0.34790000000000  -0.45430000000000  As  (4i)
  0.26130000000000   0.26130000000000   0.20760000000000  Si  (4i)
 -0.26130000000000  -0.26130000000000  -0.20760000000000  Si  (4i)
  0.36970000000000   0.36970000000000   0.08380000000000  Si  (4i)
 -0.36970000000000  -0.36970000000000  -0.08380000000000  Si  (4i)
 -0.06610000000000  -0.06610000000000   0.41160000000000  Si  (4i)
  0.06610000000000   0.06610000000000  -0.41160000000000  Si  (4i)
```

High-Temperature Mo_8O_{23}: A8B23_mP62_13_4g_c11g - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Mo8O23'
_chemical_formula_sum 'Mo8 O23'

loop_
_publ_author_name
 'H. Fujishita'
 'M. Sato'
 'S. Sato'
 'S. Hoshino'
_journal_name_full_name
;
 Journal of Solid State Chemistry
;
_journal_volume 66
_journal_year 1987
_journal_page_first 40
_journal_page_last 46
_publ_Section_title
;
 Structure Determination of low-dimensional conductor Mo${8}$O${23}$
;

_aflow_title 'High-Temperature Mo${8}$O${23}$ Structure'
_aflow_proto 'A8B23_mP62_13_4g_c11g'
_aflow_params 'a,b/a,c/a,\beta,x_{2},y_{2},z_{2},x_{3},y_{3},z_{3},x_{4
    ↪ },y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{6},x_{7},y_{7},
    ↪ z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{10},y_{10},z_{10},
    ↪ x_{11},y_{11},z_{11},x_{12},y_{12},z_{12},x_{13},y_{13},z_{13},
    ↪ x_{14},y_{14},z_{14},x_{15},y_{15},z_{15},x_{16},y_{16},z_{16}'
_aflow_params_values '13.384,0.303466826061,1.26143156007,106.27,0.06397
    ↪ ,0.5858,0.41655,0.18514,0.4131,0.24568,0.31501,0.59,0.07903,
    ↪ 0.44659,0.4079,0.40204,0.0658,-0.0012,0.4164,0.1911,-0.0029,
    ↪ 0.2458,0.3191,0.0069,0.077,0.4469,-0.0082,0.4074,0.0645,0.4943,
    ↪ 0.1643,0.1293,0.493,0.3303,0.1989,0.4945,0.496,0.2612,0.5144,
    ↪ 0.1632,0.3261,0.4891,0.3282,0.4113,0.5015,0.0045,0.457,0.4908,
    ↪ 0.1532'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mP62'

_symmetry_space_group_name_H-M "P 1 2/c 1"
_symmetry_Int_Tables_number 13

_cell_length_a    13.38400
_cell_length_b    4.06160
_cell_length_c    16.88300
_cell_angle_alpha 90.00000
_cell_angle_beta  106.27000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z+1/2
3 -x,-y,-z
4 x,-y,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 2 c 0.00000 0.50000 0.00000 1.00000
Mo1 Mo 4 g 0.06397 0.58580 0.41655 1.00000
Mo2 Mo 4 g 0.18514 0.41310 0.24568 1.00000
Mo3 Mo 4 g 0.31501 0.59000 0.07903 1.00000
Mo4 Mo 4 g 0.44659 0.40790 0.40204 1.00000
O2 O 4 g 0.06580 -0.00120 0.41640 1.00000
O3 O 4 g 0.19110 -0.00290 0.24580 1.00000
O4 O 4 g 0.31910 0.00690 0.07700 1.00000
```

```
O5 O 4 g 0.44690 -0.00820 0.40740 1.00000
O6 O 4 g 0.06450 0.49430 0.16430 1.00000
O7 O 4 g 0.12930 0.49300 0.33030 1.00000
O8 O 4 g 0.19890 0.49450 0.49600 1.00000
O9 O 4 g 0.26120 0.51440 0.16320 1.00000
O10 O 4 g 0.32610 0.48910 0.32820 1.00000
O11 O 4 g 0.41130 0.50150 0.00450 1.00000
O12 O 4 g 0.45700 0.49080 0.15320 1.00000
```

High-Temperature Mo_8O_{23}: A8B23_mP62_13_4g_c11g - POSCAR

```
A8B23_mP62_13_4g_c11g & a,b/a,c/a,beta,x2,y2,z2,x3,y3,z3,x4,y4,z4,x5,y5,
    ↪ z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10,x11,y11,z11,
    ↪ x12,y12,z12,x13,y13,z13,x14,y14,z14,x15,y15,z15,x16,y16,z16 --
    ↪ params=13.384,0.303466826061,1.26143156007,106.27,0.06397,
    ↪ 0.5858,0.41655,0.18514,0.4131,0.24568,0.31501,0.59,0.07903,
    ↪ 0.44659,0.4079,0.40204,0.0658,-0.0012,0.4164,0.1911,-0.0029,
    ↪ 0.2458,0.3191,0.0069,0.077,0.4469,-0.0082,0.4074,0.0645,0.4943,
    ↪ 0.1643,0.1293,0.493,0.3303,0.1989,0.4945,0.496,0.2612,0.5144,
    ↪ 0.1632,0.3261,0.4891,0.3282,0.4113,0.5015,0.0045,0.457,0.4908,
    ↪ 0.1532 & P2/c C_{2h}^{4} #13 (cg^15) & mP62 & None & Mo8O23 &
    ↪ Mo8O23 & H. Fujishita et al., J. Solid State Chem. 66, 40-46 (
    ↪ 1987)
   1.00000000000000
  13.38400000000000    0.00000000000000    0.00000000000000
   0.00000000000000    4.06160000000000    0.00000000000000
  -4.73001079735427    0.00000000000000   16.20687159376890
     Mo      O
     16     46
Direct
   0.06397000000000    0.58580000000000    0.41655000000000  Mo   (4g)
  -0.06397000000000    0.58580000000000    0.08345000000000  Mo   (4g)
  -0.06397000000000   -0.58580000000000   -0.41655000000000  Mo   (4g)
   0.06397000000000   -0.58580000000000    0.91655000000000  Mo   (4g)
   0.18514000000000    0.41310000000000    0.24568000000000  Mo   (4g)
  -0.18514000000000    0.41310000000000    0.25432000000000  Mo   (4g)
  -0.18514000000000   -0.41310000000000   -0.24568000000000  Mo   (4g)
   0.18514000000000   -0.41310000000000    0.74568000000000  Mo   (4g)
   0.31501000000000    0.59000000000000    0.07903000000000  Mo   (4g)
  -0.31501000000000    0.59000000000000    0.42097000000000  Mo   (4g)
  -0.31501000000000   -0.59000000000000   -0.07903000000000  Mo   (4g)
   0.31501000000000   -0.59000000000000    0.57903000000000  Mo   (4g)
   0.44659000000000    0.40790000000000    0.40204000000000  Mo   (4g)
  -0.44659000000000    0.40790000000000    0.09796000000000  Mo   (4g)
  -0.44659000000000   -0.40790000000000   -0.40204000000000  Mo   (4g)
   0.44659000000000   -0.40790000000000    0.90204000000000  Mo   (4g)
   0.00000000000000    0.50000000000000    0.00000000000000  O    (2c)
   0.00000000000000    0.50000000000000    0.50000000000000  O    (2c)
   0.06580000000000   -0.00120000000000    0.41640000000000  O    (4g)
  -0.06580000000000   -0.00120000000000    0.08360000000000  O    (4g)
  -0.06580000000000    0.00120000000000   -0.41640000000000  O    (4g)
   0.06580000000000    0.00120000000000    0.91640000000000  O    (4g)
   0.19110000000000   -0.00290000000000    0.24580000000000  O    (4g)
  -0.19110000000000   -0.00290000000000    0.25420000000000  O    (4g)
  -0.19110000000000    0.00290000000000   -0.24580000000000  O    (4g)
   0.19110000000000    0.00290000000000    0.74580000000000  O    (4g)
   0.31910000000000    0.00690000000000    0.07700000000000  O    (4g)
  -0.31910000000000    0.00690000000000    0.42300000000000  O    (4g)
  -0.31910000000000   -0.00690000000000   -0.07700000000000  O    (4g)
   0.31910000000000   -0.00690000000000    0.57700000000000  O    (4g)
   0.44690000000000   -0.00820000000000    0.40740000000000  O    (4g)
  -0.44690000000000   -0.00820000000000    0.09260000000000  O    (4g)
  -0.44690000000000    0.00820000000000   -0.40740000000000  O    (4g)
   0.44690000000000    0.00820000000000    0.90740000000000  O    (4g)
   0.06450000000000    0.49430000000000    0.16430000000000  O    (4g)
  -0.06450000000000    0.49430000000000    0.33570000000000  O    (4g)
  -0.06450000000000   -0.49430000000000   -0.16430000000000  O    (4g)
   0.06450000000000   -0.49430000000000    0.66430000000000  O    (4g)
   0.12930000000000    0.49300000000000    0.33030000000000  O    (4g)
  -0.12930000000000    0.49300000000000    0.16970000000000  O    (4g)
  -0.12930000000000   -0.49300000000000   -0.33030000000000  O    (4g)
   0.12930000000000   -0.49300000000000    0.83030000000000  O    (4g)
   0.19890000000000    0.49450000000000    0.49600000000000  O    (4g)
  -0.19890000000000    0.49450000000000    0.00400000000000  O    (4g)
  -0.19890000000000   -0.49450000000000   -0.49600000000000  O    (4g)
   0.19890000000000   -0.49450000000000    0.99600000000000  O    (4g)
   0.26120000000000    0.51440000000000    0.16320000000000  O    (4g)
  -0.26120000000000    0.51440000000000    0.33680000000000  O    (4g)
  -0.26120000000000   -0.51440000000000   -0.16320000000000  O    (4g)
   0.26120000000000   -0.51440000000000    0.66320000000000  O    (4g)
   0.32610000000000    0.48910000000000    0.32820000000000  O    (4g)
  -0.32610000000000    0.48910000000000    0.17180000000000  O    (4g)
  -0.32610000000000   -0.48910000000000   -0.32820000000000  O    (4g)
   0.32610000000000   -0.48910000000000    0.82820000000000  O    (4g)
   0.41130000000000    0.50150000000000    0.00450000000000  O    (4g)
  -0.41130000000000    0.50150000000000    0.49550000000000  O    (4g)
  -0.41130000000000   -0.50150000000000   -0.00450000000000  O    (4g)
   0.41130000000000   -0.50150000000000    0.50450000000000  O    (4g)
   0.45700000000000    0.49080000000000    0.15320000000000  O    (4g)
  -0.45700000000000    0.49080000000000    0.34680000000000  O    (4g)
  -0.45700000000000   -0.49080000000000   -0.15320000000000  O    (4g)
   0.45700000000000   -0.49080000000000    0.65320000000000  O    (4g)
```

Huanzalaite ($MgWO_4$, $H0_6$): AB4C_mP12_13_f_2g_e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Huanzalaite'
_chemical_formula_sum 'Mg O4 W'

loop_
_publ_author_name
 'V. B. Kravchenko'
_journal_name_full_name
;
 Journal of Structural Chemistry
;
_journal_volume 10
_journal_year 1969
_journal_page_first 139
_journal_page_last 140
_publ_Section_title
;
 Crystal structure of the monoclinic form of magnesium tungstate MgWO$_{
    ↪ 4}$
;

_aflow_title 'Huanzalaite ($MgWO$_{4}$, $H0_{6}$) Structure'
_aflow_proto 'AB4C_mP12_13_f_2g_e'
_aflow_params 'a,b/a,c/a,\beta,y_{1},y_{2},x_{3},y_{3},z_{3},x_{4},y_{4
    ↪ },z_{4}'
_aflow_params_values '4.68,1.2094017094,1.05128205128,90.33333,0.1818,
    ↪ 0.6592,0.221,0.101,-0.07,0.266,0.384,0.391'
_aflow_Strukturbericht '$H0_{6}$'
_aflow_Pearson 'mP12'

_symmetry_space_group_name_H-M "P 1 2/c 1"
_symmetry_Int_Tables_number 13

_cell_length_a    4.68000
_cell_length_b    5.66000
_cell_length_c    4.92000
_cell_angle_alpha 90.00000
_cell_angle_beta  90.33333
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
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2 -x,y,-z+1/2
3 -x,-y,-z
4 x,-y,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
W1 W 2 e 0.00000 0.18180 0.25000 1.00000
Mg1 Mg 2 f 0.50000 0.65920 0.25000 1.00000
O1 O 4 g 0.22100 0.10100 -0.07000 1.00000
O2 O 4 g 0.26600 0.38400 0.39100 1.00000
```

Huanzalaite ($MgWO_4$, $H0_6$): AB4C_mP12_13_f_2g_e - POSCAR

```
AB4C_mP12_13_f_2g_e & a,b/a,c/a,beta,y1,y2,x3,y3,z3,x4,y4,z4 --params=
    ↪ 4.68,1.2094017094,1.05128205128,90.33333,0.1818,0.6592,0.221,
    ↪ 0.101,-0.07,0.266,0.384,0.391 & P2/c C_{2h}^{4} #13 (efg^2) &
    ↪ mP12 & $H0_{6}$ & MgO4W & Huanzalaite & V. B. Kravchenko, J.
    ↪ Struct. Chem. 10, 139-140 (1969)
   1.00000000000000
   4.68000000000000    0.00000000000000    0.00000000000000
   0.00000000000000    5.66000000000000    0.00000000000000
  -0.02862295203742    0.00000000000000    4.91991673980533
     Mg      O      W
      2      8      2
Direct
   0.50000000000000    0.65920000000000    0.25000000000000  Mg   (2f)
   0.50000000000000   -0.65920000000000    0.75000000000000  Mg   (2f)
   0.22100000000000    0.10100000000000   -0.07000000000000  O    (4g)
  -0.22100000000000    0.10100000000000    0.57000000000000  O    (4g)
  -0.22100000000000   -0.10100000000000    0.07000000000000  O    (4g)
   0.22100000000000   -0.10100000000000    0.43000000000000  O    (4g)
   0.26600000000000    0.38400000000000    0.39100000000000  O    (4g)
  -0.26600000000000    0.38400000000000    0.10900000000000  O    (4g)
  -0.26600000000000   -0.38400000000000   -0.39100000000000  O    (4g)
   0.26600000000000   -0.38400000000000    0.89100000000000  O    (4g)
   0.00000000000000    0.18180000000000    0.25000000000000  W    (2e)
   0.00000000000000   -0.18180000000000    0.75000000000000  W    (2e)
```

$Cs_{11}O_3$: A11B3_mP56_14_11e_3e - CIF

```
# CIF file
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_audit_creation_method FINDSYM

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_chemical_formula_sum 'Cs11 O3'

loop_
_publ_author_name
 'A. Simon'
 'E. Westerbeck'
_journal_name_full_name
;
 Zeitschrift fur Anorganische und Allgemeine Chemie
;
_journal_volume 428
_journal_year 1977
_journal_page_first 187
_journal_page_last 198
_publ_Section_title
;
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```
 \"{U}ber Suboxide der Alkalimetalle. 10. Das ``komplexe Metall\''\' Cs$_
      ↪ {11}$O$_{3}$
;

# Found in Binary Alloy Phase Diagrams, 1990 Found in Binary Alloy Phase
      ↪  Diagrams, {Cd-Ce to Hf-Rb}}

_aflow_title 'Cs$_{11}$O$_{3}$ Structure'
_aflow_proto 'A11B3_mP56_14_11e_3e'
_aflow_params 'a,b/a,c/a,\beta,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3
      ↪ },y_{3},z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},
      ↪ z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{
      ↪ 10},y_{10},z_{10},x_{11},y_{11},z_{11},x_{12},y_{12},z_{12},x_{
      ↪ 13},y_{13},z_{13},x_{14},y_{14},z_{14}'
_aflow_params_values '17.61,0.523452583759,1.36553094832,100.14,0.3796,
      ↪ 0.7206,0.2438,0.092,-0.058,0.2834,0.2204,0.6157,0.3231,0.303,-
      ↪ 0.0205,0.3408,0.4356,0.4309,0.3821,0.5254,0.8324,0.4013,0.0663,
      ↪ 0.6974,0.4316,0.1617,0.1037,0.4538,0.3158,0.7371,0.4684,0.1517,
      ↪ 0.6716,0.1514,0.2388,0.4198,0.6692,0.2234,0.8414,0.2354,0.1795,
      ↪ 0.8554,0.3955,0.3835,0.7036,0.3654'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mP56'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

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_cell_length_b      9.21800
_cell_length_c      24.04700
_cell_angle_alpha   90.00000
_cell_angle_beta    100.14000
_cell_angle_gamma   90.00000

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_space_group_symop_operation_xyz
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2 -x,y+1/2,-z+1/2
3 -x,-y,-z
4 x,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
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Cs2 Cs 4 e 0.09200 -0.05800 0.28340 1.00000
Cs3 Cs 4 e 0.22040 0.61570 0.32310 1.00000
Cs4 Cs 4 e 0.30300 -0.02050 0.34080 1.00000
Cs5 Cs 4 e 0.43560 0.43090 0.38210 1.00000
Cs6 Cs 4 e 0.52540 0.83240 0.40130 1.00000
Cs7 Cs 4 e 0.06630 0.69740 0.43160 1.00000
Cs8 Cs 4 e 0.16170 0.10370 0.45380 1.00000
Cs9 Cs 4 e 0.31580 0.73710 0.46840 1.00000
Cs10 Cs 4 e 0.15170 0.67160 0.15140 1.00000
Cs11 Cs 4 e 0.23880 0.41980 0.66920 1.00000
O1 O 4 e 0.22340 0.84140 0.23540 1.00000
O2 O 4 e 0.17950 0.85540 0.39550 1.00000
O3 O 4 e 0.38350 0.70360 0.36540 1.00000
```

$Cs_{11}O_3$: A11B3_mP56_14_11e_3e - POSCAR

```
A11B3_mP56_14_11e_3e & a,b/a,c/a,beta,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,
      ↪ z4,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10,x11
      ↪ ,y11,z11,x12,y12,z12,x13,y13,z13,x14,y14,z14 --params=17.61,
      ↪ 0.523452583759,1.36553094832,100.14,0.3796,0.7206,0.2438,0.092
      ↪ ,-0.058,0.2834,0.2204,0.6157,0.3231,0.303,-0.0205,0.3408,0.4356
      ↪ ,0.4309,0.3821,0.5254,0.8324,0.4013,0.0663,0.6974,0.4316,0.1617
      ↪ ,0.1037,0.4538,0.3158,0.7371,0.4684,0.1517,0.6716,0.1514,0.2388
      ↪ ,0.4198,0.6692,0.2234,0.8414,0.2354,0.1795,0.8554,0.3955,0.3835
      ↪ ,0.7036,0.3654 & P2_{1}/c C_{2h}^{5} #14 (e^14) & mP56 & None &
      ↪  Cs11O3 & Cs11O3 & A. Simon and E. Westerbeck, Z. Anorg. Allg.
      ↪ Chem. 428, 187-198 (1977)
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  17.61000000000000    0.00000000000000    0.00000000000000
   0.00000000000000    9.21800000000000    0.00000000000000
  -4.23357044613628    0.00000000000000   23.67139814792530
     Cs      O
     44      12
Direct
   0.37960000000000    0.72060000000000    0.24380000000000   Cs   (4e)
  -0.37960000000000    1.22060000000000    0.25620000000000   Cs   (4e)
  -0.37960000000000   -0.72060000000000   -0.24380000000000   Cs   (4e)
   0.37960000000000   -0.22060000000000    0.74380000000000   Cs   (4e)
   0.09200000000000   -0.05800000000000    0.28340000000000   Cs   (4e)
  -0.09200000000000    0.44200000000000    0.21660000000000   Cs   (4e)
  -0.09200000000000    0.05800000000000   -0.28340000000000   Cs   (4e)
   0.09200000000000    0.55800000000000    0.78340000000000   Cs   (4e)
   0.22040000000000    0.61570000000000    0.32310000000000   Cs   (4e)
  -0.22040000000000    1.11570000000000    0.17690000000000   Cs   (4e)
  -0.22040000000000   -0.61570000000000   -0.32310000000000   Cs   (4e)
   0.22040000000000   -0.11570000000000    0.82310000000000   Cs   (4e)
   0.30300000000000   -0.02050000000000    0.34080000000000   Cs   (4e)
  -0.30300000000000    0.47950000000000    0.15920000000000   Cs   (4e)
  -0.30300000000000    0.02050000000000   -0.34080000000000   Cs   (4e)
   0.30300000000000    0.52050000000000    0.84080000000000   Cs   (4e)
   0.43560000000000    0.43090000000000    0.38210000000000   Cs   (4e)
  -0.43560000000000    0.93090000000000    0.11790000000000   Cs   (4e)
  -0.43560000000000   -0.43090000000000   -0.38210000000000   Cs   (4e)
   0.43560000000000    0.06910000000000    0.88210000000000   Cs   (4e)
   0.52540000000000    0.83240000000000    0.40130000000000   Cs   (4e)
  -0.52540000000000    1.33240000000000    0.09870000000000   Cs   (4e)
  -0.52540000000000   -0.83240000000000   -0.40130000000000   Cs   (4e)
   0.52540000000000   -0.33240000000000    0.90130000000000   Cs   (4e)
   0.06630000000000    0.69740000000000    0.43160000000000   Cs   (4e)
  -0.06630000000000    1.19740000000000    0.06840000000000   Cs   (4e)
  -0.06630000000000   -0.69740000000000   -0.43160000000000   Cs   (4e)
   0.06630000000000   -0.19740000000000    0.93160000000000   Cs   (4e)
   0.16170000000000    0.10370000000000    0.45380000000000   Cs   (4e)
  -0.16170000000000    0.60370000000000    0.04620000000000   Cs   (4e)
  -0.16170000000000   -0.10370000000000   -0.45380000000000   Cs   (4e)
   0.16170000000000    0.39630000000000    0.95380000000000   Cs   (4e)
   0.31580000000000    0.73710000000000    0.46840000000000   Cs   (4e)
  -0.31580000000000    1.23710000000000    0.03160000000000   Cs   (4e)
  -0.31580000000000   -0.73710000000000   -0.46840000000000   Cs   (4e)
   0.31580000000000   -0.23710000000000    0.96840000000000   Cs   (4e)
   0.15170000000000    0.67160000000000    0.15140000000000   Cs   (4e)
  -0.15170000000000    1.17160000000000    0.34860000000000   Cs   (4e)
  -0.15170000000000   -0.67160000000000   -0.15140000000000   Cs   (4e)
   0.15170000000000   -0.17160000000000    0.65140000000000   Cs   (4e)
   0.23880000000000    0.41980000000000    0.66920000000000   Cs   (4e)
  -0.23880000000000    0.91980000000000   -0.16920000000000   Cs   (4e)
  -0.23880000000000   -0.41980000000000   -0.66920000000000   Cs   (4e)
   0.23880000000000    0.08020000000000    1.16920000000000   Cs   (4e)
   0.22340000000000    0.84140000000000    0.23540000000000   O    (4e)
  -0.22340000000000    1.34140000000000    0.26460000000000   O    (4e)
  -0.22340000000000   -0.84140000000000   -0.23540000000000   O    (4e)
   0.22340000000000   -0.34140000000000    0.73540000000000   O    (4e)
   0.17950000000000    0.85540000000000    0.39550000000000   O    (4e)
  -0.17950000000000    1.35540000000000    0.10450000000000   O    (4e)
  -0.17950000000000   -0.85540000000000   -0.39550000000000   O    (4e)
   0.17950000000000   -0.35540000000000    0.89550000000000   O    (4e)
   0.38350000000000    0.70360000000000    0.36540000000000   O    (4e)
  -0.38350000000000    1.20360000000000    0.13460000000000   O    (4e)
  -0.38350000000000   -0.70360000000000   -0.36540000000000   O    (4e)
   0.38350000000000   -0.20360000000000    0.86540000000000   O    (4e)
```

Azurite [$Cu_3(CO_3)_2(OH)_2$, $G7_4$]: A2B3C2D8_mP30_14_e_ce_e_4e - CIF

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_audit_creation_method FINDSYM

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_chemical_formula_sum 'C2 Cu3 H2 O8'

loop_
_publ_author_name
 'K. C. Rule'
 'M. Reehuis'
 'M. C. R. Gibson'
 'B. Ouladdiaf'
 'M. J. Gutmann'
 'J.-U. Hoffmann'
 'S. Gerischer'
 'D. A. Tennant'
 'S. S\"{u}llow'
 'M. Lang'
_journal_name_full_name
;
 Physical Review B
;
_journal_volume 83
_journal_year 2011
_journal_page_first 104401
_journal_page_last 104401
_publ_Section_title
;
 Magnetic and crystal structure of azurite Cu$_{3}$(CO$_{3}$)$_{2}$(OH)
      ↪ $_{2}$ as determined by neutron diffraction
;

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      ↪ Structure'
_aflow_proto 'A2B3C2D8_mP30_14_e_ce_e_4e'
_aflow_params 'a,b/a,c/a,\beta,x_{2},y_{2},z_{2},x_{3},y_{3},z_{3},x_{4
      ↪ },y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{6},x_{7},y_{7},
      ↪ z_{7},x_{8},y_{8},z_{8}'
_aflow_params_values '4.99995,1.16452364524,2.06746667467,92.2103,0.3308
      ↪ ,0.7994,0.3192,0.2508,-0.0033,0.0834,0.182,0.3,0.3709,0.0975,
      ↪ 0.8972,0.3318,0.0762,0.3126,0.4451,0.4518,0.7098,0.4183,0.4339,
      ↪ 0.7949,0.2065'
_aflow_Strukturbericht '$G7_{4}$'
_aflow_Pearson 'mP30'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

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_cell_angle_beta    92.21030
_cell_angle_gamma   90.00000

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_space_group_symop_operation_xyz
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2 -x,y+1/2,-z+1/2
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4 x,-y+1/2,z+1/2

loop_
_atom_site_label
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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
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C1 C 4 e 0.33080 0.79940 0.31920 1.00000
Cu2 Cu 4 e 0.25080 -0.00330 0.08340 1.00000
H1 H 4 e 0.18200 0.30000 0.37090 1.00000
O1 O 4 e 0.09750 0.89720 0.33180 1.00000
O2 O 4 e 0.07620 0.31260 0.44510 1.00000
O3 O 4 e 0.45180 0.70980 0.41830 1.00000
O4 O 4 e 0.43390 0.79490 0.20650 1.00000
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Azurite [$Cu_3(CO_3)_2(OH)_2$, $G7_4$]: A2B3C2D8_mP30_14_e_ce_e_4e - POSCAR

```
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    ↪ x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8 --params=4.99995,
    ↪ 1.16452364524,2.06746667467,92.2103,0.3308,0.7994,0.3192,0.2508
    ↪ ,-0.0033,0.0834,0.182,0.3,0.3709,0.0975,0.8972,0.3318,0.0762,
    ↪ 0.3126,0.4451,0.4518,0.7098,0.4183,0.4339,0.7949,0.2065 & P2_{1
    ↪ }/c C_{2h}^{5} #14 (ce^7) & mP30 & $G7_{4}$ & C2Cu3H2O8 &
    ↪ Azurite & K. C. Rule et al., Phys. Rev. B 83, 104401(2011)
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   4.99995000000000    0.00000000000000    0.00000000000000
   0.00000000000000    5.82256000000000    0.00000000000000
  -0.39868054818008    0.00000000000000   10.32953909394810
     C     Cu     H     O
     4      6      4     16
Direct
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  -0.33080000000000   -0.79940000000000   -0.31920000000000    C    (4e)
   0.33080000000000   -0.29940000000000    0.81920000000000    C    (4e)
   0.00000000000000    0.00000000000000    0.50000000000000   Cu    (2c)
   0.00000000000000    0.50000000000000    0.00000000000000   Cu    (2c)
   0.25080000000000   -0.00330000000000    0.08340000000000   Cu    (4e)
  -0.25080000000000    0.49670000000000    0.41660000000000   Cu    (4e)
  -0.25080000000000    0.00330000000000   -0.08340000000000   Cu    (4e)
   0.25080000000000    0.50330000000000    0.58340000000000   Cu    (4e)
   0.18200000000000    0.30000000000000    0.37090000000000    H    (4e)
  -0.18200000000000    0.80000000000000    0.12910000000000    H    (4e)
  -0.18200000000000   -0.30000000000000   -0.37090000000000    H    (4e)
   0.18200000000000    0.20000000000000    0.87090000000000    H    (4e)
   0.09750000000000    0.89720000000000    0.33180000000000    O    (4e)
  -0.09750000000000    1.39720000000000    0.16820000000000    O    (4e)
  -0.09750000000000   -0.89720000000000   -0.33180000000000    O    (4e)
   0.09750000000000   -0.39720000000000    0.83180000000000    O    (4e)
   0.07620000000000    0.31260000000000    0.44510000000000    O    (4e)
  -0.07620000000000    0.81260000000000    0.05490000000000    O    (4e)
  -0.07620000000000   -0.31260000000000   -0.44510000000000    O    (4e)
   0.07620000000000    0.18740000000000    0.94510000000000    O    (4e)
   0.45180000000000    0.70980000000000    0.41830000000000    O    (4e)
  -0.45180000000000    1.20980000000000    0.08170000000000    O    (4e)
  -0.45180000000000   -0.70980000000000   -0.41830000000000    O    (4e)
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  -0.43390000000000    1.29490000000000    0.29350000000000    O    (4e)
  -0.43390000000000   -0.79490000000000   -0.20650000000000    O    (4e)
   0.43390000000000   -0.29490000000000    0.70650000000000    O    (4e)
```

$HgCl_2{\cdot}2HgO$: A2B3C2_mP14_14_e_ae_e - CIF

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_audit_creation_method FINDSYM

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_chemical_formula_sum 'Cl2 Hg3 O2'

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 'S. \v{S}\\'{c}avni\\'{c}ar'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 8
_journal_year 1955
_journal_page_first 379
_journal_page_last 383
_publ_Section_title
;
 The crystal structure of trimercuric oxychloride, HgCl$_{2}$$\cdot$2HgO
;

_aflow_title 'HgCl$_{2}$$\cdot$2HgO Structure'
_aflow_proto 'A2B3C2_mP14_14_e_ae_e'
_aflow_params 'a,b/a,c/a,\beta,x_{2},y_{2},z_{2},x_{3},y_{3},z_{3},x_{4
    ↪ },y_{4},z_{4}'
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    ↪ ,0.62,0.233,0.412,0.365,0.151,0.397,0.115,0.331'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mP14'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

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_cell_angle_beta  126.16667
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z+1/2
3 -x,-y,-z
4 x,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Hg1 Hg 2 a 0.00000 0.00000 0.00000 1.00000
Cl1 Cl 4 e 0.18500 0.62000 0.23300 1.00000
Hg2 Hg 4 e 0.41200 0.36500 0.15100 1.00000
O1 O 4 e 0.39700 0.11500 0.33100 1.00000
```

$HgCl_2{\cdot}2HgO$: A2B3C2_mP14_14_e_ae_e - POSCAR

```
A2B3C2_mP14_14_e_ae_e & a,b/a,c/a,beta,x2,y2,z2,x3,y3,z3,x4,y4,z4 --
    ↪ params=7.16,0.959497206704,0.958100558659,126.16667,0.185,0.62,
    ↪ 0.233,0.412,0.365,0.151,0.397,0.115,0.331 & P2_{1}/c C_{2h}^{5}
    ↪ #14 (ae^3) & mP14 & None & Cl2Hg3O2 & Cl2Hg3O2 & S. \v{S}\'{c}
    ↪ avni\'{c}ar, Acta Cryst. 8, 379-383 (1955)
   1.00000000000000
   7.16000000000000    0.00000000000000    0.00000000000000
   0.00000000000000    6.87000000000000    0.00000000000000
  -4.04833394910768    0.00000000000000    5.53810366790856
     Cl     Hg     O
      4      6      4
Direct
   0.18500000000000    0.62000000000000    0.23300000000000   Cl    (4e)
  -0.18500000000000    1.12000000000000    0.26700000000000   Cl    (4e)
  -0.18500000000000   -0.62000000000000   -0.23300000000000   Cl    (4e)
   0.18500000000000   -0.12000000000000    0.73300000000000   Cl    (4e)
   0.00000000000000    0.00000000000000    0.00000000000000   Hg    (2a)
   0.00000000000000    0.50000000000000    0.50000000000000   Hg    (2a)
   0.41200000000000    0.36500000000000    0.15100000000000   Hg    (4e)
  -0.41200000000000    0.86500000000000    0.34900000000000   Hg    (4e)
  -0.41200000000000   -0.36500000000000   -0.15100000000000   Hg    (4e)
   0.41200000000000    0.13500000000000    0.65100000000000   Hg    (4e)
   0.39700000000000    0.11500000000000    0.33100000000000    O    (4e)
  -0.39700000000000    0.61500000000000    0.16900000000000    O    (4e)
  -0.39700000000000   -0.11500000000000   -0.33100000000000    O    (4e)
   0.39700000000000    0.38500000000000    0.83100000000000    O    (4e)
```

Orpiment (As_2S_3, $D5_f$): A2B3_mP20_14_2e_3e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Orpiment'
_chemical_formula_sum 'As2 S3'

loop_
_publ_author_name
 'N. Morimoto'
_journal_name_full_name
;
 Mineralogical Journal
;
_journal_volume 1
_journal_year 1954
_journal_page_first 160
_journal_page_last 169
_publ_Section_title
;
 The Crystal Structure of Orpiment (As$_{2}$S$_{3}$) Refined
;

_aflow_title 'Orpiment (As$_{2}$S$_{3}$, $D5_{f}$) Structure'
_aflow_proto 'A2B3_mP20_14_2e_3e'
_aflow_params 'a,b/a,c/a,\beta,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3
    ↪ },y_{3},z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{5}'
_aflow_params_values '4.22,2.26777251185,2.88570616114,109.77475,0.124,
    ↪ 0.19,0.267,0.841,0.323,0.484,0.895,0.12,0.395,0.342,0.397,0.355
    ↪ ,0.715,0.293,0.125'
_aflow_Strukturbericht '$D5_{f}$'
_aflow_Pearson 'mP20'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

_cell_length_a    4.22000
_cell_length_b    9.57000
_cell_length_c    12.17768
_cell_angle_alpha 90.00000
_cell_angle_beta  109.77475
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z+1/2
3 -x,-y,-z
4 x,-y+1/2,z+1/2
```

```
loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
As1 As 4 e 0.12400 0.19000 0.26700 1.00000
As2 As 4 e 0.84100 0.32300 0.48400 1.00000
S1 S 4 e 0.89500 0.12000 0.39500 1.00000
S2 S 4 e 0.34200 0.39700 0.35500 1.00000
S3 S 4 e 0.71500 0.29300 0.12500 1.00000
```

Orpiment (As_2S_3, $D5_f$): A2B3_mP20_14_2e_3e - POSCAR

```
A2B3_mP20_14_2e_3e & a,b/a,c/a,beta,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,z4,
    ↪ x5,y5,z5 --params=4.22,2.26777251185,2.88570616114,109.77475,
    ↪ 0.124,0.19,0.267,0.841,0.323,0.484,0.895,0.12,0.395,0.342,0.397
    ↪ ,0.355,0.715,0.293,0.125 & P2_{1}/c C_{2h}^{5} #14 (e^5) & mP20
    ↪  & $D5_{f}$ & As2S3 & Orpiment & N. Morimoto, Mineral. J. 1,
    ↪ 160-169 (1954)
   1.00000000000000
   4.22000000000000    0.00000000000000    0.00000000000000
   0.00000000000000    9.57000000000000    0.00000000000000
  -4.11999221802126    0.00000000000000   11.45956169780700
     As      S
      8     12
Direct
   0.12400000000000    0.19000000000000    0.26700000000000   As   (4e)
  -0.12400000000000    0.69000000000000    0.23300000000000   As   (4e)
  -0.12400000000000   -0.19000000000000   -0.26700000000000   As   (4e)
   0.12400000000000    0.31000000000000    0.76700000000000   As   (4e)
   0.84100000000000    0.32300000000000    0.48400000000000   As   (4e)
  -0.84100000000000    0.82300000000000    0.01600000000000   As   (4e)
  -0.84100000000000   -0.32300000000000   -0.48400000000000   As   (4e)
   0.84100000000000    0.17700000000000    0.98400000000000   As   (4e)
   0.89500000000000    0.12000000000000    0.39500000000000    S   (4e)
  -0.89500000000000    0.62000000000000    0.10500000000000    S   (4e)
  -0.89500000000000   -0.12000000000000   -0.39500000000000    S   (4e)
   0.89500000000000    0.38000000000000    0.89500000000000    S   (4e)
   0.34200000000000    0.39700000000000    0.35500000000000    S   (4e)
  -0.34200000000000    0.89700000000000    0.14500000000000    S   (4e)
  -0.34200000000000   -0.39700000000000   -0.35500000000000    S   (4e)
   0.34200000000000    0.10300000000000    0.85500000000000    S   (4e)
   0.71500000000000    0.29300000000000    0.12500000000000    S   (4e)
  -0.71500000000000    0.79300000000000    0.37500000000000    S   (4e)
  -0.71500000000000   -0.29300000000000   -0.12500000000000    S   (4e)
   0.71500000000000    0.20700000000000    0.62500000000000    S   (4e)
```

Monoclinic Cu_2OSeO_3: A2B4C_mP28_14_abe_4e_e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Cu2O4Se'
_chemical_formula_sum 'Cu2 O4 Se'

loop_
_publ_author_name
 'H. Effenberger'
 'F. Pertlik'
_journal_name_full_name
;
 Monatshefte f{\"u}r Chemie - Chemical Monthly
;
_journal_volume 117
_journal_year 1986
_journal_page_first 887
_journal_page_last 896
_publ_Section_title
;
 Die Kristallstrukturen der Kupfer(II)-oxo-selenite Cu$_{2}$O(SeO$_{3}$)
    ↪  (kubisch und monoklin) und Cu$_{4}$O(SeO$_{3}$)$_{3}$ (
    ↪ monoklin und triklin)
;

# Found in Magnon spectrum of the helimagnetic insulator Cu$_{2}$OSeO$_{
    ↪ 3}$, 2016

_aflow_title 'Monoclinic Cu$_{2}$OSeO$_{3}$ Structure'
_aflow_proto 'A2B4C_mP28_14_abe_4e_e'
_aflow_params 'a,b/a,c/a,\beta,x_{3},y_{3},z_{3},x_{4},y_{4},z_{4},x_{5
    ↪ },y_{5},z_{5},x_{6},y_{6},z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},
    ↪ z_{8}'
_aflow_params_values '6.987,0.852010877344,1.53753112924,128.36609,
    ↪ 0.5082,0.705,0.7605,0.6549,-0.031,-0.0981,0.408,0.5641,0.0995,-
    ↪ 0.051,0.717,0.896,0.261,0.7826,0.8405,0.1649,0.5773,-0.0999'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mP28'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

_cell_length_a     6.98700
_cell_length_b     5.95300
_cell_length_c     10.74273
_cell_angle_alpha  90.00000
_cell_angle_beta   128.36609
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z+1/2
3 -x,-y,-z
4 x,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cu1 Cu 2 a 0.00000 0.00000 0.00000 1.00000
Cu2 Cu 2 b 0.50000 0.00000 0.00000 1.00000
Cu3 Cu 4 e 0.50820 0.70500 0.76050 1.00000
O1 O 4 e 0.65490 -0.03100 -0.09810 1.00000
O2 O 4 e 0.40800 0.56410 0.09950 1.00000
O3 O 4 e -0.05100 0.71700 0.89600 1.00000
O4 O 4 e 0.26100 0.78260 0.84050 1.00000
Se1 Se 4 e 0.16490 0.57730 -0.09990 1.00000
```

Monoclinic Cu_2OSeO_3: A2B4C_mP28_14_abe_4e_e - POSCAR

```
A2B4C_mP28_14_abe_4e_e & a,b/a,c/a,beta,x3,y3,z3,x4,y4,z4,x5,y5,z5,x6,y6
    ↪ ,z6,x7,y7,z7,x8,y8,z8 --params=6.987,0.852010877344,
    ↪ 1.53753112924,128.36609,0.5082,0.705,0.7605,0.6549,-0.031,-
    ↪ 0.0981,0.408,0.5641,0.0995,-0.051,0.717,0.896,0.261,0.7826,
    ↪ 0.8405,0.1649,0.5773,-0.0999 & P2_{1}/c C_{2h}^{5} #14 (abe^6)
    ↪ & mP28 & None & Cu2O4Se & Cu2O4Se & H. Effenberger and F.
    ↪ Pertlik, Monatshefte f{"u}r Chemie - Chemical Monthly 117,
    ↪ 887-896 (1986)
   1.00000000000000
   6.98700000000000    0.00000000000000    0.00000000000000
   0.00000000000000    5.95300000000000    0.00000000000000
  -6.66783901034905    0.00000000000000    8.42295499127043
     Cu      O     Se
      8     16      4
Direct
   0.00000000000000    0.00000000000000    0.00000000000000   Cu   (2a)
   0.00000000000000    0.50000000000000    0.50000000000000   Cu   (2a)
   0.50000000000000    0.00000000000000    0.00000000000000   Cu   (2b)
   0.50000000000000    0.50000000000000    0.50000000000000   Cu   (2b)
   0.50820000000000    0.70500000000000    0.76050000000000   Cu   (4e)
  -0.50820000000000    1.20500000000000   -0.26050000000000   Cu   (4e)
  -0.50820000000000   -0.70500000000000   -0.76050000000000   Cu   (4e)
   0.50820000000000   -0.20500000000000    1.26050000000000   Cu   (4e)
   0.65490000000000   -0.03100000000000   -0.09810000000000    O   (4e)
  -0.65490000000000    0.46900000000000    0.59810000000000    O   (4e)
  -0.65490000000000    0.03100000000000    0.09810000000000    O   (4e)
   0.65490000000000    0.53100000000000    0.40190000000000    O   (4e)
   0.40800000000000    0.56410000000000    0.09950000000000    O   (4e)
  -0.40800000000000    1.06410000000000    0.40050000000000    O   (4e)
  -0.40800000000000   -0.56410000000000   -0.09950000000000    O   (4e)
   0.40800000000000   -0.06410000000000    0.59950000000000    O   (4e)
  -0.05100000000000    0.71700000000000    0.89600000000000    O   (4e)
   0.05100000000000    1.21700000000000   -0.39600000000000    O   (4e)
   0.05100000000000   -0.71700000000000   -0.89600000000000    O   (4e)
  -0.05100000000000   -0.21700000000000    1.39600000000000    O   (4e)
   0.26100000000000    0.78260000000000    0.84050000000000    O   (4e)
  -0.26100000000000    1.28260000000000   -0.34050000000000    O   (4e)
  -0.26100000000000   -0.78260000000000   -0.84050000000000    O   (4e)
   0.26100000000000   -0.28260000000000    1.34050000000000    O   (4e)
   0.16490000000000    0.57730000000000   -0.09990000000000   Se   (4e)
  -0.16490000000000    1.07730000000000    0.59990000000000   Se   (4e)
  -0.16490000000000   -0.57730000000000    0.09990000000000   Se   (4e)
   0.16490000000000   -0.07730000000000    0.40010000000000   Se   (4e)
```

$Sb_4O_5Cl_2$: A2B5C4_mP22_14_e_c2e_2e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Cl2O5Sb4'
_chemical_formula_sum 'Cl2 O5 Sb4'

loop_
_publ_author_name
 'C. S\"{a}rnstrand'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 34
_journal_year 1978
_journal_page_first 2402
_journal_page_last 2407
_publ_Section_title
;
 The crystal structure of antimony(III) chloride oxide Sb$_{4}$O$_{5}
    ↪ $Cl$_{2}$
;

_aflow_title 'Sb$_{4}$O$_{5}$Cl$_{2}$ Structure'
_aflow_proto 'A2B5C4_mP22_14_e_c2e_2e'
_aflow_params 'a,b/a,c/a,\beta,x_{2},y_{2},z_{2},x_{3},y_{3},z_{3},x_{4
    ↪ },y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{6}'
_aflow_params_values '6.238,0.819365181148,2.170246874,97.217,0.51,0.703
    ↪ ,0.115,0.13,0.35,0.185,0.07,0.05,-0.085,0.186,0.225,0.049,0.796
    ↪ ,0.113,0.203'
_aflow_Strukturbericht 'None'
```

```
_aflow_Pearson 'mP22'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

_cell_length_a 6.23800
_cell_length_b 5.11120
_cell_length_c 13.53800
_cell_angle_alpha 90.00000
_cell_angle_beta 97.21700
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z+1/2
3 -x,-y,-z
4 x,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 2 c 0.00000 0.00000 0.50000 1.00000
Cl1 Cl 4 e 0.51000 0.70300 0.11500 1.00000
O2 O 4 e 0.13000 0.35000 0.18500 1.00000
O3 O 4 e 0.07000 0.05000 -0.08500 1.00000
Sb1 Sb 4 e 0.18600 0.22500 0.04900 1.00000
Sb2 Sb 4 e 0.79600 0.11300 0.20300 1.00000
```

$Sb_4O_5Cl_2$: A2B5C4_mP22_14_e_c2e_2e - POSCAR

```
A2B5C4_mP22_14_e_c2e_2e & a,b/a,c/a,beta,x2,y2,z2,x3,y3,z3,x4,y4,z4,x5,
    ↪ y5,z5,x6,y6,z6 --params=6.238,0.819365181148,2.170246874,97.217
    ↪ ,0.51,0.703,0.115,0.13,0.35,0.185,0.07,0.05,-0.085,0.186,0.225,
    ↪ 0.049,0.796,0.113,0.203 & P2_{1}/c C_{2h}^{5} #14 (ce^5) & mP22
    ↪ & None & Cl2O5Sb4 & Cl2O5Sb4 & C. S\"{a}rnstrand, Acta
    ↪ Crystallogr. Sect. B Struct. Sci. 34, 2402-2407 (1978)
   1.00000000000000
   6.23800000000000    0.00000000000000    0.00000000000000
   0.00000000000000    5.11120000000000    0.00000000000000
  -1.70074637299785    0.00000000000000   13.43074479597970
     Cl      O     Sb
      4     10      8
Direct
   0.51000000000000    0.70300000000000    0.11500000000000   Cl    (4e)
  -0.51000000000000    1.20300000000000    0.38500000000000   Cl    (4e)
  -0.51000000000000   -0.70300000000000   -0.11500000000000   Cl    (4e)
   0.51000000000000   -0.20300000000000    0.61500000000000   Cl    (4e)
   0.00000000000000    0.00000000000000    0.50000000000000   O     (2c)
   0.00000000000000    0.50000000000000    0.00000000000000   O     (2c)
   0.13000000000000    0.35000000000000    0.18500000000000   O     (4e)
  -0.13000000000000    0.85000000000000    0.31500000000000   O     (4e)
  -0.13000000000000   -0.35000000000000   -0.18500000000000   O     (4e)
   0.13000000000000    0.15000000000000    0.68500000000000   O     (4e)
   0.07000000000000    0.05000000000000   -0.08500000000000   O     (4e)
  -0.07000000000000    0.55000000000000    0.58500000000000   O     (4e)
  -0.07000000000000   -0.05000000000000    0.08500000000000   O     (4e)
   0.07000000000000    0.45000000000000    0.41500000000000   O     (4e)
   0.18600000000000    0.22500000000000    0.04900000000000   Sb    (4e)
  -0.18600000000000    0.72500000000000    0.45100000000000   Sb    (4e)
  -0.18600000000000   -0.22500000000000   -0.04900000000000   Sb    (4e)
   0.18600000000000    0.27500000000000    0.54900000000000   Sb    (4e)
   0.79600000000000    0.11300000000000    0.20300000000000   Sb    (4e)
  -0.79600000000000    0.61300000000000    0.29700000000000   Sb    (4e)
  -0.79600000000000   -0.11300000000000   -0.20300000000000   Sb    (4e)
   0.79600000000000    0.38700000000000    0.70300000000000   Sb    (4e)
```

Ca_2UO_5: A2B5C_mP32_14_2e_5e_ab - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Ca2O5U'
_chemical_formula_sum 'Ca2 O5 U'

loop_
_publ_author_name
 'B. O. Loopstra'
 'H. M. Rietveld'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 25
_journal_year 1969
_journal_page_first 787
_journal_page_last 791
_publ_Section_title
;
 The structure of some alkaline-earth metal uranates
;

_aflow_title 'Ca$_{2}$UO$_{5}$ Structure'
_aflow_proto 'A2B5C_mP32_14_2e_5e_ab'
_aflow_params 'a,b/a,c/a,\beta,x_{3},y_{3},z_{3},x_{4},y_{4},z_{4},x_{5
    ↪ },y_{5},z_{5},x_{6},y_{6},z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},
    ↪ z_{8},x_{9},y_{9},z_{9}'
_aflow_params_values '7.9137,0.687529221477,1.44663052681,108.803,0.1592
    ↪ ,0.0621,0.32,0.3462,-0.0015,0.6414,0.4447,0.3202,0.3267,0.8802,
    ↪ 0.2831,0.3532,0.6197,0.3043,0.072,0.0163,0.261,0.1295,0.2507,
    ↪ 0.1873,-0.0047'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mP32'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

_cell_length_a 7.91370
_cell_length_b 5.44090
_cell_length_c 11.44820
_cell_angle_alpha 90.00000
_cell_angle_beta 108.80300
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z+1/2
3 -x,-y,-z
4 x,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
U1 U 2 a 0.00000 0.00000 0.00000 1.00000
U2 U 2 b 0.50000 0.00000 0.00000 1.00000
Ca1 Ca 4 e 0.15920 0.06210 0.32000 1.00000
Ca2 Ca 4 e 0.34620 -0.00150 0.64140 1.00000
O1 O 4 e 0.44470 0.32020 0.32670 1.00000
O2 O 4 e 0.88020 0.28310 0.35320 1.00000
O3 O 4 e 0.61970 0.30430 0.07200 1.00000
O4 O 4 e 0.01630 0.26100 0.12950 1.00000
O5 O 4 e 0.25070 0.18730 -0.00470 1.00000
```

Ca_2UO_5: A2B5C_mP32_14_2e_5e_ab - POSCAR

```
A2B5C_mP32_14_2e_5e_ab & a,b/a,c/a,beta,x3,y3,z3,x4,y4,z4,x5,y5,z5,x6,y6
    ↪ ,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9 --params=7.9137,0.687529221477,
    ↪ 1.44663052681,108.803,0.1592,0.0621,0.32,0.3462,-0.0015,0.6414,
    ↪ 0.4447,0.3202,0.3267,0.8802,0.2831,0.3532,0.6197,0.3043,0.072,
    ↪ 0.0163,0.261,0.1295,0.2507,0.1873,-0.0047 & P2_{1}/c C_{2h}^{5}
    ↪ #14 (abe^7) & mP32 & None & Ca2O5U & Ca2O5U & B. O. Loopstra
    ↪ and H. M. Rietveld, Acta Crystallogr. Sect. B Struct. Sci. 25,
    ↪ 787-791 (1969)
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   7.91370000000000    0.00000000000000    0.00000000000000
   0.00000000000000    5.44090000000000    0.00000000000000
  -3.68992957359137    0.00000000000000   10.83723687025140
     Ca      O      U
      8     20      4
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  -0.15920000000000    0.56210000000000    0.18000000000000   Ca    (4e)
  -0.15920000000000   -0.06210000000000   -0.32000000000000   Ca    (4e)
   0.15920000000000    0.43790000000000    0.82000000000000   Ca    (4e)
   0.34620000000000   -0.00150000000000    0.64140000000000   Ca    (4e)
  -0.34620000000000    0.49850000000000   -0.14140000000000   Ca    (4e)
  -0.34620000000000    0.00150000000000   -0.64140000000000   Ca    (4e)
   0.34620000000000    0.50150000000000    1.14140000000000   Ca    (4e)
   0.44470000000000    0.32020000000000    0.32670000000000   O     (4e)
  -0.44470000000000    0.82020000000000    0.17330000000000   O     (4e)
  -0.44470000000000   -0.32020000000000   -0.32670000000000   O     (4e)
   0.44470000000000    0.17980000000000    0.82670000000000   O     (4e)
   0.88020000000000    0.28310000000000    0.35320000000000   O     (4e)
  -0.88020000000000    0.78310000000000    0.14680000000000   O     (4e)
  -0.88020000000000   -0.28310000000000   -0.35320000000000   O     (4e)
   0.88020000000000    0.21690000000000    0.85320000000000   O     (4e)
   0.61970000000000    0.30430000000000    0.07200000000000   O     (4e)
  -0.61970000000000    0.80430000000000    0.42800000000000   O     (4e)
  -0.61970000000000   -0.30430000000000   -0.07200000000000   O     (4e)
   0.61970000000000    0.19570000000000    0.57200000000000   O     (4e)
   0.01630000000000    0.26100000000000    0.12950000000000   O     (4e)
  -0.01630000000000    0.76100000000000    0.37050000000000   O     (4e)
  -0.01630000000000   -0.26100000000000   -0.12950000000000   O     (4e)
   0.01630000000000    0.23900000000000    0.62950000000000   O     (4e)
   0.25070000000000    0.18730000000000   -0.00470000000000   O     (4e)
  -0.25070000000000    0.68730000000000    0.50470000000000   O     (4e)
  -0.25070000000000   -0.18730000000000    0.00470000000000   O     (4e)
   0.25070000000000    0.31270000000000    0.49530000000000   O     (4e)
   0.00000000000000    0.00000000000000    0.00000000000000   U     (2a)
   0.00000000000000    0.50000000000000    0.50000000000000   U     (2a)
   0.50000000000000    0.00000000000000    0.00000000000000   U     (2b)
   0.50000000000000    0.50000000000000    0.50000000000000   U     (2b)
```

Gd_2SiO_5 (RE_2SiO_5 X1): A2B5C_mP32_14_2e_5e_e - CIF

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_audit_creation_method FINDSYM

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_chemical_formula_sum 'Gd2 O5 Si'

loop_
_publ_author_name
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  'G. V. Anan\'eva'
  'A. M. Korovkin'
  'T. I. Merkulyaeva'
  'A. M. Morozova'
  'M. V. Petrov'
  'I. R. Savinova'
  'V. R. Startsev'
  'P. P. Feofilov'
_journal_name_full_name
;
 Inorganic Materials
;
_journal_volume 17
_journal_year 1981
_journal_page_first 754
_journal_page_last 758
_publ_Section_title
;
 Growth of lanthanide oxyorthosilicate single crystals, and their
      ↪ structural and optical characteristics
;

# Found in Gd$_{2}$SiO$_{5}$ (Gd$_{2}$[SiO$_{4}$]O) Crystal Structure, {
      ↪ PAULING FILE in: Inorganic Solid Phases, SpringerMaterials (
      ↪ online database), Springer, Heidelberg (ed.) SpringerMaterials
      ↪ }, 2016

_aflow_title 'Gd$_{2}$SiO$_{5}$ ($RE$_{2}$SiO$_{5}$ X1) Structure'
_aflow_proto 'A2B5C_mP32_14_2e_5e_e'
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      ↪ },y_{3},z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},
      ↪ z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},z_{8}'
_aflow_params_values '9.16,0.770742358079,0.745633187773,107.58,0.02458,
      ↪ 0.12451,0.23428,0.38547,0.146,0.08372,0.1163,0.3782,0.4513,
      ↪ 0.2968,0.0698,0.3547,0.3683,0.4587,0.248,0.5941,0.2681,0.4507,
      ↪ 0.8839,0.3639,0.0059,0.298,0.5876,0.0402'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mP32'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

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_cell_length_c     6.83000
_cell_angle_alpha  90.00000
_cell_angle_beta   107.58000
_cell_angle_gamma  90.00000

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_space_group_symop_operation_xyz
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2 -x,y+1/2,-z+1/2
3 -x,-y,-z
4 x,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Gd1 Gd 4 e 0.02458 0.12451 0.23428 1.00000
Gd2 Gd 4 e 0.38547 0.14600 0.08372 1.00000
O1 O 4 e 0.11630 0.37820 0.45130 1.00000
O2 O 4 e 0.29680 0.06980 0.35470 1.00000
O3 O 4 e 0.36830 0.45870 0.24800 1.00000
O4 O 4 e 0.59410 0.26810 0.45070 1.00000
O5 O 4 e 0.88390 0.36390 0.00590 1.00000
Si1 Si 4 e 0.29800 0.58760 0.04020 1.00000
```

Gd_2SiO_5 (RE_2SiO_5 X1): A2B5C_mP32_14_2e_5e_e - POSCAR

```
A2B5C_mP32_14_2e_5e_e & a,b/a,c/a,beta,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,
      ↪ z4,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8 --params=9.16,
      ↪ 0.770742358079,0.745633187773,107.58,0.02458,0.12451,0.23428,
      ↪ 0.38547,0.146,0.08372,0.1163,0.3782,0.4513,0.2968,0.0698,0.3547
      ↪ ,0.3683,0.4587,0.248,0.5941,0.2681,0.4507,0.8839,0.3639,0.0059,
      ↪ 0.298,0.5876,0.0402 & P2_{1}/c C_{2h}^{5} #14 (e^8) & mP32 &
      ↪ None & Gd2O5Si & Gd2O5Si & G. V. Anan'eva et al., Inorg. Mat.
      ↪ 17, 754-758 (1981)
   1.00000000000000
   9.16000000000000    0.00000000000000    0.00000000000000
   0.00000000000000    7.06000000000000    0.00000000000000
  -2.06291370734890    0.00000000000000    6.51101275041234
     Gd      O     Si
      8     20      4
Direct
   0.02458000000000    0.12451000000000    0.23428000000000   Gd   (4e)
  -0.02458000000000    0.62451000000000    0.26572000000000   Gd   (4e)
  -0.02458000000000   -0.12451000000000   -0.23428000000000   Gd   (4e)
   0.02458000000000    0.37549000000000    0.73428000000000   Gd   (4e)
   0.38547000000000    0.14600000000000    0.08372000000000   Gd   (4e)
  -0.38547000000000    0.64600000000000    0.41628000000000   Gd   (4e)
  -0.38547000000000   -0.14600000000000   -0.08372000000000   Gd   (4e)
   0.38547000000000    0.35400000000000    0.58372000000000   Gd   (4e)
   0.11630000000000    0.37820000000000    0.45130000000000    O   (4e)
  -0.11630000000000    0.87820000000000    0.04870000000000    O   (4e)
  -0.11630000000000   -0.37820000000000   -0.45130000000000    O   (4e)
   0.11630000000000    0.12180000000000    0.95130000000000    O   (4e)
   0.29680000000000    0.06980000000000    0.35470000000000    O   (4e)
  -0.29680000000000    0.56980000000000    0.14530000000000    O   (4e)
  -0.29680000000000   -0.06980000000000   -0.35470000000000    O   (4e)
   0.29680000000000    0.43020000000000    0.85470000000000    O   (4e)
   0.36830000000000    0.45870000000000    0.24800000000000    O   (4e)
  -0.36830000000000    0.95870000000000    0.25200000000000    O   (4e)
  -0.36830000000000   -0.45870000000000   -0.24800000000000    O   (4e)
   0.36830000000000    0.04130000000000    0.74800000000000    O   (4e)
   0.59410000000000    0.26810000000000    0.45070000000000    O   (4e)
  -0.59410000000000    0.76810000000000    0.04930000000000    O   (4e)
  -0.59410000000000   -0.26810000000000   -0.45070000000000    O   (4e)
   0.59410000000000    0.23190000000000    0.95070000000000    O   (4e)
   0.88390000000000    0.36390000000000    0.00590000000000    O   (4e)
  -0.88390000000000    0.86390000000000    0.49410000000000    O   (4e)
  -0.88390000000000   -0.36390000000000   -0.00590000000000    O   (4e)
   0.88390000000000    0.13610000000000    0.50590000000000    O   (4e)
   0.29800000000000    0.58760000000000    0.04020000000000   Si   (4e)
  -0.29800000000000    1.08760000000000    0.45980000000000   Si   (4e)
  -0.29800000000000   -0.58760000000000   -0.04020000000000   Si   (4e)
   0.29800000000000   -0.08760000000000    0.54020000000000   Si   (4e)
```

Sanguite ($KCuCl_3$): A3BC_mP20_14_3e_e_e - CIF

```
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_audit_creation_method FINDSYM

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_chemical_formula_sum 'Cl3 Cu K'

loop_
_publ_author_name
 'R. D. Willett'
 'C. {Dwiggins, Jr.}'
 'R. F. Kruh'
 'R. E. Rundle'
_journal_name_full_name
;
 Journal of Chemical Physics
;
_journal_volume 38
_journal_year 1963
_journal_page_first 2429
_journal_page_last 2436
_publ_Section_title
;
 Crystal Structures of KCuCl$_{3}$ and NH$_{4}$CuCl$_{3}$
;

_aflow_title 'Sanguite (KCuCl$_{3}$) Structure'
_aflow_proto 'A3BC_mP20_14_3e_e_e'
_aflow_params 'a,b/a,c/a,\beta,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3
      ↪ },y_{3},z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{5}'
_aflow_params_values '4.029,3.42144452718,2.16827997022,97.33333,0.2754,
      ↪ 0.19875,0.263,0.6782,-0.00745,0.32171,0.8203,0.09875,-0.03369,
      ↪ 0.2408,0.04976,0.1575,0.7825,0.17081,0.55692'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mP20'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

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_cell_length_c     8.73600
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_cell_angle_beta   97.33333
_cell_angle_gamma  90.00000

loop_
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_space_group_symop_operation_xyz
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3 -x,-y,-z
4 x,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
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Cl2 Cl 4 e 0.67820 -0.00745 0.32171 1.00000
Cl3 Cl 4 e 0.82030 0.09875 -0.03369 1.00000
Cu1 Cu 4 e 0.24080 0.04976 0.15750 1.00000
K1 K 4 e 0.78250 0.17081 0.55692 1.00000
```

Sanguite ($KCuCl_3$): A3BC_mP20_14_3e_e_e - POSCAR

```
A3BC_mP20_14_3e_e_e & a,b/a,c/a,beta,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,z4
      ↪ ,x5,y5,z5 --params=4.029,3.42144452718,2.16827997022,97.33333,
      ↪ 0.2754,0.19875,0.263,0.6782,-0.00745,0.32171,0.8203,0.09875,-
      ↪ 0.03369,0.2408,0.04976,0.1575,0.7825,0.17081,0.55692 & P2_{1}/c
      ↪  C_{2h}^{5} #14 (e^5) & mP20 & None & Cl3CuK & Sanguite & R. D.
      ↪  Willett et al., J. Chem. Phys. 38, 2429-2436 (1963)
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   4.02900000000000    0.00000000000000    0.00000000000000
   0.00000000000000   13.78500000000000    0.00000000000000
  -1.11507693162750    0.00000000000000    8.66454265593702
     Cl     Cu      K
     12      4      4
```

```
Direct
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   -0.27540000000000   -0.19875000000000   -0.26300000000000   Cl   (4e)
    0.27540000000000    0.30125000000000    0.76300000000000   Cl   (4e)
    0.67820000000000   -0.00745000000000    0.32171000000000   Cl   (4e)
   -0.67820000000000    0.49255000000000    0.17829000000000   Cl   (4e)
   -0.67820000000000    0.00745000000000   -0.32171000000000   Cl   (4e)
    0.67820000000000    0.50745000000000    0.82171000000000   Cl   (4e)
    0.82030000000000    0.09875000000000   -0.03369000000000   Cl   (4e)
   -0.82030000000000    0.59875000000000    0.53369000000000   Cl   (4e)
   -0.82030000000000   -0.09875000000000    0.03369000000000   Cl   (4e)
    0.82030000000000    0.40125000000000    0.46631000000000   Cl   (4e)
    0.24080000000000    0.04976000000000    0.15750000000000   Cu   (4e)
   -0.24080000000000    0.54976000000000    0.34250000000000   Cu   (4e)
   -0.24080000000000   -0.04976000000000   -0.15750000000000   Cu   (4e)
    0.24080000000000    0.45024000000000    0.65750000000000   Cu   (4e)
    0.78250000000000    0.17081000000000    0.55692000000000   K    (4e)
   -0.78250000000000    0.67081000000000   -0.05692000000000   K    (4e)
   -0.78250000000000   -0.17081000000000   -0.55692000000000   K    (4e)
    0.78250000000000    0.32919000000000    1.05692000000000   K    (4e)
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γ-WO_3: A3B_mP32_14_6e_2e - CIF

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_chemical_formula_sum 'O3 W'

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 'A. W. Sleight'
 'T. Vogt'
_journal_name_full_name
;
 Journal of Solid State Chemistry
;
_journal_volume 131
_journal_year 1997
_journal_page_first 9
_journal_page_last 17
_publ_Section_title
;
 Ferroelectric Tungsten Trioxide
;

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_aflow_proto 'A3B_mP32_14_6e_2e'
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      ↪ },y_{3},z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},
      ↪ z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},z_{8}'
_aflow_params_values '7.3271,1.03238661954,1.44716463539,133.21984,0.782
      ↪ ,-0.03,0.782,0.781,0.536,0.779,0.005,0.736,0.723,0.472,0.743,
      ↪ 0.258,0.2768,-0.028,-0.0002,0.288,0.502,0.0,-0.03,-0.026,0.717,
      ↪ 0.4649,-0.033,0.2189'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mP32'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

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_cell_length_c    10.60352
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_cell_angle_beta   133.21984
_cell_angle_gamma  90.00000

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_space_group_symop_operation_xyz
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3 -x,-y,-z
4 x,-y+1/2,z+1/2

loop_
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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
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O2 O 4 e 0.78100 0.53600 0.77900 1.00000
O3 O 4 e 0.00500 0.73600 0.72300 1.00000
O4 O 4 e 0.47200 0.74300 0.25800 1.00000
O5 O 4 e 0.27680 -0.02800 -0.00020 1.00000
O6 O 4 e 0.28800 0.50200 0.00000 1.00000
W1 W 4 e -0.03000 -0.02600 0.71700 1.00000
W2 W 4 e 0.46490 -0.03300 0.21890 1.00000
```

γ-WO_3: A3B_mP32_14_6e_2e - POSCAR

```
A3B_mP32_14_6e_2e & a,b/a,c/a,beta,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,z4,
      ↪ x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8 --params=7.3271,
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      ↪ 0.536,0.779,0.005,0.736,0.723,0.472,0.743,0.258,0.2768,-0.028,-
      ↪ 0.0002,0.288,0.502,0.0,-0.03,-0.026,0.717,0.4649,-0.033,0.2189
      ↪ & P2_{1}/c C_{2h}^{5} #14 (e^8) & mP32 & None & O3W & O3W & P.
      ↪ M. Woodward and A. W. Sleight and T. Vogt, J. Solid State Chem.
      ↪ 131, 9-17 (1997)
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   7.32710000000000    0.00000000000000    0.00000000000000
   0.00000000000000    7.56440000000000    0.00000000000000
  -7.26128505928556    0.00000000000000    7.72711949423563
     O     W
    24     8
Direct
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   -0.78200000000000    0.47000000000000   -0.28200000000000   O   (4e)
   -0.78200000000000    0.03000000000000   -0.78200000000000   O   (4e)
    0.78200000000000    0.53000000000000    1.28200000000000   O   (4e)
    0.78100000000000    0.53600000000000    0.77900000000000   O   (4e)
   -0.78100000000000    1.03600000000000   -0.27900000000000   O   (4e)
   -0.78100000000000   -0.53600000000000   -0.77900000000000   O   (4e)
    0.78100000000000   -0.03600000000000    1.27900000000000   O   (4e)
    0.00500000000000    0.73600000000000    0.72300000000000   O   (4e)
   -0.00500000000000    1.23600000000000   -0.22300000000000   O   (4e)
   -0.00500000000000   -0.73600000000000   -0.72300000000000   O   (4e)
    0.00500000000000   -0.23600000000000    1.22300000000000   O   (4e)
    0.47200000000000    0.74300000000000    0.25800000000000   O   (4e)
   -0.47200000000000    1.24300000000000    0.24200000000000   O   (4e)
   -0.47200000000000   -0.74300000000000   -0.25800000000000   O   (4e)
    0.47200000000000   -0.24300000000000    0.75800000000000   O   (4e)
    0.27680000000000   -0.02800000000000   -0.00020000000000   O   (4e)
   -0.27680000000000    0.47200000000000    0.50020000000000   O   (4e)
   -0.27680000000000    0.02800000000000    0.00020000000000   O   (4e)
    0.27680000000000    0.52800000000000    0.49980000000000   O   (4e)
    0.28800000000000    0.50200000000000    0.00000000000000   O   (4e)
   -0.28800000000000    1.00200000000000    0.50000000000000   O   (4e)
   -0.28800000000000   -0.50200000000000    0.00000000000000   O   (4e)
    0.28800000000000   -0.00200000000000    0.50000000000000   O   (4e)
   -0.03000000000000   -0.02600000000000    0.71700000000000   W   (4e)
    0.03000000000000    0.47400000000000   -0.21700000000000   W   (4e)
    0.03000000000000    0.02600000000000   -0.71700000000000   W   (4e)
   -0.03000000000000    0.52600000000000    1.21700000000000   W   (4e)
    0.46490000000000   -0.03300000000000    0.21890000000000   W   (4e)
   -0.46490000000000    0.46700000000000    0.28110000000000   W   (4e)
   -0.46490000000000    0.03300000000000   -0.21890000000000   W   (4e)
    0.46490000000000    0.53300000000000    0.71890000000000   W   (4e)
```

$K_2Ni(CN)_4$: A4B2C4D_mP22_14_2e_e_2e_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'C4K2N4Ni'
_chemical_formula_sum 'C4 K2 N4 Ni'

loop_
_publ_author_name
 'N.-G. Vannerberg'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 18
_journal_year 1964
_journal_page_first 2385
_journal_page_last 2391
_publ_Section_title
;
 The Crystal Structure of K$_{2}$Ni(CN)$_{4}$
;

_aflow_title 'K$_{2}$Ni(CN)$_{4}$ Structure'
_aflow_proto 'A4B2C4D_mP22_14_2e_e_2e_a'
_aflow_params 'a,b/a,c/a,\beta,x_{2},y_{2},z_{2},x_{3},y_{3},z_{3},x_{4
      ↪ },y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{6}'
_aflow_params_values '4.294,1.78854215184,3.03446669772,87.26667,0.167,
      ↪ 0.207,0.055,0.849,0.382,0.39,0.281,0.046,0.341,0.261,0.338,
      ↪ 0.082,0.764,0.3,0.327'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mP22'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

_cell_length_a    4.29400
_cell_length_b    7.68000
_cell_length_c    13.03000
_cell_angle_alpha  90.00000
_cell_angle_beta   87.26667
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z+1/2
3 -x,-y,-z
4 x,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ni1 Ni 2 a 0.00000 0.00000 0.00000 1.00000
```

```
C1 C 4 e 0.16700 0.20700 0.05500 1.00000
C2 C 4 e 0.84900 0.38200 0.39000 1.00000
K1 K 4 e 0.28100 0.04600 0.34100 1.00000
N1 N 4 e 0.26100 0.33800 0.08200 1.00000
N2 N 4 e 0.76400 0.30000 0.32700 1.00000
```

$K_2Ni(CN)_4$: A4B2C4D_mP22_14_2e_e_2e_a - POSCAR

```
A4B2C4D_mP22_14_2e_e_2e_a & a,b/a,c/a,beta,x2,y2,z2,x3,y3,z3,x4,y4,z4,x5
    ↪ ,y5,z5,x6,y6,z6 --params=4.294,1.78854215184,3.03446669772,
    ↪ 87.26667,0.167,0.207,0.055,0.849,0.382,0.39,0.281,0.046,0.341,
    ↪ 0.261,0.338,0.082,0.764,0.3,0.327 & P2_{1}/c C_{2h}^{5} #14 (ae
    ↪ ^5) & mP22 & None & C4K2N4Ni & C4K2N4Ni & N.-G. Vannerberg,
    ↪ Acta Chem. Scand. 18, 2385-2391 (1964)
   1.00000000000000
   4.29400000000000   0.00000000000000   0.00000000000000
   0.00000000000000   7.68000000000000   0.00000000000000
   0.62136832261634   0.00000000000000  13.01517581163040
     C     K     N     Ni
     8     4     8     2
Direct
   0.16700000000000   0.20700000000000   0.05500000000000   C   (4e)
  -0.16700000000000   0.70700000000000   0.44500000000000   C   (4e)
  -0.16700000000000  -0.20700000000000  -0.05500000000000   C   (4e)
   0.16700000000000   0.29300000000000   0.55500000000000   C   (4e)
   0.84900000000000   0.38200000000000   0.39000000000000   C   (4e)
  -0.84900000000000   0.88200000000000   0.11000000000000   C   (4e)
  -0.84900000000000  -0.38200000000000  -0.39000000000000   C   (4e)
   0.84900000000000   0.11800000000000   0.89000000000000   C   (4e)
   0.28100000000000   0.04600000000000   0.34100000000000   K   (4e)
  -0.28100000000000   0.54600000000000   0.15900000000000   K   (4e)
  -0.28100000000000  -0.04600000000000  -0.34100000000000   K   (4e)
   0.28100000000000   0.45400000000000   0.84100000000000   K   (4e)
   0.26100000000000   0.33800000000000   0.08200000000000   N   (4e)
  -0.26100000000000   0.83800000000000   0.41800000000000   N   (4e)
  -0.26100000000000  -0.33800000000000  -0.08200000000000   N   (4e)
   0.26100000000000   0.16200000000000   0.58200000000000   N   (4e)
   0.76400000000000   0.30000000000000   0.32700000000000   N   (4e)
  -0.76400000000000   0.80000000000000   0.17300000000000   N   (4e)
  -0.76400000000000  -0.30000000000000  -0.32700000000000   N   (4e)
   0.76400000000000   0.20000000000000   0.82700000000000   N   (4e)
   0.00000000000000   0.00000000000000   0.00000000000000   Ni  (2a)
   0.00000000000000   0.50000000000000   0.50000000000000   Ni  (2a)
```

$KICl_4{\cdot}H_2O$ ($H0_{10}$): A4BCD_mP28_14_4e_e_e_e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Cl4IK \cdot H2O'
_chemical_formula_sum 'Cl4 (H2O) I K'

loop_
_publ_author_name
 'R. J. Elema'
 'J. L. {de Boer}'
 'A. Vos'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 16
_journal_year 1963
_journal_page_first 243
_journal_page_last 247
_publ_Section_title
;
 The refinement of the crystal structure of KICl$_{4}$ $\cdot$ H$_{2}$O
;

_aflow_title 'KICl$_{4}$$\cdot$H$_{2}$O ($H0_{10}$) Structure'
_aflow_proto 'A4BCD_mP28_14_4e_e_e_e'
_aflow_params 'a,b/a,c/a,\beta,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3
    ↪ },y_{3},z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},
    ↪ z_{6},x_{7},y_{7},z_{7}'
_aflow_params_values '4.284,3.34990662932,3.16137721755,102.62529,0.4857
    ↪ ,0.7807,0.2292,-0.06,0.0955,0.1866,0.625,-0.0116,0.3824,0.7875,
    ↪ 0.8953,0.0398,0.264,0.708,-0.043,0.7052,-0.0592,0.2041,0.1442,
    ↪ 0.3409,-0.0791'
_aflow_Strukturbericht '$H0_{10}$'
_aflow_Pearson 'mP28'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

_cell_length_a    4.28400
_cell_length_b    14.35100
_cell_length_c    13.54334
_cell_angle_alpha 90.00000
_cell_angle_beta  102.62529
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z+1/2
3 -x,-y,-z
4 x,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cl1 Cl 4 e 0.48570 0.78070 0.22920 1.00000
Cl2 Cl 4 e -0.06000 0.09550 0.18660 1.00000
Cl3 Cl 4 e 0.62500 -0.01160 0.38240 1.00000
Cl4 Cl 4 e 0.78750 0.89530 0.03980 1.00000
H2O1 H2O 4 e 0.26400 0.70800 -0.04300 1.00000
I1 I 4 e 0.70520 -0.05920 0.20410 1.00000
K1 K 4 e 0.14420 0.34090 -0.07910 1.00000
```

$KICl_4{\cdot}H_2O$ ($H0_{10}$): A4BCD_mP28_14_4e_e_e_e - POSCAR

```
A4BCD_mP28_14_4e_e_e_e & a,b/a,c/a,beta,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4
    ↪ ,z4,x5,y5,z5,x6,y6,z6,x7,y7,z7 --params=4.284,3.34990662932,
    ↪ 3.16137721755,102.62529,0.4857,0.7807,0.2292,-0.06,0.0955,
    ↪ 0.1866,0.625,-0.0116,0.3824,0.7875,0.8953,0.0398,0.264,0.708,-
    ↪ 0.043,0.7052,-0.0592,0.2041,0.1442,0.3409,-0.0791 & P2_{1}/c C_
    ↪ {2h}^{5} #14 (e^7) & mP28 & $H0_{10}$ & Cl4IK \cdot H2O & Cl4IK
    ↪  \cdot H2O & R. J. Elema and J. L. {de Boer} and A. Vos, Acta
    ↪ Cryst. 16, 243-247 (1963)
   1.00000000000000
   4.28400000000000   0.00000000000000   0.00000000000000
   0.00000000000000  14.35100000000000   0.00000000000000
  -2.96022177652463   0.00000000000000  13.21586718264790
     Cl    H2O     I     K
     16     4     4     4
Direct
   0.48570000000000   0.78070000000000   0.22920000000000   Cl   (4e)
  -0.48570000000000   1.28070000000000   0.27080000000000   Cl   (4e)
  -0.48570000000000  -0.78070000000000  -0.22920000000000   Cl   (4e)
   0.48570000000000  -0.28070000000000   0.72920000000000   Cl   (4e)
  -0.06000000000000   0.09550000000000   0.18660000000000   Cl   (4e)
   0.06000000000000   0.59550000000000   0.31340000000000   Cl   (4e)
   0.06000000000000  -0.09550000000000  -0.18660000000000   Cl   (4e)
  -0.06000000000000   0.40450000000000   0.68660000000000   Cl   (4e)
   0.62500000000000  -0.01160000000000   0.38240000000000   Cl   (4e)
  -0.62500000000000   0.48840000000000   0.11760000000000   Cl   (4e)
  -0.62500000000000   0.01160000000000  -0.38240000000000   Cl   (4e)
   0.62500000000000   0.51160000000000   0.88240000000000   Cl   (4e)
   0.78750000000000   0.89530000000000   0.03980000000000   Cl   (4e)
  -0.78750000000000   1.39530000000000   0.46020000000000   Cl   (4e)
  -0.78750000000000  -0.89530000000000  -0.03980000000000   Cl   (4e)
   0.78750000000000  -0.39530000000000   0.53980000000000   Cl   (4e)
   0.26400000000000   0.70800000000000  -0.04300000000000   H2O  (4e)
  -0.26400000000000   1.20800000000000   0.54300000000000   H2O  (4e)
  -0.26400000000000  -0.70800000000000   0.04300000000000   H2O  (4e)
   0.26400000000000  -0.20800000000000   0.45700000000000   H2O  (4e)
   0.70520000000000  -0.05920000000000   0.20410000000000   I    (4e)
  -0.70520000000000   0.44080000000000   0.29590000000000   I    (4e)
  -0.70520000000000   0.05920000000000  -0.20410000000000   I    (4e)
   0.70520000000000   0.55920000000000   0.70410000000000   I    (4e)
   0.14420000000000   0.34090000000000  -0.07910000000000   K    (4e)
  -0.14420000000000   0.84090000000000   0.57910000000000   K    (4e)
  -0.14420000000000  -0.34090000000000   0.07910000000000   K    (4e)
   0.14420000000000   0.15910000000000   0.42090000000000   K    (4e)
```

γ-$Y_2Si_2O_7$: A4BC_mP24_14_4e_e_e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'O7Si2Y2'
_chemical_formula_sum 'O4 Si Y'

loop_
_publ_author_name
 'A. N. Christensen'
 'R. G. Hazell'
 'A. W. Hewat'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 51
_journal_year 1997
_journal_page_first 37
_journal_page_last 43
_publ_Section_title
;
 Synthesis, Crystal Growth and Structure Investigations of Rare-Earth
    ↪ Disilicates and Rare-Earth Oxyapatites
;

# Found in Revision of the crystallographic data of polymorphic Y$_{2}
    ↪ $Si$_{2}$O$_{7}$ and Y$_{2}$SiO$_{5}$ compounds, 2004

_aflow_title '$\gamma$-Y$_{2}$Si$_{2}$O$_{7}$ Structure'
_aflow_proto 'A4BC_mP24_14_4e_e_e'
_aflow_params 'a,b/a,c/a,\beta,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3
    ↪ },y_{3},z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},
    ↪ z_{6}'
_aflow_params_values '4.6916,2.31309148265,1.1908943644,96.04,0.7933,
    ↪ 0.0502,0.1358,0.867,0.1822,0.5371,0.534,-0.0106,0.4976,0.3792,
    ↪ 0.2017,0.2513,0.6445,0.1136,0.3696,0.8904,0.8495,0.0941'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mP24'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

_cell_length_a    4.69160
```

```
_cell_length_b      10.85210
_cell_length_c      5.58720
_cell_angle_alpha   90.00000
_cell_angle_beta    96.04000
_cell_angle_gamma   90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z+1/2
3 -x,-y,-z
4 x,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 4 e 0.79330 0.05020 0.13580 1.00000
O2 O 4 e 0.86700 0.18220 0.53710 1.00000
O3 O 4 e 0.53400 -0.01060 0.49760 0.50000
O4 O 4 e 0.37920 0.20170 0.25130 1.00000
Si1 Si 4 e 0.64450 0.11360 0.36960 1.00000
Y1 Y 4 e 0.89040 0.84950 0.09410 1.00000
```

γ-$Y_2Si_2O_7$: A4BC_mP24_14_4e_e_e - POSCAR

```
A4BC_mP24_14_4e_e_e & a,b/a,c/a,beta,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,z4
    ↪ ,x5,y5,z5,x6,y6,z6 --params=4.6916,2.31309148265,1.1908943644,
    ↪ 96.04,0.7933,0.0502,0.1358,0.867,0.1822,0.5371,0.534,-0.0106,
    ↪ 0.4976,0.3792,0.2017,0.2513,0.6445,0.1136,0.3696,0.8904,0.8495,
    ↪ 0.0941 & P2_{1}/c C_{2h}^{5} #14 (e^6) & mP24 & None & O7Si2Y2
    ↪ & O7Si2Y2 & A. N. Christensen and R. G. Hazell and A. W. Hewat,
    ↪  Acta Chem. Scand. 51, 37-43 (1997)
   1.00000000000000
   4.69160000000000    0.00000000000000    0.00000000000000
   0.00000000000000   10.85210000000000    0.00000000000000
  -0.58790052086733    0.00000000000000    5.55618365585263
     O     Si     Y
    16      4      4
Direct
   0.79330000000000    0.05020000000000    0.13580000000000    O    (4e)
  -0.79330000000000    0.55020000000000    0.36420000000000    O    (4e)
  -0.79330000000000   -0.05020000000000   -0.13580000000000    O    (4e)
   0.79330000000000    0.44980000000000    0.63580000000000    O    (4e)
   0.86700000000000    0.18220000000000    0.53710000000000    O    (4e)
  -0.86700000000000    0.68220000000000   -0.03710000000000    O    (4e)
  -0.86700000000000   -0.18220000000000   -0.53710000000000    O    (4e)
   0.86700000000000    0.31780000000000    1.03710000000000    O    (4e)
   0.53400000000000   -0.01060000000000    0.49760000000000    O    (4e)
  -0.53400000000000    0.48940000000000    0.00240000000000    O    (4e)
  -0.53400000000000    0.01060000000000   -0.49760000000000    O    (4e)
   0.53400000000000    0.51060000000000    0.99760000000000    O    (4e)
   0.37920000000000    0.20170000000000    0.25130000000000    O    (4e)
  -0.37920000000000    0.70170000000000    0.24870000000000    O    (4e)
  -0.37920000000000   -0.20170000000000   -0.25130000000000    O    (4e)
   0.37920000000000    0.29830000000000    0.75130000000000    O    (4e)
   0.64450000000000    0.11360000000000    0.36960000000000   Si    (4e)
  -0.64450000000000    0.61360000000000    0.13040000000000   Si    (4e)
  -0.64450000000000   -0.11360000000000   -0.36960000000000   Si    (4e)
   0.64450000000000    0.38640000000000    0.86960000000000   Si    (4e)
   0.89040000000000    0.84950000000000    0.09410000000000    Y    (4e)
  -0.89040000000000    1.34950000000000    0.40590000000000    Y    (4e)
  -0.89040000000000   -0.84950000000000   -0.09410000000000    Y    (4e)
   0.89040000000000   -0.34950000000000    0.59410000000000    Y    (4e)
```

$K_2Pt(SCN)_6{\cdot}2H_2O$: A6B4C2D6E2FG6_mP54_14_3e_2e_e_3e_e_a_3e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'C6H4K2N6O2PtS6'
_chemical_formula_sum 'C6 H4 K2 N6 O2 Pt S6'

loop_
_publ_author_name
 'J. Arpalahti'
 'J. H\"{o}ls\"{a}'
 'R. Sillanp\"{a}\"{a}'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 47
_journal_year 1993
_journal_page_first 1078
_journal_page_last 1082
_publ_Section_title
;
 Studies on Potassium Thiocyanatoplatinates. II. Crystal Structure of
    ↪ Potassium Hexathiocyanatoplatinate(IV) Dihydrate, K$_{2}$Pt(
    ↪ SCN)$_{6}$ $\cdot$ 2H$_{2}$O
;

_aflow_title 'K$_{2}$Pt(SCN)$_{6}$$\cdot$2H$_{2}$O Structure'
_aflow_proto 'A6B4C2D6E2FG6_mP54_14_3e_2e_e_3e_e_a_3e'
_aflow_params 'a,b/a,c/a,\beta,x_{2},y_{2},z_{2},x_{3},y_{3},z_{3},x_{4
    ↪ },y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{6},x_{7},y_{7},
    ↪ z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{10},y_{10},z_{10},
    ↪ x_{11},y_{11},z_{11},x_{12},y_{12},z_{12},x_{13},y_{13},z_{13},
    ↪ x_{14},y_{14},z_{14}'
_aflow_params_values '11.33,0.983495145631,0.644571932921,94.76,0.7648,
    ↪ 0.8251,-0.0619,0.2402,-0.0953,0.8473,0.2213,-0.0589,0.3316,
    ↪ 0.4514,-0.0915,0.1233,0.5826,-0.0741,0.2021,0.4986,0.8473,
    ↪ 0.6064,0.6695,0.8204,-0.0303,0.327,0.8561,0.8963,0.2991,-0.0101
    ↪ ,0.4229,0.5087,-0.0876,0.2358,-0.0988,0.8216,0.8743,0.1151,-
    ↪ 0.0193,0.7552,0.1129,0.8611,0.2058'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mP54'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

_cell_length_a      11.33000
_cell_length_b      11.14300
_cell_length_c      7.30300
_cell_angle_alpha   90.00000
_cell_angle_beta    94.76000
_cell_angle_gamma   90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z+1/2
3 -x,-y,-z
4 x,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Pt1 Pt 2 a 0.00000 0.00000 0.00000 1.00000
C1 C 4 e 0.76480 0.82510 -0.06190 1.00000
C2 C 4 e 0.24020 -0.09530 0.84730 1.00000
C3 C 4 e 0.22130 -0.05890 0.33160 1.00000
H1 H 4 e 0.45140 -0.09150 0.12330 1.00000
H2 H 4 e 0.58260 -0.07410 0.20210 1.00000
K1 K 4 e 0.49860 0.84730 0.60640 1.00000
N1 N 4 e 0.66950 0.82040 -0.03030 1.00000
N2 N 4 e 0.32700 0.85610 0.89630 1.00000
N3 N 4 e 0.29910 -0.01010 0.42290 1.00000
O1 O 4 e 0.50870 -0.08760 0.23580 1.00000
S1 S 4 e -0.09880 0.82160 0.87430 1.00000
S2 S 4 e 0.11510 -0.01930 0.75520 1.00000
S3 S 4 e 0.11290 0.86110 0.20580 1.00000
```

$K_2Pt(SCN)_6{\cdot}2H_2O$: A6B4C2D6E2FG6_mP54_14_3e_2e_e_3e_e_a_3e - POSCAR

```
A6B4C2D6E2FG6_mP54_14_3e_2e_e_3e_e_a_3e & a,b/a,c/a,beta,x2,y2,z2,x3,y3,
    ↪ z3,x4,y4,z4,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,
    ↪ y10,z10,x11,y11,z11,x12,y12,z12,x13,y13,z13,x14,y14,z14 --
    ↪ params=11.33,0.983495145631,0.644571932921,94.76,0.7648,0.8251
    ↪ ,-0.0619,0.2402,-0.0953,0.8473,0.2213,-0.0589,0.3316,0.4514,-
    ↪ 0.0915,0.1233,0.5826,-0.0741,0.2021,0.4986,0.8473,0.6064,0.6695
    ↪ ,0.8204,-0.0303,0.327,0.8561,0.8963,0.2991,-0.0101,0.4229,
    ↪ 0.5087,-0.0876,0.2358,-0.0988,0.8216,0.8743,0.1151,-0.0193,
    ↪ 0.7552,0.1129,0.8611,0.2058 & P2_{1}/c C_{2h}^{5} #14 (ae^13) &
    ↪  mP54 & None & C6H4K2N6O2PtS6 & C6H4K2N6O2PtS6 & J. Arpalahti
    ↪ and J. H\"{o}ls\"{a} and R. Sillanp\"{a}\"{a}, Acta Chem.
    ↪ Scand. 47, 1078-1082 (1993)
   1.00000000000000
  11.33000000000000    0.00000000000000    0.00000000000000
   0.00000000000000   11.14300000000000    0.00000000000000
  -0.60601856653863    0.00000000000000    7.27781220539596
     C     H     K     N     O    Pt     S
    12     8     4     6     4     2    12
Direct
   0.76480000000000    0.82510000000000   -0.06190000000000    C    (4e)
  -0.76480000000000    1.32510000000000    0.56190000000000    C    (4e)
  -0.76480000000000   -0.82510000000000    0.06190000000000    C    (4e)
   0.76480000000000   -0.32510000000000    0.43810000000000    C    (4e)
   0.24020000000000   -0.09530000000000    0.84730000000000    C    (4e)
  -0.24020000000000    0.40470000000000   -0.34730000000000    C    (4e)
  -0.24020000000000    0.09530000000000   -0.84730000000000    C    (4e)
   0.24020000000000    0.59530000000000    1.34730000000000    C    (4e)
   0.22130000000000   -0.05890000000000    0.33160000000000    C    (4e)
  -0.22130000000000    0.44110000000000    0.16840000000000    C    (4e)
  -0.22130000000000    0.05890000000000   -0.33160000000000    C    (4e)
   0.22130000000000    0.55890000000000    0.83160000000000    C    (4e)
   0.45140000000000   -0.09150000000000    0.12330000000000    H    (4e)
  -0.45140000000000    0.40850000000000    0.37670000000000    H    (4e)
  -0.45140000000000    0.09150000000000   -0.12330000000000    H    (4e)
   0.45140000000000    0.59150000000000    0.62330000000000    H    (4e)
   0.58260000000000   -0.07410000000000    0.20210000000000    H    (4e)
  -0.58260000000000    0.42590000000000    0.29790000000000    H    (4e)
  -0.58260000000000    0.07410000000000   -0.20210000000000    H    (4e)
   0.58260000000000    0.57410000000000    0.70210000000000    H    (4e)
   0.49860000000000    0.84730000000000    0.60640000000000    K    (4e)
  -0.49860000000000    1.34730000000000   -0.10640000000000    K    (4e)
  -0.49860000000000   -0.84730000000000   -0.60640000000000    K    (4e)
   0.49860000000000   -0.34730000000000    1.10640000000000    K    (4e)
   0.66950000000000    0.82040000000000   -0.03030000000000    N    (4e)
  -0.66950000000000    1.32040000000000    0.53030000000000    N    (4e)
  -0.66950000000000   -0.82040000000000    0.03030000000000    N    (4e)
   0.66950000000000   -0.32040000000000    0.46970000000000    N    (4e)
   0.32700000000000    0.85610000000000    0.89630000000000    N    (4e)
  -0.32700000000000    1.35610000000000   -0.39630000000000    N    (4e)
  -0.32700000000000   -0.85610000000000   -0.89630000000000    N    (4e)
```

```
  0.32700000000000  −0.35610000000000   1.39630000000000   N   (4e)
  0.29910000000000  −0.01010000000000   0.42290000000000   N   (4e)
 −0.29910000000000   0.48990000000000   0.07710000000000   N   (4e)
 −0.29910000000000   0.01010000000000  −0.42290000000000   N   (4e)
  0.29910000000000   0.51010000000000   0.92290000000000   N   (4e)
  0.50870000000000  −0.08760000000000   0.23580000000000   O   (4e)
 −0.50870000000000   0.41240000000000   0.26420000000000   O   (4e)
 −0.50870000000000   0.08760000000000  −0.23580000000000   O   (4e)
  0.50870000000000   0.58760000000000   0.73580000000000   O   (4e)
  0.00000000000000   0.00000000000000   0.00000000000000   Pt  (2a)
  0.00000000000000   0.50000000000000   0.50000000000000   Pt  (2a)
 −0.09880000000000   0.82160000000000   0.87430000000000   S   (4e)
  0.09880000000000   1.32160000000000  −0.37430000000000   S   (4e)
  0.09880000000000  −0.82160000000000  −0.87430000000000   S   (4e)
 −0.09880000000000  −0.32160000000000   1.37430000000000   S   (4e)
  0.11510000000000  −0.01930000000000   0.75520000000000   S   (4e)
 −0.11510000000000   0.48070000000000  −0.25520000000000   S   (4e)
 −0.11510000000000   0.01930000000000  −0.75520000000000   S   (4e)
  0.11510000000000   0.51930000000000   1.25520000000000   S   (4e)
  0.11290000000000   0.86110000000000   0.20580000000000   S   (4e)
 −0.11290000000000   1.36110000000000   0.29420000000000   S   (4e)
 −0.11290000000000  −0.86110000000000  −0.20580000000000   S   (4e)
  0.11290000000000  −0.36110000000000   0.70580000000000   S   (4e)
```

K_2NbF_7 ($K6_2$): A7B2C_mP40_14_7e_2e_e - CIF

```
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_chemical_formula_sum 'F7 K2 Nb'

loop_
_publ_author_name
 'G. M. Brown'
 'L. A. Walker'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 20
_journal_year 1966
_journal_page_first 220
_journal_page_last 229
_publ_Section_title
;
 Refinement of the structure of potassium heptafluoroniobate, K$_{2}
      ↪ $NbF$_{7}$, from neutron-diffraction data
;

# Found in Structure and luminescence of K$_{2}$TaF$_{7}$ and K$_{2}
      ↪ $NbF$_{7}$, 1987

_aflow_title 'K$_{2}$NbF$_{7}$ ($K6_{2}$) Structure'
_aflow_proto 'A7B2C_mP40_14_7e_2e_e'
_aflow_params 'a,b/a,c/a,\beta,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3
      ↪ },y_{3},z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},
      ↪ z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{
      ↪ 10},y_{10},z_{10}'
_aflow_params_values '5.846,2.17122819022,1.4565514882,90.0,0.0463,
      ↪ 0.2431,0.1994,0.4573,0.2567,0.2114,0.0448,0.1072,0.388,0.4599,
      ↪ 0.1107,0.4101,0.2166,-0.0207,0.2159,0.1875,0.1216,0.0063,0.5572
      ↪ ,0.0792,0.1317,0.2397,0.4404,0.1833,0.7626,0.2845,0.4446,0.2718
      ↪ ,0.1288,0.2229'
_aflow_Strukturbericht '$K6_{2}$'
_aflow_Pearson 'mP40'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

_cell_length_a    5.84600
_cell_length_b    12.69300
_cell_length_c    8.51500
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

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_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z+1/2
3 -x,-y,-z
4 x,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
F1 F 4 e 0.04630 0.24310 0.19940 1.00000
F2 F 4 e 0.45730 0.25670 0.21140 1.00000
F3 F 4 e 0.04480 0.10720 0.38800 1.00000
F4 F 4 e 0.45990 0.11070 0.41010 1.00000
F5 F 4 e 0.21660 -0.02070 0.21590 1.00000
F6 F 4 e 0.18750 0.12160 0.00630 1.00000
F7 F 4 e 0.55720 0.07920 0.13170 1.00000
K1 K 4 e 0.23970 0.44040 0.18330 1.00000
K2 K 4 e 0.76260 0.28450 0.44460 1.00000
Nb1 Nb 4 e 0.27180 0.12880 0.22290 1.00000
```

K_2NbF_7 ($K6_2$): A7B2C_mP40_14_7e_2e_e - POSCAR

```
A7B2C_mP40_14_7e_2e_e & a,b/a,c/a,beta,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,
      ↪ z4,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10 --
      ↪ params=5.846,2.17122819022,1.4565514882,90.0,0.0463,0.2431,
      ↪ 0.1994,0.4573,0.2567,0.2114,0.0448,0.1072,0.388,0.4599,0.1107,
      ↪ 0.4101,0.2166,-0.0207,0.2159,0.1875,0.1216,0.0063,0.5572,0.0792
      ↪ ,0.1317,0.2397,0.4404,0.1833,0.7626,0.2845,0.4446,0.2718,0.1288
      ↪ ,0.2229 & P2_{1}/c C_{2h}^{5} #14 (e^10) & mP40 & $K6_{2}$ &
      ↪ F7K2Nb & F7K2Nb & G. M. Brown and L. A. Walker, Acta Cryst. 20,
      ↪ 220-229 (1966)
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   5.84600000000000   0.00000000000000   0.00000000000000
   0.00000000000000  12.69300000000000   0.00000000000000
   0.00000000000000   0.00000000000000   8.51500000000000
     F     K    Nb
    28     8     4
Direct
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  -0.04630000000000   0.74310000000000   0.30060000000000   F   (4e)
  -0.04630000000000  -0.24310000000000  -0.19940000000000   F   (4e)
   0.04630000000000   0.25690000000000   0.69940000000000   F   (4e)
   0.45730000000000   0.25670000000000   0.21140000000000   F   (4e)
  -0.45730000000000   0.75670000000000   0.28860000000000   F   (4e)
  -0.45730000000000  -0.25670000000000  -0.21140000000000   F   (4e)
   0.45730000000000   0.24330000000000   0.71140000000000   F   (4e)
   0.04480000000000   0.10720000000000   0.38800000000000   F   (4e)
  -0.04480000000000   0.60720000000000   0.11200000000000   F   (4e)
  -0.04480000000000  -0.10720000000000  -0.38800000000000   F   (4e)
   0.04480000000000   0.39280000000000   0.88800000000000   F   (4e)
   0.45990000000000   0.11070000000000   0.41010000000000   F   (4e)
  -0.45990000000000   0.61070000000000   0.08990000000000   F   (4e)
  -0.45990000000000  -0.11070000000000  -0.41010000000000   F   (4e)
   0.45990000000000   0.38930000000000   0.91010000000000   F   (4e)
   0.21660000000000  -0.02070000000000   0.21590000000000   F   (4e)
  -0.21660000000000   0.47930000000000   0.28410000000000   F   (4e)
  -0.21660000000000   0.02070000000000  -0.21590000000000   F   (4e)
   0.21660000000000   0.52070000000000   0.71590000000000   F   (4e)
   0.18750000000000   0.12160000000000   0.00630000000000   F   (4e)
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  -0.18750000000000  -0.12160000000000  -0.00630000000000   F   (4e)
   0.18750000000000   0.37840000000000   0.50630000000000   F   (4e)
   0.55720000000000   0.07920000000000   0.13170000000000   F   (4e)
  -0.55720000000000   0.57920000000000   0.36830000000000   F   (4e)
  -0.55720000000000  -0.07920000000000  -0.13170000000000   F   (4e)
   0.55720000000000   0.42080000000000   0.63170000000000   F   (4e)
   0.23970000000000   0.44040000000000   0.18330000000000   K   (4e)
  -0.23970000000000   0.94040000000000   0.31670000000000   K   (4e)
  -0.23970000000000  -0.44040000000000  -0.18330000000000   K   (4e)
   0.23970000000000   0.05960000000000   0.68330000000000   K   (4e)
   0.76260000000000   0.28450000000000   0.44460000000000   K   (4e)
  -0.76260000000000   0.78450000000000   0.05540000000000   K   (4e)
  -0.76260000000000  -0.28450000000000  -0.44460000000000   K   (4e)
   0.76260000000000   0.21550000000000   0.94460000000000   K   (4e)
   0.27180000000000   0.12880000000000   0.22290000000000   Nb  (4e)
  -0.27180000000000   0.62880000000000   0.27710000000000   Nb  (4e)
  -0.27180000000000  -0.12880000000000  -0.22290000000000   Nb  (4e)
   0.27180000000000   0.37120000000000   0.72290000000000   Nb  (4e)
```

Manganese-leonite 110 K [$K_2Mn(SO_4)_2{\cdot}4H_2O$]: A8B2CD12E2_mP100_14_8e_2e_ad_12e_2e - CIF

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_chemical_formula_sum 'H8 K2 Mn O12 S2'

loop_
_publ_author_name
 'B. Hertweck'
 'G. Giester'
 'E. Libowitzky'
_journal_name_full_name
;
 American Mineralogist
;
_journal_volume 86
_journal_year 2001
_journal_page_first 1282
_journal_page_last 1292
_publ_Section_title
;
 The crystal structures of the low-temperature phases of leonite-type
      ↪ compounds, K$_{2}$$Me$(SO$_{4}$)$_{2}$$\cdot$4H$_{2}$O ($Me^{2
      ↪ +}$ = Mg, Mn, Fe)
;

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      ↪ z_{8},x_{9},y_{9},z_{9},x_{10},y_{10},z_{10},x_{11},y_{11},z_{
      ↪ 11},x_{12},y_{12},z_{12},x_{13},y_{13},z_{13},x_{14},y_{14},z_{
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      ↪ 20},x_{21},y_{21},z_{21},x_{22},y_{22},z_{22},x_{23},y_{23},z_{
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      ↪ 26}'
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      ↪ 0.081,0.025,0.72,-0.049,-0.037,0.726,0.043,0.733,-0.008,0.06,
      ↪ 0.729,0.014,-0.044,0.29373,0.73232,0.16965,0.25852,0.25782,
      ↪ 0.16992,0.1163,0.6276,0.8466,0.8888,0.6261,0.1633,0.5998,-
```

```
      ↪ 0.0908 , 0.1493 , 0.5455 , 0.1528 , 0.1791 , 0.0611 , −0.0127 , 0.1725 , 0.0973
      ↪ , 0.4938 , 0.1853 , 0.4342 , 0.4911 , 0.1674 , 0.3741 , −0.0227 , 0.1931 ,
      ↪ 0.3537 , 0.6684 , −0.0342 , 0.6484 , 0.6653 , 0.0431 , −0.0131 , 0.7721 , −
      ↪ 0.0051 , 0.7802 , 0.0103 , 0.0111 , −0.04519 , 0.49523 , 0.20732 , 0.51973 ,
      ↪ 0.00747 , 0.21252 '
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mP100'

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_symmetry_Int_Tables_number 14

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_cell_angle_beta 95.02000
_cell_angle_gamma 90.00000

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_space_group_symop_operation_xyz
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3 -x,-y,-z
4 x,-y+1/2,z+1/2

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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
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Mn2 Mn 2 d 0.50000 0.00000 0.50000 1.00000
H1 H 4 e 0.38700 0.72900 -0.07000 1.00000
H2 H 4 e 0.27500 0.64900 -0.07400 1.00000
H3 H 4 e 0.62000 0.73500 0.07400 1.00000
H4 H 4 e 0.72100 0.64800 0.08100 1.00000
H5 H 4 e 0.02500 0.72000 -0.04900 1.00000
H6 H 4 e -0.03700 0.72600 0.04300 1.00000
H7 H 4 e 0.73300 -0.00800 0.06000 1.00000
H8 H 4 e 0.72900 0.01400 -0.04400 1.00000
K1 K 4 e 0.29373 0.73232 0.16965 1.00000
K2 K 4 e 0.25852 0.25782 0.16992 1.00000
O1 O 4 e 0.11630 0.62760 0.84660 1.00000
O2 O 4 e 0.88880 0.62610 0.16330 1.00000
O3 O 4 e 0.59980 -0.09080 0.14930 1.00000
O4 O 4 e 0.54550 0.15280 0.17910 1.00000
O5 O 4 e 0.06110 -0.01270 0.17250 1.00000
O6 O 4 e 0.09730 0.49380 0.18530 1.00000
O7 O 4 e 0.43420 0.49110 0.16740 1.00000
O8 O 4 e 0.37410 -0.02270 0.19310 1.00000
O9 O 4 e 0.35370 0.66840 -0.03420 1.00000
O10 O 4 e 0.64840 0.66530 0.04310 1.00000
O11 O 4 e -0.01310 0.77210 -0.00510 1.00000
O12 O 4 e 0.78020 0.01030 0.01110 1.00000
S1 S 4 e -0.04519 0.49523 0.20732 1.00000
S2 S 4 e 0.51973 0.00747 0.21252 1.00000
```

Manganese-leonite 110 K [$K_2Mn(SO_4)_2{\cdot}4H_2O$]: A8B2CD12E2_mP100_14_8e_2e_ad_12e_2e - POSCAR

```
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      ↪ y11,z11,x12,y12,z12,x13,y13,z13,x14,y14,z14,x15,y15,z15,x16,y16
      ↪ ,z16,x17,y17,z17,x18,y18,z18,x19,y19,z19,x20,y20,z20,x21,y21,
      ↪ z21,x22,y22,z22,x23,y23,z23,x24,y24,z24,x25,y25,z25,x26,y26,z26
      ↪  --params=9.902,0.962532821652,1.21500706928,95.02,0.387,0.729
      ↪ ,-0.07,0.275,0.649,-0.074,0.62,0.735,0.074,0.721,0.648,0.081,
      ↪ 0.025,0.72,-0.049,-0.037,0.726,0.043,0.733,-0.008,0.06,0.729,
      ↪ 0.014,-0.044,0.29373,0.73232,0.16965,0.25852,0.25782,0.16992,
      ↪ 0.1163,0.6276,0.8466,0.8888,0.6261,0.1633,0.5998,-0.0908,0.1493
      ↪ ,0.5455,0.1528,0.1791,0.0611,-0.0127,0.1725,0.0973,0.4938,
      ↪ 0.1853,0.4342,0.4911,0.1674,0.3741,-0.0227,0.1931,0.3537,0.6684
      ↪ ,-0.0342,0.6484,0.6653,0.0431,-0.0131,0.7721,-0.0051,0.7802,
      ↪ 0.0103,0.0111,-0.04519,0.49523,0.20732,0.51973,0.00747,0.21252
      ↪ & P2_{1}/c C_{2h}^{5} #14 (ade^24) & mP100 & None & H8K2MnO12S2
      ↪  & Manganese-leonite & B. Hertweck and G. Giester and E.
      ↪ Libowitzky, Am. Mineral. 86, 1282-1292 (2001)
   1.00000000000000
   9.90200000000000    0.00000000000000    0.00000000000000
   0.00000000000000    9.53100000000000    0.00000000000000
  -1.05275430748699    0.00000000000000   11.98485166233060
     H     K    Mn     O     S
    32     8     4    48     8
Direct
   0.38700000000000    0.72900000000000   -0.07000000000000    H    (4e)
  -0.38700000000000    1.22900000000000    0.57000000000000    H    (4e)
  -0.38700000000000   -0.72900000000000    0.07000000000000    H    (4e)
   0.38700000000000   -0.22900000000000    0.43000000000000    H    (4e)
   0.27500000000000    0.64900000000000   -0.07400000000000    H    (4e)
  -0.27500000000000    1.14900000000000    0.57400000000000    H    (4e)
  -0.27500000000000   -0.64900000000000    0.07400000000000    H    (4e)
   0.27500000000000   -0.14900000000000    0.42600000000000    H    (4e)
   0.62000000000000    0.73500000000000    0.07400000000000    H    (4e)
  -0.62000000000000    1.23500000000000    0.42600000000000    H    (4e)
  -0.62000000000000   -0.73500000000000   -0.07400000000000    H    (4e)
   0.62000000000000   -0.23500000000000    0.57400000000000    H    (4e)
   0.72100000000000    0.64800000000000    0.08100000000000    H    (4e)
  -0.72100000000000    1.14800000000000    0.41900000000000    H    (4e)
  -0.72100000000000   -0.64800000000000   -0.08100000000000    H    (4e)
   0.72100000000000   -0.14800000000000    0.58100000000000    H    (4e)
   0.02500000000000    0.72000000000000   -0.04900000000000    H    (4e)
  -0.02500000000000    1.22000000000000    0.54900000000000    H    (4e)
  -0.02500000000000   -0.72000000000000    0.04900000000000    H    (4e)
   0.02500000000000   -0.22000000000000    0.45100000000000    H    (4e)
  -0.03700000000000    0.72600000000000    0.04300000000000    H    (4e)
   0.03700000000000    1.22600000000000    0.45700000000000    H    (4e)
   0.03700000000000   -0.72600000000000   -0.04300000000000    H    (4e)
  -0.03700000000000   -0.22600000000000    0.54300000000000    H    (4e)
   0.73300000000000   -0.00800000000000    0.06000000000000    H    (4e)
  -0.73300000000000    0.49200000000000    0.44000000000000    H    (4e)
  -0.73300000000000    0.00800000000000   -0.06000000000000    H    (4e)
   0.73300000000000    0.50800000000000    0.56000000000000    H    (4e)
   0.72900000000000    0.01400000000000   -0.04400000000000    H    (4e)
  -0.72900000000000    0.51400000000000    0.54400000000000    H    (4e)
  -0.72900000000000   -0.01400000000000    0.04400000000000    H    (4e)
   0.72900000000000    0.48600000000000    0.45600000000000    H    (4e)
   0.29373000000000    0.73232000000000    0.16965000000000    K    (4e)
  -0.29373000000000    1.23232000000000    0.33035000000000    K    (4e)
  -0.29373000000000   -0.73232000000000   -0.16965000000000    K    (4e)
   0.29373000000000   -0.23232000000000    0.66965000000000    K    (4e)
   0.25852000000000    0.25782000000000    0.16992000000000    K    (4e)
  -0.25852000000000    0.75782000000000    0.33008000000000    K    (4e)
  -0.25852000000000   -0.25782000000000   -0.16992000000000    K    (4e)
   0.25852000000000    0.24218000000000    0.66992000000000    K    (4e)
   0.00000000000000    0.00000000000000    0.00000000000000   Mn    (2a)
   0.00000000000000    0.50000000000000    0.50000000000000   Mn    (2a)
   0.50000000000000    0.00000000000000    0.50000000000000   Mn    (2d)
   0.50000000000000    0.50000000000000    0.00000000000000   Mn    (2d)
   0.11630000000000    0.62760000000000    0.84660000000000    O    (4e)
  -0.11630000000000    1.12760000000000   -0.34660000000000    O    (4e)
  -0.11630000000000   -0.62760000000000   -0.84660000000000    O    (4e)
   0.11630000000000   -0.12760000000000    1.34660000000000    O    (4e)
   0.88880000000000    0.62610000000000    0.16330000000000    O    (4e)
  -0.88880000000000    1.12610000000000    0.33670000000000    O    (4e)
  -0.88880000000000   -0.62610000000000   -0.16330000000000    O    (4e)
   0.88880000000000   -0.12610000000000    0.66330000000000    O    (4e)
   0.59980000000000   -0.09080000000000    0.14930000000000    O    (4e)
  -0.59980000000000    0.40920000000000    0.35070000000000    O    (4e)
  -0.59980000000000    0.09080000000000   -0.14930000000000    O    (4e)
   0.59980000000000    0.59080000000000    0.64930000000000    O    (4e)
   0.54550000000000    0.15280000000000    0.17910000000000    O    (4e)
  -0.54550000000000    0.65280000000000    0.32090000000000    O    (4e)
  -0.54550000000000   -0.15280000000000   -0.17910000000000    O    (4e)
   0.54550000000000    0.34720000000000    0.67910000000000    O    (4e)
   0.06110000000000   -0.01270000000000    0.17250000000000    O    (4e)
  -0.06110000000000    0.48730000000000    0.32750000000000    O    (4e)
  -0.06110000000000    0.01270000000000   -0.17250000000000    O    (4e)
   0.06110000000000    0.51270000000000    0.67250000000000    O    (4e)
   0.09730000000000    0.49380000000000    0.18530000000000    O    (4e)
  -0.09730000000000    0.99380000000000    0.31470000000000    O    (4e)
  -0.09730000000000   -0.49380000000000   -0.18530000000000    O    (4e)
   0.09730000000000    0.00620000000000    0.68530000000000    O    (4e)
   0.43420000000000    0.49110000000000    0.16740000000000    O    (4e)
  -0.43420000000000    0.99110000000000    0.33260000000000    O    (4e)
  -0.43420000000000   -0.49110000000000   -0.16740000000000    O    (4e)
   0.43420000000000    0.00890000000000    0.66740000000000    O    (4e)
   0.37410000000000   -0.02270000000000    0.19310000000000    O    (4e)
  -0.37410000000000    0.47730000000000    0.30690000000000    O    (4e)
  -0.37410000000000    0.02270000000000   -0.19310000000000    O    (4e)
   0.37410000000000    0.52270000000000    0.69310000000000    O    (4e)
   0.35370000000000    0.66840000000000   -0.03420000000000    O    (4e)
  -0.35370000000000    1.16840000000000    0.53420000000000    O    (4e)
  -0.35370000000000   -0.66840000000000    0.03420000000000    O    (4e)
   0.35370000000000   -0.16840000000000    0.46580000000000    O    (4e)
   0.64840000000000    0.66530000000000    0.04310000000000    O    (4e)
  -0.64840000000000    1.16530000000000    0.45690000000000    O    (4e)
  -0.64840000000000   -0.66530000000000   -0.04310000000000    O    (4e)
   0.64840000000000   -0.16530000000000    0.54310000000000    O    (4e)
  -0.01310000000000    0.77210000000000   -0.00510000000000    O    (4e)
   0.01310000000000    1.27210000000000    0.50510000000000    O    (4e)
   0.01310000000000   -0.77210000000000    0.00510000000000    O    (4e)
  -0.01310000000000   -0.27210000000000    0.49490000000000    O    (4e)
   0.78020000000000    0.01030000000000    0.01110000000000    O    (4e)
  -0.78020000000000    0.51030000000000    0.48890000000000    O    (4e)
  -0.78020000000000   -0.01030000000000   -0.01110000000000    O    (4e)
   0.78020000000000    0.48970000000000    0.51110000000000    O    (4e)
  -0.04519000000000    0.49523000000000    0.20732000000000    S    (4e)
   0.04519000000000    0.99523000000000    0.29268000000000    S    (4e)
   0.04519000000000   -0.49523000000000   -0.20732000000000    S    (4e)
  -0.04519000000000    0.00477000000000    0.70732000000000    S    (4e)
   0.51973000000000    0.00747000000000    0.21252000000000    S    (4e)
  -0.51973000000000    0.50747000000000    0.28748000000000    S    (4e)
  -0.51973000000000   -0.00747000000000   -0.21252000000000    S    (4e)
   0.51973000000000    0.49253000000000    0.71252000000000    S    (4e)
```

Co_2Al_9 ($D8_d$): A9B2_mP22_14_a4e_e - CIF

```
# CIF file
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_audit_creation_method FINDSYM

_chemical_name_mineral 'Co2Al9'
_chemical_formula_sum 'Al9 Co2'

loop_
_publ_author_name
 'A. M. B. Douglas'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 3
_journal_year 1950
_journal_page_first 19
_journal_page_last 24
_publ_Section_title
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```
;
 The Structure of Co$_{2}$Al$_{9}$
;

_aflow_title 'Co$_{2}$Al$_{9}$ ($D8_{d}$) Structure'
_aflow_proto 'A9B2_mP22_14_a4e_e'
_aflow_params 'a,b/a,c/a,\beta,x_{2},y_{2},z_{2},x_{3},y_{3},z_{3},x_{4
      ↪ },y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{6}'
_aflow_params_values '6.213,1.01239336874,1.37719298246,94.76,0.4044,
      ↪ 0.5381,0.2682,0.0889,0.2101,0.293,0.3891,0.3069,-0.0014,0.2159,
      ↪ 0.8852,0.0417,0.2646,0.8851,0.3335'
_aflow_Strukturbericht '$D8_{d}$'
_aflow_Pearson 'mP22'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

_cell_length_a     6.21300
_cell_length_b     6.29000
_cell_length_c     8.55650
_cell_angle_alpha  90.00000
_cell_angle_beta   94.76000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z+1/2
3 -x,-y,-z
4 x,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Al1  Al  2  a  0.00000  0.00000  0.00000  1.00000
Al2  Al  4  e  0.40440  0.53810  0.26820  1.00000
Al3  Al  4  e  0.08890  0.21010  0.29300  1.00000
Al4  Al  4  e  0.38910  0.30690 -0.00140  1.00000
Al5  Al  4  e  0.21590  0.88520  0.04170  1.00000
Co1  Co  4  e  0.26460  0.88510  0.33350  1.00000
```

Co_2Al_9 ($D8_d$): A9B2_mP22_14_a4e_e - POSCAR

```
A9B2_mP22_14_a4e_e & a,b/a,c/a,beta,x2,y2,z2,x3,y3,z3,x4,y4,z4,x5,y5,z5,
      ↪ x6,y6,z6 --params=6.213,1.01239336874,1.37719298246,94.76,
      ↪ 0.4044,0.5381,0.2682,0.0889,0.2101,0.293,0.3891,0.3069,-0.0014,
      ↪ 0.2159,0.8852,0.0417,0.2646,0.8851,0.3335 & P2_{1}/c C_{2h}^{5}
      ↪  #14 (ae^5) & mP22 & $D8_{d}$ & Co2Al9 & Co2Al9 & A. M. B.
      ↪ Douglas, Acta Crystallogr. Sect. B Struct. Sci. 3, 19-24 (1950)
   1.00000000000000
   6.21300000000000    0.00000000000000    0.00000000000000
   0.00000000000000    6.29000000000000    0.00000000000000
  -0.71003667870571    0.00000000000000    8.52698892721766
     Al      Co
     18       4
Direct
   0.00000000000000    0.00000000000000    0.00000000000000    Al    (2a)
   0.00000000000000    0.50000000000000    0.50000000000000    Al    (2a)
   0.40440000000000    0.53810000000000    0.26820000000000    Al    (4e)
  -0.40440000000000    1.03810000000000    0.23180000000000    Al    (4e)
  -0.40440000000000   -0.53810000000000   -0.26820000000000    Al    (4e)
   0.40440000000000   -0.03810000000000    0.76820000000000    Al    (4e)
   0.08890000000000    0.21010000000000    0.29300000000000    Al    (4e)
  -0.08890000000000    0.71010000000000    0.20700000000000    Al    (4e)
  -0.08890000000000   -0.21010000000000   -0.29300000000000    Al    (4e)
   0.08890000000000    0.28990000000000    0.79300000000000    Al    (4e)
   0.38910000000000    0.30690000000000   -0.00140000000000    Al    (4e)
  -0.38910000000000    0.80690000000000    0.50140000000000    Al    (4e)
  -0.38910000000000   -0.30690000000000    0.00140000000000    Al    (4e)
   0.38910000000000    0.19310000000000    0.49860000000000    Al    (4e)
   0.21590000000000    0.88520000000000    0.04170000000000    Al    (4e)
  -0.21590000000000    1.38520000000000    0.45830000000000    Al    (4e)
  -0.21590000000000   -0.88520000000000   -0.04170000000000    Al    (4e)
   0.21590000000000   -0.38520000000000    0.54170000000000    Al    (4e)
   0.26460000000000    0.88510000000000    0.33350000000000    Co    (4e)
  -0.26460000000000    1.38510000000000    0.16650000000000    Co    (4e)
  -0.26460000000000   -0.88510000000000   -0.33350000000000    Co    (4e)
   0.26460000000000   -0.38510000000000    0.83350000000000    Co    (4e)
```

Tutton salt $[Cu(NH_4)_2(SO_4)_2{\cdot}6H_2O, H4_4]$: AB20C2D14E2_mP78_14_a_10e_e_7e_e - CIF

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_audit_creation_method FINDSYM

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_chemical_formula_sum 'Cu H20 N2 O14 S2'

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_publ_author_name
 'G. M. Brown'
 'R. Chidambaram'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 25
_journal_year 1969
_journal_page_first 676
_journal_page_last 687
_publ_Section_title
;
 The structure of copper ammonium sulfate hexahydrate from neutron
      ↪ diffraction data
;

_aflow_title 'Tutton salt [Cu(NH$_{4}$)$_{2}$(SO$_{4}$)$_{2}$$\cdot$6H$_
      ↪ {2}$O, $H4_{4}$] Structure'
_aflow_proto 'AB20C2D14E2_mP78_14_a_10e_e_7e_e'
_aflow_params 'a,b/a,c/a,\beta,x_{2},y_{2},z_{2},x_{3},y_{3},z_{3},x_{4
      ↪ },y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{6},x_{7},y_{7},
      ↪ z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{10},y_{10},z_{10},
      ↪ x_{11},y_{11},z_{11},x_{12},y_{12},z_{12},x_{13},y_{13},z_{13},
      ↪ x_{14},y_{14},z_{14},x_{15},y_{15},z_{15},x_{16},y_{16},z_{16},
      ↪ x_{17},y_{17},z_{17},x_{18},y_{18},z_{18},x_{19},y_{19},z_{19},
      ↪ x_{20},y_{20},z_{20}'
_aflow_params_values '6.3016,1.96450107909,1.4616129237,106.112,0.20966,
      ↪ 0.66679,0.06554,0.40162,0.70084,0.22515,0.46879,0.66573,0.07617
      ↪ ,0.36184,0.57533,0.16861,0.32899,-0.0962,0.22103,0.1068,0.87132
      ↪ ,0.25555,-0.0597,-0.09535,0.73282,0.00217,0.81595,0.85798,
      ↪ 0.31563,0.05877,0.89729,0.3095,0.14082,0.02457,0.35953,0.65214,
      ↪ 0.13458,0.60057,0.76827,0.41576,0.78118,-0.0758,0.54888,0.63349
      ↪ ,-0.07032,0.28057,-0.04314,0.82138,0.39147,0.17726,0.88327,
      ↪ 0.17658,0.03043,0.89108,0.83608,0.2821,0.06534,-0.00531,0.74517
      ↪ ,0.86102,0.41069'
_aflow_Strukturbericht '$H4_{4}$'
_aflow_Pearson 'mP78'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

_cell_length_a     6.30160
_cell_length_b     12.37950
_cell_length_c     9.21050
_cell_angle_alpha  90.00000
_cell_angle_beta   106.11200
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z+1/2
3 -x,-y,-z
4 x,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cu1  Cu  2  a  0.00000  0.00000  0.00000  1.00000
H1   H   4  e  0.20966  0.66679  0.06554  1.00000
H2   H   4  e  0.40162  0.70084  0.22515  1.00000
H3   H   4  e  0.46879  0.66573  0.07617  1.00000
H4   H   4  e  0.36184  0.57533  0.16861  1.00000
H5   H   4  e  0.32899 -0.09620  0.22103  1.00000
H6   H   4  e  0.10680  0.87132  0.25555  1.00000
H7   H   4  e -0.05970 -0.09535  0.73282  1.00000
H8   H   4  e  0.00217  0.81595  0.85798  1.00000
H9   H   4  e  0.31563  0.05877  0.89729  1.00000
H10  H   4  e  0.30950  0.14082  0.02457  1.00000
N1   N   4  e  0.35953  0.65214  0.13458  1.00000
O1   O   4  e  0.60057  0.76827  0.41576  1.00000
O2   O   4  e  0.78118 -0.07580  0.54888  1.00000
O3   O   4  e  0.63349 -0.07032  0.28057  1.00000
O4   O   4  e -0.04314  0.82138  0.39147  1.00000
O5   O   4  e  0.17726  0.88327  0.17658  1.00000
O6   O   4  e  0.03043  0.89108  0.83608  1.00000
O7   O   4  e  0.28210  0.06534 -0.00531  1.00000
S1   S   4  e  0.74517  0.86102  0.41069  1.00000
```

Tutton salt $[Cu(NH_4)_2(SO_4)_2{\cdot}6H_2O, H4_4]$: AB20C2D14E2_mP78_14_a_10e_e_7e_e - POSCAR

```
AB20C2D14E2_mP78_14_a_10e_e_7e_e & a,b/a,c/a,beta,x2,y2,z2,x3,y3,z3,x4,
      ↪ y4,z4,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10,
      ↪ x11,y11,z11,x12,y12,z12,x13,y13,z13,x14,y14,z14,x15,y15,z15,x16
      ↪ ,y16,z16,x17,y17,z17,x18,y18,z18,x19,y19,z19,x20,y20,z20 --
      ↪ params=6.3016,1.96450107909,1.4616129237,106.112,0.20966,
      ↪ 0.66679,0.06554,0.40162,0.70084,0.22515,0.46879,0.66573,0.07617
      ↪ ,0.36184,0.57533,0.16861,0.32899,-0.0962,0.22103,0.1068,0.87132
      ↪ ,0.25555,-0.0597,-0.09535,0.73282,0.00217,0.81595,0.85798,
      ↪ 0.31563,0.05877,0.89729,0.3095,0.14082,0.02457,0.35953,0.65214,
      ↪ 0.13458,0.60057,0.76827,0.41576,0.78118,-0.0758,0.54888,0.63349
      ↪ ,-0.07032,0.28057,-0.04314,0.82138,0.39147,0.17726,0.88327,
      ↪ 0.17658,0.03043,0.89108,0.83608,0.2821,0.06534,-0.00531,0.74517
      ↪ ,0.86102,0.41069 & mP78 &
      ↪ $H4_{4}$ & CuH20N2O14S2 & Tutton salt & G. M. Brown and R.
      ↪ Chidambaram, Acta Crystallogr. Sect. B Struct. Sci. 25, 676-687
      ↪ (1969)
   1.00000000000000
   6.30160000000000    0.00000000000000    0.00000000000000
   0.00000000000000   12.37950000000000    0.00000000000000
  -2.55605994213425    0.00000000000000    8.84872125350419
     Cu       H       N       O       S
      2      40       4      28       4
Direct
   0.00000000000000    0.00000000000000    0.00000000000000    Cu    (2a)
   0.00000000000000    0.50000000000000    0.50000000000000    Cu    (2a)
```

```
  0.20966000000000   0.66679000000000   0.06554000000000   H   (4e)
 -0.20966000000000   1.16679000000000   0.43446000000000   H   (4e)
 -0.20966000000000  -0.66679000000000  -0.06554000000000   H   (4e)
  0.20966000000000  -0.16679000000000   0.56554000000000   H   (4e)
  0.40162000000000   0.70084000000000   0.22515000000000   H   (4e)
 -0.40162000000000   1.20084000000000   0.27485000000000   H   (4e)
 -0.40162000000000  -0.70084000000000  -0.22515000000000   H   (4e)
  0.40162000000000  -0.20084000000000   0.72515000000000   H   (4e)
  0.46879000000000   0.66573000000000   0.07617000000000   H   (4e)
 -0.46879000000000   1.16573000000000   0.42383000000000   H   (4e)
 -0.46879000000000  -0.66573000000000  -0.07617000000000   H   (4e)
  0.46879000000000  -0.16573000000000   0.57617000000000   H   (4e)
  0.36184000000000   0.57533000000000   0.16861000000000   H   (4e)
 -0.36184000000000   1.07533000000000   0.33139000000000   H   (4e)
 -0.36184000000000  -0.57533000000000  -0.16861000000000   H   (4e)
  0.36184000000000  -0.07533000000000   0.66861000000000   H   (4e)
  0.32899000000000  -0.09620000000000   0.22103000000000   H   (4e)
 -0.32899000000000   0.40380000000000   0.27897000000000   H   (4e)
 -0.32899000000000   0.09620000000000  -0.22103000000000   H   (4e)
  0.32899000000000   0.59620000000000   0.72103000000000   H   (4e)
  0.10680000000000   0.87132000000000   0.25555000000000   H   (4e)
 -0.10680000000000   1.37132000000000   0.24445000000000   H   (4e)
 -0.10680000000000  -0.87132000000000  -0.25555000000000   H   (4e)
  0.10680000000000  -0.37132000000000   0.75555000000000   H   (4e)
 -0.05970000000000  -0.09535000000000   0.73282000000000   H   (4e)
  0.05970000000000   0.40465000000000  -0.23282000000000   H   (4e)
  0.05970000000000   0.09535000000000  -0.73282000000000   H   (4e)
 -0.05970000000000   0.59535000000000   1.23282000000000   H   (4e)
  0.00217000000000   0.81595000000000   0.85798000000000   H   (4e)
 -0.00217000000000   1.31595000000000  -0.35798000000000   H   (4e)
 -0.00217000000000  -0.81595000000000  -0.85798000000000   H   (4e)
  0.00217000000000  -0.31595000000000   1.35798000000000   H   (4e)
  0.31563000000000   0.05877000000000   0.89729000000000   H   (4e)
 -0.31563000000000   0.55877000000000  -0.39729000000000   H   (4e)
 -0.31563000000000  -0.05877000000000  -0.89729000000000   H   (4e)
  0.31563000000000   0.44123000000000   1.39729000000000   H   (4e)
  0.30950000000000   0.14082000000000   0.02457000000000   H   (4e)
 -0.30950000000000   0.64082000000000   0.47543000000000   H   (4e)
 -0.30950000000000  -0.14082000000000  -0.02457000000000   H   (4e)
  0.30950000000000   0.35918000000000   0.52457000000000   H   (4e)
  0.35953000000000   0.65214000000000   0.13458000000000   N   (4e)
 -0.35953000000000   1.15214000000000   0.36542000000000   N   (4e)
 -0.35953000000000  -0.65214000000000  -0.13458000000000   N   (4e)
  0.35953000000000  -0.15214000000000   0.63458000000000   N   (4e)
  0.60057000000000   0.76827000000000   0.41576000000000   O   (4e)
 -0.60057000000000   1.26827000000000   0.08424000000000   O   (4e)
 -0.60057000000000  -0.76827000000000  -0.41576000000000   O   (4e)
  0.60057000000000  -0.26827000000000   0.91576000000000   O   (4e)
  0.78118000000000  -0.07588000000000   0.54888000000000   O   (4e)
 -0.78118000000000   0.42420000000000  -0.04888000000000   O   (4e)
 -0.78118000000000   0.07588000000000  -0.54888000000000   O   (4e)
  0.78118000000000   0.57588000000000   1.04888000000000   O   (4e)
  0.63349000000000  -0.07032000000000   0.28057000000000   O   (4e)
 -0.63349000000000   0.42968000000000   0.21943000000000   O   (4e)
 -0.63349000000000   0.07032000000000  -0.28057000000000   O   (4e)
  0.63349000000000   0.57032000000000   0.78057000000000   O   (4e)
 -0.04314000000000   0.82138000000000   0.39147000000000   O   (4e)
  0.04314000000000   1.32138000000000   0.10853000000000   O   (4e)
  0.04314000000000  -0.82138000000000  -0.39147000000000   O   (4e)
 -0.04314000000000  -0.32138000000000   0.89147000000000   O   (4e)
  0.17726000000000   0.88327000000000   0.17658000000000   O   (4e)
 -0.17726000000000   1.38327000000000   0.32342000000000   O   (4e)
 -0.17726000000000  -0.88327000000000  -0.17658000000000   O   (4e)
  0.17726000000000  -0.38327000000000   0.67658000000000   O   (4e)
  0.03043000000000   0.89108000000000   0.83608000000000   O   (4e)
 -0.03043000000000   1.39108000000000  -0.33608000000000   O   (4e)
 -0.03043000000000  -0.89108000000000  -0.83608000000000   O   (4e)
  0.03043000000000  -0.39108000000000   1.33608000000000   O   (4e)
  0.28210000000000   0.06534000000000  -0.00531000000000   O   (4e)
 -0.28210000000000   0.56534000000000   0.50531000000000   O   (4e)
 -0.28210000000000  -0.06534000000000   0.00531000000000   O   (4e)
  0.28210000000000   0.43466000000000   0.49469000000000   O   (4e)
  0.74517000000000   0.86102000000000   0.41069000000000   S   (4e)
 -0.74517000000000   1.36102000000000   0.08931000000000   S   (4e)
 -0.74517000000000  -0.86102000000000  -0.41069000000000   S   (4e)
  0.74517000000000  -0.36102000000000   0.91069000000000   S   (4e)
```

Parawollastonite ($CaSiO_3$, $S3_3$ (II)): AB3C_mP60_14_3e_9e_3e - CIF

```
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_audit_creation_method FINDSYM

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_chemical_formula_sum 'Ca O3 Si'

loop_
_publ_author_name
 'F. J. Trojer'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 127
_journal_year 1968
_journal_page_first 291
_journal_page_last 308
_publ_Section_title
;
 The crystal structure of parawollastonite
;

_aflow_title 'Parawollastonite (CaSiO$_{3}$, $S3_{3}$ (II)) Structure'
_aflow_proto 'AB3C_mP60_14_3e_9e_3e'
_aflow_params 'a,b/a,c/a,\beta,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3
      ↪ },y_{3},z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},
      ↪ z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{
      ↪ 10},y_{10},z_{10},x_{11},y_{11},z_{11},x_{12},y_{12},z_{12},x_{
      ↪ 13},y_{13},z_{13},x_{14},y_{14},z_{14},x_{15},y_{15},z_{15}'
_aflow_params_values '7.066,1.03594678743,2.18313048401,95.40417,-0.0288
      ↪ ,0.6242,0.2482,0.7397,0.3735,0.4011,0.7364,0.8791,0.3987,0.6685
      ↪ ,0.6253,0.3,0.3031,0.6241,0.2156,0.0328,0.8603,0.349,0.0348,
      ↪ 0.3843,0.3473,0.2388,0.8774,0.5086,0.2347,0.3824,0.5078,0.406,
      ↪ 0.8038,0.3642,0.4067,0.4467,0.3633,0.2767,0.1245,0.3906,0.2313
      ↪ ,-0.0907,0.4076,0.2313,0.3402,0.4075,0.4432,0.6239,0.3016'
_aflow_Strukturbericht '$S3_{3}$ (II)'
_aflow_Pearson 'mP60'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

_cell_length_a      7.06600
_cell_length_b      7.32000
_cell_length_c      15.42600
_cell_angle_alpha   90.00000
_cell_angle_beta    95.40417
_cell_angle_gamma   90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z+1/2
3 -x,-y,-z
4 x,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ca1 Ca 4 e -0.02880 0.62420 0.24820 1.00000
Ca2 Ca 4 e 0.73970 0.37350 0.40110 1.00000
Ca3 Ca 4 e 0.73640 0.87910 0.39870 1.00000
O1 O 4 e 0.66850 0.62530 0.30000 1.00000
O2 O 4 e 0.30310 0.62410 0.21560 1.00000
O3 O 4 e 0.03280 0.86030 0.34900 1.00000
O4 O 4 e 0.03480 0.38430 0.34730 1.00000
O5 O 4 e 0.23880 0.87740 0.50860 1.00000
O6 O 4 e 0.23470 0.38240 0.50780 1.00000
O7 O 4 e 0.40600 0.80380 0.36420 1.00000
O8 O 4 e 0.40670 0.44670 0.36330 1.00000
O9 O 4 e 0.27670 0.12450 0.39060 1.00000
Si1 Si 4 e 0.23130 -0.09070 0.40760 1.00000
Si2 Si 4 e 0.23130 0.34020 0.40750 1.00000
Si3 Si 4 e 0.44320 0.62390 0.30160 1.00000
```

Parawollastonite ($CaSiO_3$, $S3_3$ (II)): AB3C_mP60_14_3e_9e_3e - POSCAR

```
AB3C_mP60_14_3e_9e_3e & a,b/a,c/a,beta,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,
      ↪ z4,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10,x11
      ↪ ,y11,z11,x12,y12,z12,x13,y13,z13,x14,y14,z14,x15,y15,z15 --
      ↪ params=7.066,1.03594678743,2.18313048401,95.40417,-0.0288,
      ↪ 0.6242,0.2482,0.7397,0.3735,0.4011,0.7364,0.8791,0.3987,0.6685,
      ↪ 0.6253,0.3,0.3031,0.6241,0.2156,0.0328,0.8603,0.349,0.0348,
      ↪ 0.3843,0.3473,0.2388,0.8774,0.5086,0.2347,0.3824,0.5078,0.406,
      ↪ 0.8038,0.3642,0.4067,0.4467,0.3633,0.2767,0.1245,0.3906,0.2313
      ↪ ,-0.0907,0.4076,0.2313,0.3402,0.4075,0.4432,0.6239,0.3016 & P2_
      ↪ {1}/c C_{2h}^{5} #14 (e^15) & mP60 & $S3_{3}$ (II) & CaO3Si &
      ↪ Parawollastonite & F. J. Trojer, Zeitschrift f{"u}r
      ↪ Kristallographie - Crystalline Materials 127, 291-308 (1968)
   1.00000000000000
   7.06600000000000     0.00000000000000     0.00000000000000
   0.00000000000000     7.32000000000000     0.00000000000000
  -1.45283256261353     0.00000000000000    15.35743316915330
    Ca    O    Si
    12    36    12
Direct
  -0.02880000000000    0.62420000000000    0.24820000000000   Ca   (4e)
   0.02880000000000    1.12420000000000    0.25180000000000   Ca   (4e)
   0.02880000000000   -0.62420000000000   -0.24820000000000   Ca   (4e)
  -0.02880000000000   -0.12420000000000    0.74820000000000   Ca   (4e)
   0.73970000000000    0.37350000000000    0.40110000000000   Ca   (4e)
  -0.73970000000000    0.87350000000000    0.09890000000000   Ca   (4e)
  -0.73970000000000   -0.37350000000000   -0.40110000000000   Ca   (4e)
   0.73970000000000    0.12650000000000    0.90110000000000   Ca   (4e)
   0.73640000000000    0.87910000000000    0.39870000000000   Ca   (4e)
  -0.73640000000000    1.37910000000000    0.10130000000000   Ca   (4e)
  -0.73640000000000   -0.87910000000000   -0.39870000000000   Ca   (4e)
   0.73640000000000   -0.37910000000000    0.89870000000000   Ca   (4e)
   0.66850000000000    0.62530000000000    0.30000000000000   O    (4e)
  -0.66850000000000    1.12530000000000    0.20000000000000   O    (4e)
  -0.66850000000000   -0.62530000000000   -0.30000000000000   O    (4e)
   0.66850000000000   -0.12530000000000    0.80000000000000   O    (4e)
   0.30310000000000    0.62410000000000    0.21560000000000   O    (4e)
  -0.30310000000000    1.12410000000000    0.28440000000000   O    (4e)
  -0.30310000000000   -0.62410000000000   -0.21560000000000   O    (4e)
   0.30310000000000   -0.12410000000000    0.71560000000000   O    (4e)
   0.03280000000000    0.86030000000000    0.34900000000000   O    (4e)
  -0.03280000000000    1.36030000000000    0.15100000000000   O    (4e)
  -0.03280000000000   -0.86030000000000   -0.34900000000000   O    (4e)
   0.03280000000000   -0.36030000000000    0.84900000000000   O    (4e)
   0.03480000000000    0.38430000000000    0.34730000000000   O    (4e)
  -0.03480000000000    0.88430000000000    0.15270000000000   O    (4e)
  -0.03480000000000   -0.38430000000000   -0.34730000000000   O    (4e)
   0.03480000000000    0.11570000000000    0.84730000000000   O    (4e)
```

```
   0.23880000000000   0.87740000000000   0.50860000000000   O   (4e)
  -0.23880000000000   1.37740000000000  -0.00860000000000   O   (4e)
  -0.23880000000000  -0.87740000000000  -0.50860000000000   O   (4e)
   0.23880000000000  -0.37740000000000   1.00860000000000   O   (4e)
   0.23470000000000   0.38240000000000   0.50780000000000   O   (4e)
  -0.23470000000000   0.88240000000000  -0.00780000000000   O   (4e)
  -0.23470000000000  -0.38240000000000  -0.50780000000000   O   (4e)
   0.23470000000000   0.11760000000000   1.00780000000000   O   (4e)
   0.40600000000000   0.80380000000000   0.36420000000000   O   (4e)
  -0.40600000000000   1.30380000000000   0.13580000000000   O   (4e)
  -0.40600000000000  -0.80380000000000  -0.36420000000000   O   (4e)
   0.40600000000000  -0.30380000000000   0.86420000000000   O   (4e)
   0.40670000000000   0.44670000000000   0.36330000000000   O   (4e)
  -0.40670000000000   0.94670000000000   0.13670000000000   O   (4e)
  -0.40670000000000  -0.44670000000000  -0.36330000000000   O   (4e)
   0.40670000000000   0.05330000000000   0.86330000000000   O   (4e)
   0.27670000000000   0.12450000000000   0.39060000000000   O   (4e)
  -0.27670000000000   0.62450000000000   0.10940000000000   O   (4e)
  -0.27670000000000  -0.12450000000000  -0.39060000000000   O   (4e)
   0.27670000000000   0.37550000000000   0.89060000000000   O   (4e)
   0.23130000000000  -0.09070000000000   0.40760000000000   Si  (4e)
  -0.23130000000000   0.40930000000000   0.09240000000000   Si  (4e)
  -0.23130000000000   0.09070000000000  -0.40760000000000   Si  (4e)
   0.23130000000000   0.59070000000000   0.90760000000000   Si  (4e)
   0.23130000000000   0.34020000000000   0.40750000000000   Si  (4e)
  -0.23130000000000   0.84020000000000   0.09250000000000   Si  (4e)
  -0.23130000000000  -0.34020000000000  -0.40750000000000   Si  (4e)
   0.23130000000000   0.15980000000000   0.90750000000000   Si  (4e)
   0.44320000000000   0.62390000000000   0.30160000000000   Si  (4e)
  -0.44320000000000   1.12390000000000   0.19840000000000   Si  (4e)
  -0.44320000000000  -0.62390000000000  -0.30160000000000   Si  (4e)
   0.44320000000000  -0.12390000000000   0.80160000000000   Si  (4e)
```

β-B_2H_6: AB3_mP16_14_e_3e - CIF

```
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_chemical_formula_sum 'B H3'

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 'H. W. Smith'
 'W. N. Lipscomb'
_journal_name_full_name
;
 Journal of Chemical Physics
;
_journal_volume 43
_journal_year 1965
_journal_page_first 1060
_journal_page_last 1064
_publ_Section_title
;
 Single-Crystal X-Ray Diffraction Study of $\beta$-Diborane
;

_aflow_title '$\beta$-B$_{2}$H$_{6}$ Structure'
_aflow_proto 'AB3_mP16_14_e_3e'
_aflow_params 'a,b/a,c/a,\beta,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3
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_aflow_params_values '4.4,1.47727272727,1.41869318182,117.78562,0.144,
      ↪ 0.042,0.146,0.36,0.14,0.166,0.098,-0.005,0.294,0.877,0.112,-
      ↪ 0.019'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mP16'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

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_cell_length_b     6.50000
_cell_length_c     6.24225
_cell_angle_alpha  90.00000
_cell_angle_beta   117.78562
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z+1/2
3 -x,-y,-z
4 x,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
B1 B 4 e 0.14400 0.04200 0.14600 1.00000
H1 H 4 e 0.36000 0.14000 0.16600 1.00000
H2 H 4 e 0.09800 -0.00500 0.29400 1.00000
H3 H 4 e 0.87700 0.11200 -0.01900 1.00000
```

β-B_2H_6: AB3_mP16_14_e_3e - POSCAR

```
AB3_mP16_14_e_3e & a,b/a,c/a,beta,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,z4 --
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      ↪ 0.146,0.36,0.14,0.166,0.098,-0.005,0.294,0.877,0.112,-0.019 &
      ↪ P2_{1}/c C_{2h}^{5} #14 (e^4) & mP16 & None & B2H6 & beta & H.
      ↪ W. Smith and W. N. Lipscomb, J. Chem. Phys. 43, 1060-1064 (1965
      ↪ )
   1.00000000000000
   4.40000000000000   0.00000000000000   0.00000000000000
   0.00000000000000   6.50000000000000   0.00000000000000
  -2.90991606787698   0.00000000000000   5.52250609238337
     B     H
     4    12
Direct
   0.14400000000000   0.04200000000000   0.14600000000000   B   (4e)
  -0.14400000000000   0.54200000000000   0.35400000000000   B   (4e)
  -0.14400000000000  -0.04200000000000  -0.14600000000000   B   (4e)
   0.14400000000000   0.45800000000000   0.64600000000000   B   (4e)
   0.36000000000000   0.14000000000000   0.16600000000000   H   (4e)
  -0.36000000000000   0.64000000000000   0.33400000000000   H   (4e)
  -0.36000000000000  -0.14000000000000  -0.16600000000000   H   (4e)
   0.36000000000000   0.36000000000000   0.66600000000000   H   (4e)
   0.09800000000000  -0.00500000000000   0.29400000000000   H   (4e)
  -0.09800000000000   0.49500000000000   0.20600000000000   H   (4e)
  -0.09800000000000   0.00500000000000  -0.29400000000000   H   (4e)
   0.09800000000000   0.50500000000000   0.79400000000000   H   (4e)
   0.87700000000000   0.11200000000000  -0.01900000000000   H   (4e)
  -0.87700000000000   0.61200000000000   0.51900000000000   H   (4e)
  -0.87700000000000  -0.11200000000000   0.01900000000000   H   (4e)
   0.87700000000000   0.38800000000000   0.48100000000000   H   (4e)
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B_2H_6 ($P2_1/c$): AB3_mP16_14_e_3e - CIF

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_chemical_formula_sum 'B H3'

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 'Y. Yao'
 'R. Hoffmann'
_journal_name_full_name
;
 Journal of the American Chemical Society
;
_journal_volume 133
_journal_year 2011
_journal_page_first 21002
_journal_page_last 21009
_publ_Section_title
;
 BH$_{3}$ under Pressure: Leaving the Molecular Diborane Motif
;

_aflow_title 'B$_{2}$H$_{6}$ ($P2_{1}/c$) Structure'
_aflow_proto 'AB3_mP16_14_e_3e'
_aflow_params 'a,b/a,c/a,\beta,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3
      ↪ },y_{3},z_{3},x_{4},y_{4},z_{4}'
_aflow_params_values '4.46,1.94618834081,1.0201793722,120.5,0.3698,
      ↪ 0.4285,0.0119,0.0865,0.4057,0.7725,0.4789,0.1446,0.7686,0.5981,
      ↪ 0.4237,-0.0903'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mP16'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

_cell_length_a     4.46000
_cell_length_b     8.68000
_cell_length_c     4.55000
_cell_angle_alpha  90.00000
_cell_angle_beta   120.50000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z+1/2
3 -x,-y,-z
4 x,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
B1 B 4 e 0.36980 0.42850 0.01190 1.00000
H1 H 4 e 0.08650 0.40570 0.77250 1.00000
H2 H 4 e 0.47890 0.14460 0.76860 1.00000
H3 H 4 e 0.59810 0.42370 -0.09030 1.00000
```

B_2H_6 ($P2_1/c$): AB3_mP16_14_e_3e - POSCAR

```
AB3_mP16_14_e_3e & a,b/a,c/a,beta,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,z4 --
      ↪ params=4.46,1.94618834081,1.0201793722,120.5,0.3698,0.4285,
      ↪ 0.0119,0.0865,0.4057,0.7725,0.4789,0.1446,0.7686,0.5981,0.4237
      ↪ ,-0.0903 & P2_{1}/c C_{2h}^{5} #14 (e^4) & mP16 & None & B2H6 &
      ↪  B2H6 & Y. Yao and R. Hoffmann, J. Am. Chem. Soc. 133,
      ↪ 21002-21009 (2011)
   1.00000000000000
```

```
   4.46000000000000   0.00000000000000   0.00000000000000
   0.00000000000000   8.68000000000000   0.00000000000000
  -2.30929955147120   0.00000000000000   3.92041268000894
     B     H
     4    12
Direct
   0.36980000000000   0.42850000000000   0.01190000000000  B   (4e)
  -0.36980000000000   0.92850000000000   0.48810000000000  B   (4e)
  -0.36980000000000  -0.42850000000000  -0.01190000000000  B   (4e)
   0.36980000000000   0.07150000000000   0.51190000000000  B   (4e)
   0.08650000000000   0.40570000000000   0.77250000000000  H   (4e)
  -0.08650000000000   0.90570000000000  -0.27250000000000  H   (4e)
  -0.08650000000000  -0.40570000000000  -0.77250000000000  H   (4e)
   0.08650000000000   0.09430000000000   1.27250000000000  H   (4e)
   0.47890000000000   0.14460000000000   0.76860000000000  H   (4e)
  -0.47890000000000   0.64460000000000  -0.26860000000000  H   (4e)
  -0.47890000000000  -0.14460000000000  -0.76860000000000  H   (4e)
   0.47890000000000   0.35540000000000   1.26860000000000  H   (4e)
   0.59810000000000   0.42370000000000  -0.09030000000000  H   (4e)
  -0.59810000000000   0.92370000000000   0.59030000000000  H   (4e)
  -0.59810000000000  -0.42370000000000   0.09030000000000  H   (4e)
   0.59810000000000   0.07630000000000   0.40970000000000  H   (4e)
```

KAuBr$_4$·2H$_2$O ($H4_{19}$): AB4C2D_mP32_14_e_4e_2e_e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'AuBr4(H2O)2K'
_chemical_formula_sum 'Au Br4 (H2O)2 K'

loop_
_publ_author_name
 'H. Omrani'
 'F. Th\\'{e}obald'
 'H. Vivier'
_journal_name_full_name
;
 Acta Crystallographica Section C: Structural Chemistry
;
_journal_volume 42
_journal_year 1986
_journal_page_first 1091
_journal_page_last 1092
_publ_Section_title
;
 Structure of potassium tetrabromoaurate(III) dihydrate
;

# Found in Sodium tetrabromoaurate(III) dihydrate, 2001

_aflow_title 'KAuBr$_{4}$\cdot$2H$_{2}$O ($H4_{19}$) Structure'
_aflow_proto 'AB4C2D_mP32_14_e_4e_2e_e'
_aflow_params 'a,b/a,c/a,\beta,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3
      ↪ },y_{3},z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},
      ↪ z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},z_{8}'
_aflow_params_values '8.48,1.41580188679,1.45127830189,129.09091,0.2507,
      ↪ 0.0067,0.5026,0.3935,0.8371,0.6395,0.114,0.1778,0.3679,0.2208,-
      ↪ 0.0834,0.313,0.2767,0.0954,0.6901,0.5843,0.1353,0.0613,0.0171,
      ↪ 0.364,0.5512,0.2671,0.4723,0.5181'
_aflow_Strukturbericht '$H4_{19}$'
_aflow_Pearson 'mP32'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

_cell_length_a    8.48000
_cell_length_b    12.00600
_cell_length_c    12.30684
_cell_angle_alpha 90.00000
_cell_angle_beta  129.09091
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z+1/2
3 -x,-y,-z
4 x,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Au1 Au 4 e 0.25070 0.00670 0.50260 1.00000
Br1 Br 4 e 0.39350 0.83710 0.63950 1.00000
Br2 Br 4 e 0.11400 0.17780 0.36790 1.00000
Br3 Br 4 e 0.22080 -0.08340 0.31300 1.00000
Br4 Br 4 e 0.27670 0.09540 0.69010 1.00000
H2O1 H2O 4 e 0.58430 0.13530 0.06130 1.00000
H2O2 H2O 4 e 0.01710 0.36400 0.55120 1.00000
K1 K 4 e 0.26710 0.47230 0.51810 1.00000
```

KAuBr$_4$·2H$_2$O ($H4_{19}$): AB4C2D_mP32_14_e_4e_2e_e - POSCAR

```
AB4C2D_mP32_14_e_4e_2e_e & a,b/a,c/a,beta,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,
      ↪ y4,z4,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8 --params=8.48,
      ↪ 1.41580188679,1.45127830189,129.09091,0.2507,0.0067,0.5026,
      ↪ 0.3935,0.8371,0.6395,0.114,0.1778,0.3679,0.2208,-0.0834,0.313,
      ↪ 0.2767,0.0954,0.6901,0.5843,0.1353,0.0613,0.0171,0.364,0.5512,
      ↪ 0.2671,0.4723,0.5181 & P2_{1}/c C_{2h}^{5} #14 (e^8) & mP32 &
      ↪ $H4_{19}$ & AuBr4(H2O)2K & AuBr4(H2O)2K & H. Omrani and F. Th
      ↪ \'{e}obald and H. Vivier, Acta Crystallogr. C 42, 1091-1092 (
      ↪ 1986)
   1.00000000000000
   8.48000000000000   0.00000000000000   0.00000000000000
   0.00000000000000  12.00600000000000   0.00000000000000
  -7.76011093693911   0.00000000000000   9.55191022947756
    Au    Br   H2O    K
     4    16     8     4
Direct
   0.25070000000000   0.00670000000000   0.50260000000000  Au   (4e)
  -0.25070000000000   0.50670000000000  -0.00260000000000  Au   (4e)
  -0.25070000000000  -0.00670000000000  -0.50260000000000  Au   (4e)
   0.25070000000000   0.49330000000000   1.00260000000000  Au   (4e)
   0.39350000000000   0.83710000000000   0.63950000000000  Br   (4e)
  -0.39350000000000   1.33710000000000  -0.13950000000000  Br   (4e)
  -0.39350000000000  -0.83710000000000  -0.63950000000000  Br   (4e)
   0.39350000000000  -0.33710000000000   1.13950000000000  Br   (4e)
   0.11400000000000   0.17780000000000   0.36790000000000  Br   (4e)
  -0.11400000000000   0.67780000000000   0.13210000000000  Br   (4e)
  -0.11400000000000  -0.17780000000000  -0.36790000000000  Br   (4e)
   0.11400000000000   0.32220000000000   0.86790000000000  Br   (4e)
   0.22080000000000  -0.08340000000000   0.31300000000000  Br   (4e)
  -0.22080000000000   0.41660000000000   0.18700000000000  Br   (4e)
  -0.22080000000000   0.08340000000000  -0.31300000000000  Br   (4e)
   0.22080000000000   0.58340000000000   0.81300000000000  Br   (4e)
   0.27670000000000   0.09540000000000   0.69010000000000  Br   (4e)
  -0.27670000000000   0.59540000000000  -0.19010000000000  Br   (4e)
  -0.27670000000000  -0.09540000000000  -0.69010000000000  Br   (4e)
   0.27670000000000   0.40460000000000   1.19010000000000  Br   (4e)
   0.58430000000000   0.13530000000000   0.06130000000000  H2O  (4e)
  -0.58430000000000   0.63530000000000   0.43870000000000  H2O  (4e)
  -0.58430000000000  -0.13530000000000  -0.06130000000000  H2O  (4e)
   0.58430000000000   0.36470000000000   0.56130000000000  H2O  (4e)
   0.01710000000000   0.36400000000000   0.55120000000000  H2O  (4e)
  -0.01710000000000   0.86400000000000  -0.05120000000000  H2O  (4e)
  -0.01710000000000  -0.36400000000000  -0.55120000000000  H2O  (4e)
   0.01710000000000   0.13600000000000   1.05120000000000  H2O  (4e)
   0.26710000000000   0.47230000000000   0.51810000000000  K    (4e)
  -0.26710000000000   0.97230000000000  -0.01810000000000  K    (4e)
  -0.26710000000000  -0.47230000000000  -0.51810000000000  K    (4e)
   0.26710000000000   0.02770000000000   1.01810000000000  K    (4e)
```

Anhydrous KAuBr$_4$: AB4C_mP24_14_ab_4e_e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'AuBr4K'
_chemical_formula_sum 'Au Br4 K'

loop_
_publ_author_name
 'H. Omrani'
 'R. Welter'
 'R. Vangelisti'
_journal_name_full_name
;
 Acta Crystallographica Section C: Structural Chemistry
;
_journal_volume 55
_journal_year 1999
_journal_page_first 13
_journal_page_last 14
_publ_Section_title
;
 Potassium tetrabromoaurate(III)
;

# Found in Sodium tetrabromoaurate(III) dihydrate, 2001

_aflow_title 'Anhydrous KAuBr$_{4}$ Structure'
_aflow_proto 'AB4C_mP24_14_ab_4e_e'
_aflow_params 'a,b/a,c/a,\beta,x_{3},y_{3},z_{3},x_{4},y_{4},z_{4},x_{5
      ↪ },y_{5},z_{5},x_{6},y_{6},z_{6},x_{7},y_{7},z_{7}'
_aflow_params_values '9.0306,0.736208003898,1.41740305185,96.88,0.86659,
      ↪ 0.2934,0.04766,0.09492,-0.0538,0.18234,0.5056,0.1517,0.1707,
      ↪ 0.6635,0.7336,0.07207,0.2128,0.4492,0.1813'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mP24'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

_cell_length_a    9.03060
_cell_length_b    6.64840
_cell_length_c    12.80000
_cell_angle_alpha 90.00000
_cell_angle_beta  96.88000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z+1/2
3 -x,-y,-z
4 x,-y+1/2,z+1/2

loop_
_atom_site_label
```

```
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Au1 Au 2 a 0.00000 0.00000 0.00000 1.00000
Au2 Au 2 b 0.50000 0.00000 0.00000 1.00000
Br1 Br 4 e 0.86659 0.29340 0.04766 1.00000
Br2 Br 4 e 0.09492 -0.05380 0.18234 1.00000
Br3 Br 4 e 0.50560 0.15170 0.17070 1.00000
Br4 Br 4 e 0.66350 0.73360 0.07207 1.00000
K1 K 4 e 0.21280 0.44920 0.18130 1.00000
```

Anhydrous $KAuBr_4$: AB4C_mP24_14_ab_4e_e - POSCAR

```
AB4C_mP24_14_ab_4e_e & a,b/a,c/a,beta,x3,y3,z3,x4,y4,z4,x5,y5,z5,x6,y6,
    ↪ z6,x7,y7,z7 --params=9.0306,0.736208003898,1.41740305185,96.88,
    ↪ 0.86659,0.2934,0.04766,0.09492,-0.0538,0.18234,0.5056,0.1517,
    ↪ 0.1707,0.6635,0.7336,0.07207,0.2128,0.4492,0.1813 & P2_{1}/c C_
    ↪ {2h}^{5} #14 (abe^5) & mP24 & None & AuBr4K & AuBr4K & H.
    ↪ Omrani and R. Welter and R. Vangelisti, Acta Crystallogr. C 55,
    ↪ 13-14 (1999)
   1.00000000000000
   9.03060000000000    0.00000000000000    0.00000000000000
   0.00000000000000    6.64840000000000    0.00000000000000
  -1.53331576111014    0.00000000000000   12.70782997906140
     Au     Br      K
      4     16      4
Direct
   0.00000000000000    0.00000000000000    0.00000000000000   Au   (2a)
   0.00000000000000    0.50000000000000    0.50000000000000   Au   (2a)
   0.50000000000000    0.00000000000000    0.00000000000000   Au   (2b)
   0.50000000000000    0.50000000000000    0.50000000000000   Au   (2b)
   0.86659000000000    0.29340000000000    0.04766000000000   Br   (4e)
  -0.86659000000000    0.79340000000000    0.45234000000000   Br   (4e)
  -0.86659000000000   -0.29340000000000   -0.04766000000000   Br   (4e)
   0.86659000000000    0.20660000000000    0.54766000000000   Br   (4e)
   0.09492000000000   -0.05380000000000    0.18234000000000   Br   (4e)
  -0.09492000000000    0.44620000000000    0.31766000000000   Br   (4e)
  -0.09492000000000    0.05380000000000   -0.18234000000000   Br   (4e)
   0.09492000000000    0.55380000000000    0.68234000000000   Br   (4e)
   0.50560000000000    0.15170000000000    0.17070000000000   Br   (4e)
  -0.50560000000000    0.65170000000000    0.32930000000000   Br   (4e)
  -0.50560000000000   -0.15170000000000   -0.17070000000000   Br   (4e)
   0.50560000000000    0.34830000000000    0.67070000000000   Br   (4e)
   0.66350000000000    0.73360000000000    0.07207000000000   Br   (4e)
  -0.66350000000000    1.23360000000000    0.42793000000000   Br   (4e)
  -0.66350000000000   -0.73360000000000   -0.07207000000000   Br   (4e)
   0.66350000000000   -0.23360000000000    0.57207000000000   Br   (4e)
   0.21280000000000    0.44920000000000    0.18130000000000   K    (4e)
  -0.21280000000000    0.94920000000000    0.31870000000000   K    (4e)
  -0.21280000000000   -0.44920000000000   -0.18130000000000   K    (4e)
   0.21280000000000    0.05080000000000    0.68130000000000   K    (4e)
```

Ammonium Persulfate $[(NH_4)_2S_2O_8, K4_1]$: AB4C_mP24_14_e_4e_e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Ammonium persulfate'
_chemical_formula_sum '(NH4) O4 S'

loop_
_publ_author_name
 'B. K. Sivertsen'
 'H. Sorum'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 130
_journal_year 1969
_journal_page_first 449
_journal_page_last 460
_publ_Section_title
;
 A reinvestigation of the crystal structure of ammonium persulfate, (
    ↪ NH$_{4}$)$_{2}$S$_{2}$O$_{8}$
;

_aflow_title 'Ammonium Persulfate [(NH$_{4}$)$_{2}$S$_{2}$O$_{8}$, $K4_{
    ↪ 1}$] Structure'
_aflow_proto 'AB4C_mP24_14_e_4e_e'
_aflow_params 'a,b/a,c/a,\beta,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3
    ↪ },y_{3},z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},
    ↪ z_{6}'
_aflow_params_values '7.83,1.02273307791,1.21392081737,139.39,0.0899,
    ↪ 0.8839,0.2331,0.4059,0.5186,0.3818,0.343,0.217,0.3158,0.2414,
    ↪ 0.4109,0.0667,0.6985,0.351,0.4181,0.4242,0.3572,0.2869'
_aflow_Strukturbericht '$K4_{1}$'
_aflow_Pearson 'mP24'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

_cell_length_a      7.83000
_cell_length_b      8.00800
_cell_length_c      9.50500
_cell_angle_alpha  90.00000
_cell_angle_beta   139.39000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z+1/2
3 -x,-y,-z
4 x,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
NH41 NH4 4 e 0.08990 0.88390 0.23310 1.00000
O1 O 4 e 0.40590 0.51860 0.38180 1.00000
O2 O 4 e 0.34300 0.21700 0.31580 1.00000
O3 O 4 e 0.24140 0.41090 0.06670 1.00000
O4 O 4 e 0.69850 0.35100 0.41810 1.00000
S1 S 4 e 0.42420 0.35720 0.28690 1.00000
```

Ammonium Persulfate $[(NH_4)_2S_2O_8, K4_1]$: AB4C_mP24_14_e_4e_e - POSCAR

```
AB4C_mP24_14_e_4e_e & a,b/a,c/a,beta,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,z4
    ↪ ,x5,y5,z5,x6,y6,z6 --params=7.83,1.02273307791,1.21392081737,
    ↪ 139.39,0.0899,0.8839,0.2331,0.4059,0.5186,0.3818,0.343,0.217,
    ↪ 0.3158,0.2414,0.4109,0.0667,0.6985,0.351,0.4181,0.4242,0.3572,
    ↪ 0.2869 & P2_{1}/c C_{2h}^{5} #14 (e^6) & mP24 & $K4_{1}$ & (NH4
    ↪ )O4S & Ammonium persulfate & B. K. Sivertsen and H. Sorum,
    ↪ Zeitschrift f{"u}r Kristallographie - Crystalline Materials 130
    ↪ , 449-460 (1969)
   1.00000000000000
   7.83000000000000    0.00000000000000    0.00000000000000
   0.00000000000000    8.00800000000000    0.00000000000000
  -7.21579407388266    0.00000000000000    6.18686842298427
    NH4      O      S
      4     16      4
Direct
   0.08990000000000    0.88390000000000    0.23310000000000   NH4   (4e)
  -0.08990000000000    1.38390000000000    0.26690000000000   NH4   (4e)
  -0.08990000000000   -0.88390000000000   -0.23310000000000   NH4   (4e)
   0.08990000000000   -0.38390000000000    0.73310000000000   NH4   (4e)
   0.40590000000000    0.51860000000000    0.38180000000000    O    (4e)
  -0.40590000000000    1.01860000000000    0.11820000000000    O    (4e)
  -0.40590000000000   -0.51860000000000   -0.38180000000000    O    (4e)
   0.40590000000000   -0.01860000000000    0.88180000000000    O    (4e)
   0.34300000000000    0.21700000000000    0.31580000000000    O    (4e)
  -0.34300000000000    0.71700000000000    0.18420000000000    O    (4e)
  -0.34300000000000   -0.21700000000000   -0.31580000000000    O    (4e)
   0.34300000000000    0.28300000000000    0.81580000000000    O    (4e)
   0.24140000000000    0.41090000000000    0.06670000000000    O    (4e)
  -0.24140000000000    0.91090000000000    0.43330000000000    O    (4e)
  -0.24140000000000   -0.41090000000000   -0.06670000000000    O    (4e)
   0.24140000000000    0.08910000000000    0.56670000000000    O    (4e)
   0.69850000000000    0.35100000000000    0.41810000000000    O    (4e)
  -0.69850000000000    0.85100000000000    0.08190000000000    O    (4e)
  -0.69850000000000   -0.35100000000000   -0.41810000000000    O    (4e)
   0.69850000000000    0.14900000000000    0.91810000000000    O    (4e)
   0.42420000000000    0.35720000000000    0.28690000000000    S    (4e)
  -0.42420000000000    0.85720000000000    0.21310000000000    S    (4e)
  -0.42420000000000   -0.35720000000000   -0.28690000000000    S    (4e)
   0.42420000000000    0.14280000000000    0.78690000000000    S    (4e)
```

Monazite ($LaPO_4$): AB4C_mP24_14_e_4e_e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Monazite'
_chemical_formula_sum 'La O4 P'

loop_
_publ_author_name
 'Y. Ni'
 'J. M. Hughes'
 'A. N. Mariano'
_journal_name_full_name
;
 American Mineralogist
;
_journal_volume 80
_journal_year 1995
_journal_page_first 21
_journal_page_last 26
_publ_Section_title
;
 Crystal chemistry of the monazite and xenotime structures
;

# Found in Gasparite-(La), La(AsO$_{4}$), a new mineral from Mn ores of
    ↪ the Ushkatyn-III deposit, Central Kazakhstan, and metamorphic
    ↪ rocks of the Wanni glacier, Switzerland, 2019

_aflow_title 'Monazite (LaPO$_{4}$) Structure'
_aflow_proto 'AB4C_mP24_14_e_4e_e'
_aflow_params 'a,b/a,c/a,\beta,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3
    ↪ },y_{3},z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},
    ↪ z_{6}'
_aflow_params_values '6.5034,1.08720054126,1.27324630193,126.58575,
    ↪ 0.18086,0.16033,0.28154,0.8026,0.0077,0.2503,0.8835,0.3315,
```

```
      ↪ 0.3799,0.673,0.1071,0.4748,0.4176,0.2168,0.1277,0.6926,0.1639,
      ↪ 0.3047'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mP24'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

_cell_length_a 6.50340
_cell_length_b 7.07050
_cell_length_c 8.28043
_cell_angle_alpha 90.00000
_cell_angle_beta 126.58575
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z+1/2
3 -x,-y,-z
4 x,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
La1 La 4 e 0.18086 0.16033 0.28154 1.00000
O1 O 4 e 0.80260 0.00770 0.25030 1.00000
O2 O 4 e 0.88350 0.33150 0.37990 1.00000
O3 O 4 e 0.67300 0.10710 0.47480 1.00000
O4 O 4 e 0.41760 0.21680 0.12770 1.00000
P1 P 4 e 0.69260 0.16390 0.30470 1.00000
```

Monazite ($LaPO_4$): AB4C_mP24_14_e_4e_e - POSCAR

```
AB4C_mP24_14_e_4e_e & a,b/a,c/a,beta,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,z4
      ↪ ,x5,y5,z5,x6,y6,z6 --params=6.5034,1.08720054126,1.27324630193,
      ↪ 126.58575,0.18086,0.16033,0.28154,0.8026,0.0077,0.2503,0.8835,
      ↪ 0.3315,0.3799,0.673,0.1071,0.4748,0.4176,0.2168,0.1277,0.6926,
      ↪ 0.1639,0.3047 & P2_{1}/c C_{2h}^{5} #14 (e^6) & mP24 & None &
      ↪ LaO4P & Monazite & Y. Ni and J. M. Hughes and A. N. Mariano,
      ↪ Am. Mineral. 80, 21-26 (1995)
   1.00000000000000
   6.50340000000000    0.00000000000000    0.00000000000000
   0.00000000000000    7.07050000000000    0.00000000000000
  -4.93534484962521    0.00000000000000    6.64890157846978
     La      O      P
      4     16      4
Direct
   0.18086000000000    0.16033000000000    0.28154000000000   La   (4e)
  -0.18086000000000    0.66033000000000    0.21846000000000   La   (4e)
  -0.18086000000000   -0.16033000000000   -0.28154000000000   La   (4e)
   0.18086000000000    0.33967000000000    0.78154000000000   La   (4e)
   0.80260000000000    0.00770000000000    0.25030000000000   O    (4e)
  -0.80260000000000    0.50770000000000    0.24970000000000   O    (4e)
  -0.80260000000000   -0.00770000000000   -0.25030000000000   O    (4e)
   0.80260000000000    0.49230000000000    0.75030000000000   O    (4e)
   0.88350000000000    0.33150000000000    0.37990000000000   O    (4e)
  -0.88350000000000    0.83150000000000    0.12010000000000   O    (4e)
  -0.88350000000000   -0.33150000000000   -0.37990000000000   O    (4e)
   0.88350000000000    0.16850000000000    0.87990000000000   O    (4e)
   0.67300000000000    0.10710000000000    0.47480000000000   O    (4e)
  -0.67300000000000    0.60710000000000    0.02520000000000   O    (4e)
  -0.67300000000000   -0.10710000000000   -0.47480000000000   O    (4e)
   0.67300000000000    0.39290000000000    0.97480000000000   O    (4e)
   0.41760000000000    0.21680000000000    0.12770000000000   O    (4e)
  -0.41760000000000    0.71680000000000    0.37230000000000   O    (4e)
  -0.41760000000000   -0.21680000000000   -0.12770000000000   O    (4e)
   0.41760000000000    0.28320000000000    0.62770000000000   O    (4e)
   0.69260000000000    0.16390000000000    0.30470000000000   P    (4e)
  -0.69260000000000    0.66390000000000    0.19530000000000   P    (4e)
  -0.69260000000000   -0.16390000000000   -0.30470000000000   P    (4e)
   0.69260000000000    0.33610000000000    0.80470000000000   P    (4e)
```

Sr_2MnTeO_6: AB6C2D_mP20_14_a_3e_e_d - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'MnO6Sr2Te'
_chemical_formula_sum 'Mn O6 Sr2 Te'

loop_
_publ_author_name
 'L. {Ortega-San Martin}'
 'J. P. Chapman'
 'E. {Hern\\'{a}ndez-Bocanegra}'
 'M. Insausti'
 'M. I. Arriortua'
 'T. Rojo'
_journal_name_full_name
;
 Journal of Physics: Condensed Matter
;
_journal_volume 16
_journal_year 2004
_journal_page_first 3879
_journal_page_last 3888
_publ_Section_title
;
 Structural phase transitions in the ordered double perovskite Sr${_2}
      ↪ $MnTeO$_{6}$
;

_aflow_title 'Sr$_{2}$MnTeO$_{6}$ Structure'
_aflow_proto 'AB6C2D_mP20_14_a_3e_e_d'
_aflow_params 'a,b/a,c/a,\beta,x_{3},y_{3},z_{3},x_{4},y_{4},z_{4},x_{5
      ↪ },y_{5},z_{5},x_{6},y_{6},z_{6}'
_aflow_params_values '5.7009,0.995807679489,1.72670455542,125.30487,
      ↪ 0.2103,-0.0065,0.265,0.3193,0.7621,0.0287,0.7308,0.7126,-0.0283
      ↪ ,0.2523,-0.01532,0.7508'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mP20'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

_cell_length_a 5.70090
_cell_length_b 5.67700
_cell_length_c 9.84377
_cell_angle_alpha 90.00000
_cell_angle_beta 125.30487
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z+1/2
3 -x,-y,-z
4 x,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Mn1 Mn 2 a 0.00000 0.00000 0.00000 1.00000
Te1 Te 2 d 0.50000 0.00000 0.50000 1.00000
O1 O 4 e 0.21030 -0.00650 0.26500 1.00000
O2 O 4 e 0.31930 0.76210 0.02870 1.00000
O3 O 4 e 0.73080 0.71260 -0.02830 1.00000
Sr1 Sr 4 e 0.25230 -0.01532 0.75080 1.00000
```

Sr_2MnTeO_6: AB6C2D_mP20_14_a_3e_e_d - POSCAR

```
AB6C2D_mP20_14_a_3e_e_d & a,b/a,c/a,beta,x3,y3,z3,x4,y4,z4,x5,y5,z5,x6,
      ↪ y6,z6 --params=5.7009,0.995807679489,1.72670455542,125.30487,
      ↪ 0.2103,-0.0065,0.265,0.3193,0.7621,0.0287,0.7308,0.7126,-0.0283
      ↪ ,0.2523,-0.01532,0.7508 & P2_{1}/c C_{2h}^{5} #14 (ade^4) &
      ↪ mP20 & None & MnO6Sr2Te & MnO6Sr2Te & L. {Ortega-San Martin} et
      ↪ al., J. Phys.: Condens. Matter 16, 3879-3888 (2004)
   1.00000000000000
   5.70090000000000    0.00000000000000    0.00000000000000
   0.00000000000000    5.67700000000000    0.00000000000000
  -5.68898038583372    0.00000000000000    8.03338720481586
     Mn      O     Sr     Te
      2     12      4      2
Direct
   0.00000000000000    0.00000000000000    0.00000000000000   Mn   (2a)
   0.00000000000000    0.50000000000000    0.50000000000000   Mn   (2a)
   0.21030000000000   -0.00650000000000    0.26500000000000   O    (4e)
  -0.21030000000000    0.49350000000000    0.23500000000000   O    (4e)
  -0.21030000000000    0.00650000000000   -0.26500000000000   O    (4e)
   0.21030000000000    0.50650000000000    0.76500000000000   O    (4e)
   0.31930000000000    0.76210000000000    0.02870000000000   O    (4e)
  -0.31930000000000    1.26210000000000    0.47130000000000   O    (4e)
  -0.31930000000000   -0.76210000000000   -0.02870000000000   O    (4e)
   0.31930000000000   -0.26210000000000    0.52870000000000   O    (4e)
   0.73080000000000    0.71260000000000   -0.02830000000000   O    (4e)
  -0.73080000000000    1.21260000000000    0.52830000000000   O    (4e)
  -0.73080000000000   -0.71260000000000    0.02830000000000   O    (4e)
   0.73080000000000   -0.21260000000000    0.47170000000000   O    (4e)
   0.25230000000000   -0.01532000000000    0.75080000000000   Sr   (4e)
  -0.25230000000000    0.48468000000000   -0.25080000000000   Sr   (4e)
  -0.25230000000000    0.01532000000000   -0.75080000000000   Sr   (4e)
   0.25230000000000    0.51532000000000    1.25080000000000   Sr   (4e)
   0.50000000000000    0.00000000000000    0.50000000000000   Te   (2d)
   0.50000000000000    0.50000000000000    0.00000000000000   Te   (2d)
```

Cryolite (Na_3AlF_6, $J2_6$): AB6C3_mP20_14_a_3e_de - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Cryolite'
_chemical_formula_sum 'Al F6 Na3'

loop_
_publ_author_name
 'H. Yang'
 'S. Ghose'
 'D. M. Hatch'
_journal_name_full_name
;
 Physics and Chemistry of Minerals
;
_journal_volume 19
```

```
_journal_year 1993
_journal_page_first 528
_journal_page_last 544
_publ_Section_title
;
 Ferroelastic phase transition in cryolite, Na$_{3}$AlF$_{6}$, a mixed
      ↪ fluoride perovskite: High temperature single crystal X-ray
      ↪ diffraction study and symmetry analysis of the transition
      ↪ mechanism
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'Cryolite (Na$_{3}$AlF$_{6}$, $J2_{6}$) Structure'
_aflow_proto 'AB6C3_mP20_14_a_3e_de'
_aflow_params 'a,b/a,c/a,\beta,x_{3},y_{3},z_{3},x_{4},y_{4},z_{4},x_{5
      ↪ },y_{5},z_{5},x_{6},y_{6},z_{6}'
_aflow_params_values '5.4139,1.03459613218,1.74764587451,124.72045,
      ↪ 0.8828,-0.0442,0.7812,0.6828,0.6733,0.455,0.2239,0.7313,0.0617,
      ↪ 0.7353,0.5508,0.7474'
_aflow_Strukturbericht '$J2_{6}$'
_aflow_Pearson 'mP20'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

_cell_length_a     5.41390
_cell_length_b     5.60120
_cell_length_c     9.46158
_cell_angle_alpha  90.00000
_cell_angle_beta   124.72045
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z+1/2
3 -x,-y,-z
4 x,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Al1 Al 2 a 0.00000 0.00000 0.00000 1.00000
Na1 Na 2 d 0.50000 0.00000 0.50000 1.00000
F1 F 4 e 0.88280 -0.04420 0.78120 1.00000
F2 F 4 e 0.68280 0.67330 0.45500 1.00000
F3 F 4 e 0.22390 0.73130 0.06170 1.00000
Na2 Na 4 e 0.73530 0.55080 0.74740 1.00000
```

Cryolite (Na_3AlF_6, $J2_6$): AB6C3_mP20_14_a_3e_de - POSCAR

```
AB6C3_mP20_14_a_3e_de & a,b/a,c/a,beta,x3,y3,z3,x4,y4,z4,x5,y5,z5,x6,y6,
      ↪ z6 --params=5.4139,1.03459613218,1.74764587451,124.72045,0.8828
      ↪ ,-0.0442,0.7812,0.6828,0.6733,0.455,0.2239,0.7313,0.0617,0.7353
      ↪ ,0.5508,0.7474 & P2_{1}/c C_{2h}^{5} #14 (ade^4) & mP20 & $J2_{
      ↪ 6}$ & AlF6Na3 & Cryolite & H. Yang and S. Ghose and D. M. Hatch
      ↪ , Phys. Chem. Miner. 19, 528-544 (1993)
   1.00000000000000
   5.41390000000000    0.00000000000000    0.00000000000000
   0.00000000000000    5.60120000000000    0.00000000000000
  -5.38905981158747    0.00000000000000    7.77685864880755
     Al      F      Na
      2     12       6
Direct
   0.00000000000000    0.00000000000000    0.00000000000000    Al    (2a)
   0.00000000000000    0.50000000000000    0.50000000000000    Al    (2a)
   0.88280000000000   -0.04420000000000    0.78120000000000    F     (4e)
  -0.88280000000000    0.45580000000000   -0.28120000000000    F     (4e)
  -0.88280000000000    0.04420000000000   -0.78120000000000    F     (4e)
   0.88280000000000    0.54420000000000    1.28120000000000    F     (4e)
   0.68280000000000    0.67330000000000    0.45500000000000    F     (4e)
  -0.68280000000000    1.17330000000000    0.04500000000000    F     (4e)
  -0.68280000000000   -0.67330000000000   -0.45500000000000    F     (4e)
   0.68280000000000   -0.17330000000000    0.95500000000000    F     (4e)
   0.22390000000000    0.73130000000000    0.06170000000000    F     (4e)
  -0.22390000000000    1.23130000000000    0.43830000000000    F     (4e)
  -0.22390000000000   -0.73130000000000   -0.06170000000000    F     (4e)
   0.22390000000000   -0.23130000000000    0.56170000000000    F     (4e)
   0.50000000000000    0.00000000000000    0.50000000000000    Na    (2d)
   0.50000000000000    0.50000000000000    0.00000000000000    Na    (2d)
   0.73530000000000    0.55080000000000    0.74740000000000    Na    (4e)
  -0.73530000000000    1.05080000000000   -0.24740000000000    Na    (4e)
  -0.73530000000000   -0.55080000000000   -0.74740000000000    Na    (4e)
   0.73530000000000   -0.05080000000000    1.24740000000000    Na    (4e)
```

KNO_2 III: ABC2_mP16_14_e_e_2e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Potassium nitrite'
_chemical_formula_sum 'K N O2'

loop_
_publ_author_name
 'N. {Onoda-Yamamuro}'
 'H. Honda'
 'R. Ikeda'
 'O. Yamamuro'
 'T. Matsuo'
 'K. Oikawa'
 'T. Kamiyama'
 'F. Izumi'
_journal_name_full_name
;
 Journal of Physics: Condensed Matter
;
_journal_volume 10
_journal_year 1998
_journal_page_first 3341
_journal_page_last 3351
_publ_Section_title
;
 Neutron powder diffraction study of the low-temperature phases of KNO$_
      ↪ {2}$
;

# Found in Order-disorder phase transitions in KNO$_{2}$, CsNO$_{2}$,
      ↪ and TlNO$_{2}$ crystals: A molecular dynamics study, 2001

_aflow_title 'KNO$_{2}$ III Structure'
_aflow_proto 'ABC2_mP16_14_e_e_2e'
_aflow_params 'a,b/a,c/a,\beta,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3
      ↪ },y_{3},z_{3},x_{4},y_{4},z_{4}'
_aflow_params_values '4.401,2.18057259714,1.58638945694,108.274,0.4105,
      ↪ 0.1407,0.2101,-0.0732,0.4107,0.1839,0.2034,0.4127,0.3067,0.8385
      ↪ ,0.2999,0.0865'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mP16'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

_cell_length_a     4.40100
_cell_length_b     9.59670
_cell_length_c     6.98170
_cell_angle_alpha  90.00000
_cell_angle_beta   108.27400
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z+1/2
3 -x,-y,-z
4 x,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
K1 K 4 e 0.41050 0.14070 0.21010 1.00000
N1 N 4 e -0.07320 0.41070 0.18390 1.00000
O1 O 4 e 0.20340 0.41270 0.30670 1.00000
O2 O 4 e 0.83850 0.29990 0.08650 1.00000
```

KNO_2 III: ABC2_mP16_14_e_e_2e - POSCAR

```
ABC2_mP16_14_e_e_2e & a,b/a,c/a,beta,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,z4
      ↪ --params=4.401,2.18057259714,1.58638945694,108.274,0.4105,
      ↪ 0.1407,0.2101,-0.0732,0.4107,0.1839,0.2034,0.4127,0.3067,0.8385
      ↪ ,0.2999,0.0865 & P2_{1}/c C_{2h}^{5} #14 (e^4) & mP16 & None &
      ↪ KNO2 & Potassium nitrite & N. {Onoda-Yamamuro} et al., J. Phys.
      ↪ : Condens. Matter 10, 3341-3351 (1998)
   1.00000000000000
   4.40100000000000    0.00000000000000    0.00000000000000
   0.00000000000000    9.59670000000000    0.00000000000000
  -2.18919293920523    0.00000000000000    6.62959796404986
     K      N      O
     4      4      8
Direct
   0.41050000000000    0.14070000000000    0.21010000000000    K     (4e)
  -0.41050000000000    0.64070000000000    0.28990000000000    K     (4e)
  -0.41050000000000   -0.14070000000000   -0.21010000000000    K     (4e)
   0.41050000000000    0.35930000000000    0.71010000000000    K     (4e)
  -0.07320000000000    0.41070000000000    0.18390000000000    N     (4e)
   0.07320000000000    0.91070000000000    0.31610000000000    N     (4e)
   0.07320000000000   -0.41070000000000   -0.18390000000000    N     (4e)
  -0.07320000000000    0.08930000000000    0.68390000000000    N     (4e)
   0.20340000000000    0.41270000000000    0.30670000000000    O     (4e)
  -0.20340000000000    0.91270000000000    0.19330000000000    O     (4e)
  -0.20340000000000   -0.41270000000000   -0.30670000000000    O     (4e)
   0.20340000000000    0.08730000000000    0.80670000000000    O     (4e)
   0.83850000000000    0.29990000000000    0.08650000000000    O     (4e)
  -0.83850000000000    0.79990000000000    0.41350000000000    O     (4e)
  -0.83850000000000   -0.29990000000000   -0.08650000000000    O     (4e)
   0.83850000000000    0.20010000000000    0.58650000000000    O     (4e)
```

Manganite (γ-MnO(OH), $E0_6$): ABC2_mP16_14_e_e_2e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Manganite'
```

```
_chemical_formula_sum 'H Mn O2'

loop_
_publ_author_name
 'T. Kohler'
 'T. Armbruster'
 'E. Libowitzky'
_journal_name_full_name
;
 Journal of Solid State Chemistry
;
_journal_volume 133
_journal_year 1997
_journal_page_first 486
_journal_page_last 500
_publ_Section_title
;
 Hydrogen Bonding and Jahn-Teller Distortion in Groutite, $\alpha$-MnOOH
      ↪ , and Manganite, $\gamma$-MnOOH, and Their Relations to the
      ↪ Manganese Dioxides Ramsdellite and Pyrolusite
;

_aflow_title 'Manganite ($\gamma$-MnO(OH), $E0_{6}$) Structure'
_aflow_proto 'ABC2_mP16_14_e_e_2e'
_aflow_params 'a,b/a,c/a,\beta,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3
      ↪ },y_{3},z_{3},x_{4},y_{4},z_{4}'
_aflow_params_values '5.304,0.994909502262,1.0,114.38,0.284,0.027,0.725,
      ↪ 0.76316,0.01033,0.75464,0.3749,0.1238,0.6279,0.8752,0.1256,
      ↪ 0.1206'
_aflow_Strukturbericht '$E0_{6}$'
_aflow_Pearson 'mP16'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

_cell_length_a      5.30400
_cell_length_b      5.27700
_cell_length_c      5.30400
_cell_angle_alpha   90.00000
_cell_angle_beta    114.38000
_cell_angle_gamma   90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z+1/2
3 -x,-y,-z
4 x,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
H1 H 4 e 0.28400 0.02700 0.72500 1.00000
Mn1 Mn 4 e 0.76316 0.01033 0.75464 1.00000
O1 O 4 e 0.37490 0.12380 0.62790 1.00000
O2 O 4 e 0.87520 0.12560 0.12060 1.00000
```

Manganite (γ-MnO(OH), $E0_6$): ABC2_mP16_14_e_e_2e - POSCAR

```
ABC2_mP16_14_e_e_2e & a,b/a,c/a,beta,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,z4
      ↪  --params=5.304,0.994909502262,1.0,114.38,0.284,0.027,0.725,
      ↪ 0.76316,0.01033,0.75464,0.3749,0.1238,0.6279,0.8752,0.1256,
      ↪ 0.1206 & P2_{1}/c C_{2h}^{5} #14 (e^4) & mP16 & $E0_{6}$ &
      ↪ MnHO2 & Manganite & T. Kohler and T. Armbruster and E.
      ↪ Libowitzky, J. Solid State Chem. 133, 486-500 (1997)
   1.000000000000000
   5.304000000000000    0.000000000000000    0.000000000000000
   0.000000000000000    5.277000000000000    0.000000000000000
  -2.189419681377070    0.000000000000000    4.831030682866620
     H    Mn     O
     4     4     8
Direct
   0.284000000000000    0.027000000000000    0.725000000000000     H    (4e)
  -0.284000000000000    0.527000000000000   -0.225000000000000     H    (4e)
  -0.284000000000000   -0.027000000000000   -0.725000000000000     H    (4e)
   0.284000000000000    0.473000000000000    1.225000000000000     H    (4e)
   0.763160000000000    0.010330000000000    0.754640000000000    Mn    (4e)
  -0.763160000000000    0.510330000000000   -0.254640000000000    Mn    (4e)
  -0.763160000000000   -0.010330000000000   -0.754640000000000    Mn    (4e)
   0.763160000000000    0.489670000000000    1.254640000000000    Mn    (4e)
   0.374900000000000    0.123800000000000    0.627900000000000     O    (4e)
  -0.374900000000000    0.623800000000000   -0.127900000000000     O    (4e)
  -0.374900000000000   -0.123800000000000   -0.627900000000000     O    (4e)
   0.374900000000000    0.376200000000000    1.127900000000000     O    (4e)
   0.875200000000000    0.125600000000000    0.120600000000000     O    (4e)
  -0.875200000000000    0.625600000000000    0.379400000000000     O    (4e)
  -0.875200000000000   -0.125600000000000   -0.120600000000000     O    (4e)
   0.875200000000000    0.374400000000000    0.620600000000000     O    (4e)
```

$AgMnO_4$ ($H0_9$): ABC4_mP24_14_e_e_4e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'AgMnO4'
_chemical_formula_sum 'Ag Mn O4'

loop_
_publ_author_name
 'E. G. Boonstra'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 24
_journal_year 1968
_journal_page_first 1053
_journal_page_last 1062
_publ_Section_title
;
 The Crystal Structure of Silver Permanganate
;

_aflow_title 'AgMnO$_{4}$ ($H0_{9}$) Structure'
_aflow_proto 'ABC4_mP24_14_e_e_4e'
_aflow_params 'a,b/a,c/a,\beta,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3
      ↪ },y_{3},z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},
      ↪ z_{6}'
_aflow_params_values '5.64,1.47695035461,1.57941666667,126.99641,0.0792,
      ↪ 0.4672,0.8353,0.5965,0.3132,0.3392,0.7908,0.1508,0.4079,0.2708,
      ↪ 0.2767,0.2847,0.5599,0.381,0.1555,0.7557,0.441,0.5037'
_aflow_Strukturbericht '$H0_{9}$'
_aflow_Pearson 'mP24'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

_cell_length_a      5.64000
_cell_length_b      8.33000
_cell_length_c      8.90791
_cell_angle_alpha   90.00000
_cell_angle_beta    126.99641
_cell_angle_gamma   90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z+1/2
3 -x,-y,-z
4 x,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ag1 Ag 4 e 0.07920 0.46720 0.83530 1.00000
Mn1 Mn 4 e 0.59650 0.31320 0.33920 1.00000
O1 O 4 e 0.79080 0.15080 0.40790 1.00000
O2 O 4 e 0.27080 0.27670 0.28470 1.00000
O3 O 4 e 0.55990 0.38100 0.15550 1.00000
O4 O 4 e 0.75570 0.44100 0.50370 1.00000
```

$AgMnO_4$ ($H0_9$): ABC4_mP24_14_e_e_4e - POSCAR

```
ABC4_mP24_14_e_e_4e & a,b/a,c/a,beta,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,z4
      ↪ ,x5,y5,z5,x6,y6,z6 --params=5.64,1.47695035461,1.57941666667,
      ↪ 126.99641,0.0792,0.4672,0.8353,0.5965,0.3132,0.3392,0.7908,
      ↪ 0.1508,0.4079,0.2708,0.2767,0.2847,0.5599,0.381,0.1555,0.7557,
      ↪ 0.441,0.5037 & P2_{1}/c C_{2h}^{5} #14 (e^6) & mP24 & $H0_{9}$
      ↪ & AgMnO4 & AgMnO4 & E. G. Boonstra, Acta Crystallogr. Sect. B
      ↪ Struct. Sci. 24, 1053-1062 (1968)
   1.000000000000000
   5.640000000000000    0.000000000000000    0.000000000000000
   0.000000000000000    8.330000000000000    0.000000000000000
  -5.360468529733267    0.000000000000000    7.114509132848970
    Ag     Mn     O
     4     4    16
Direct
   0.079200000000000    0.467200000000000    0.835300000000000    Ag    (4e)
  -0.079200000000000    0.967200000000000   -0.335300000000000    Ag    (4e)
  -0.079200000000000   -0.467200000000000   -0.835300000000000    Ag    (4e)
   0.079200000000000    0.032800000000000    1.335300000000000    Ag    (4e)
   0.596500000000000    0.313200000000000    0.339200000000000    Mn    (4e)
  -0.596500000000000    0.813200000000000    0.160800000000000    Mn    (4e)
  -0.596500000000000   -0.313200000000000   -0.339200000000000    Mn    (4e)
   0.596500000000000    0.186800000000000    0.839200000000000    Mn    (4e)
   0.790800000000000    0.150800000000000    0.407900000000000     O    (4e)
  -0.790800000000000    0.650800000000000    0.092100000000000     O    (4e)
  -0.790800000000000   -0.150800000000000   -0.407900000000000     O    (4e)
   0.790800000000000    0.349200000000000    0.907900000000000     O    (4e)
   0.270800000000000    0.276700000000000    0.284700000000000     O    (4e)
  -0.270800000000000    0.776700000000000    0.215300000000000     O    (4e)
  -0.270800000000000   -0.276700000000000   -0.284700000000000     O    (4e)
   0.270800000000000    0.223300000000000    0.784700000000000     O    (4e)
   0.559900000000000    0.381000000000000    0.155500000000000     O    (4e)
  -0.559900000000000    0.881000000000000    0.344500000000000     O    (4e)
  -0.559900000000000   -0.381000000000000   -0.155500000000000     O    (4e)
   0.559900000000000    0.119000000000000    0.655500000000000     O    (4e)
   0.755700000000000    0.441000000000000    0.503700000000000     O    (4e)
  -0.755700000000000    0.941000000000000   -0.003700000000000     O    (4e)
  -0.755700000000000   -0.441000000000000   -0.503700000000000     O    (4e)
   0.755700000000000    0.059000000000000    1.003700000000000     O    (4e)
```

Nahcolite ($NaHCO_3$, $G0_{12}$): ABCD3_mP24_14_e_e_e_3e - CIF

```
# CIF file
```

```
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Nacolite'
_chemical_formula_sum 'C H Na O3'

loop_
_publ_author_name
 'R. L. Sass'
 'R. F. Scheuerman'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 15
_journal_year 1962
_journal_page_first 77
_journal_page_last 81
_publ_Section_title
;
 The Crystal Structure of Sodium Bicarbonate
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'Nahcolite (NaHCO$_{3}$, $G0_{12}$) Structure'
_aflow_proto 'ABCD3_mP24_14_e_e_e_3e'
_aflow_params 'a,b/a,c/a,\beta,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3
    ↪ },y_{3},z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},
    ↪ z_{6}'
_aflow_params_values '3.51,2.76638176638,2.29344729345,111.85,0.2098,
    ↪ 0.237,-0.0768,0.74205,0.2539,0.1773,0.4274,0.0047,0.7145,0.1896
    ↪ ,0.3668,-0.0709,-0.0117,0.1629,0.7946,0.4958,0.1707,0.06'
_aflow_Strukturbericht '$G0_{12}$'
_aflow_Pearson 'mP24'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

_cell_length_a 3.51000
_cell_length_b 9.71000
_cell_length_c 8.05000
_cell_angle_alpha 90.00000
_cell_angle_beta 111.85000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z+1/2
3 -x,-y,-z
4 x,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
C1 C 4 e 0.20980 0.23700 -0.07680 1.00000
H1 H 4 e 0.74205 0.25390 0.17730 1.00000
Na1 Na 4 e 0.42740 0.00470 0.71450 1.00000
O1 O 4 e 0.18960 0.36680 -0.07090 1.00000
O2 O 4 e -0.01170 0.16290 0.79460 1.00000
O3 O 4 e 0.49580 0.17070 0.06000 1.00000
```

Nahcolite ($NaHCO_3$, $G0_{12}$): ABCD3_mP24_14_e_e_e_3e - POSCAR

```
ABCD3_mP24_14_e_e_e_3e & a,b/a,c/a,beta,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4
    ↪ ,z4,x5,y5,z5,x6,y6,z6 --params=3.51,2.76638176638,2.29344729345
    ↪ ,111.85,0.2098,0.237,-0.0768,0.74205,0.2539,0.1773,0.4274,
    ↪ 0.0047,0.7145,0.1896,0.3668,-0.0709,-0.0117,0.1629,0.7946,
    ↪ 0.4958,0.1707,0.06 & P2_{1}/c C_{2h}^{5} #14 (e^6) & mP24 &
    ↪ $G0_{12}$ & CHNaO3 & Nacolite & R. L. Sass and R. F. Scheuerman
    ↪ , Acta Cryst. 15, 77-81 (1962)
   1.00000000000000
   3.51000000000000    0.00000000000000    0.00000000000000
   0.00000000000000    9.71000000000000    0.00000000000000
  -2.99603250376187    0.00000000000000    7.47169922015082
     C     H     Na     O
     4     4     4     12
Direct
   0.20980000000000    0.23700000000000   -0.07680000000000    C    (4e)
  -0.20980000000000    0.73700000000000    0.57680000000000    C    (4e)
  -0.20980000000000   -0.23700000000000    0.07680000000000    C    (4e)
   0.20980000000000    0.26300000000000    0.42320000000000    C    (4e)
   0.74205000000000    0.25390000000000    0.17730000000000    H    (4e)
  -0.74205000000000    0.75390000000000    0.32270000000000    H    (4e)
  -0.74205000000000   -0.25390000000000   -0.17730000000000    H    (4e)
   0.74205000000000    0.24610000000000    0.67730000000000    H    (4e)
   0.42740000000000    0.00470000000000    0.71450000000000    Na   (4e)
  -0.42740000000000    0.50470000000000   -0.21450000000000    Na   (4e)
  -0.42740000000000   -0.00470000000000   -0.71450000000000    Na   (4e)
   0.42740000000000    0.49530000000000    1.21450000000000    Na   (4e)
   0.18960000000000    0.36680000000000   -0.07090000000000    O    (4e)
  -0.18960000000000    0.86680000000000    0.57090000000000    O    (4e)
  -0.18960000000000   -0.36680000000000    0.07090000000000    O    (4e)
   0.18960000000000    0.13320000000000    0.42910000000000    O    (4e)
  -0.01170000000000    0.16290000000000    0.79460000000000    O    (4e)
   0.01170000000000    0.66290000000000   -0.29460000000000    O    (4e)
   0.01170000000000   -0.16290000000000   -0.79460000000000    O    (4e)
  -0.01170000000000    0.33710000000000    1.29460000000000    O    (4e)
   0.49580000000000    0.17070000000000    0.06000000000000    O    (4e)
  -0.49580000000000    0.67070000000000    0.44000000000000    O    (4e)
  -0.49580000000000   -0.17070000000000   -0.06000000000000    O    (4e)
   0.49580000000000    0.32930000000000    0.56000000000000    O    (4e)
```

Cu(OH)Cl: ABCD_mP16_14_e_e_e_e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'ClCuHO'
_chemical_formula_sum 'Cl Cu H O'

loop_
_publ_author_name
 'Y. Cudennec'
 'A. Riou'
 'Y. G\\'{e}rault'
 'A. Lecerf'
_journal_name_full_name
;
 Journal of Solid State Chemistry
;
_journal_volume 151
_journal_year 2000
_journal_page_first 308
_journal_page_last 312
_publ_Section_title
;
 Synthesis and Crystal Structures of Cd(OH)Cl and Cu(OH)Cl and
    ↪ Relationship to Brucite Type
;

_aflow_title 'Cu(OH)Cl Structure'
_aflow_proto 'ABCD_mP16_14_e_e_e_e'
_aflow_params 'a,b/a,c/a,\beta,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3
    ↪ },y_{3},z_{3},x_{4},y_{4},z_{4}'
_aflow_params_values '6.2953,1.05871046654,0.882880879386,118.138,0.3115
    ↪ ,0.0907,0.1334,0.03201,0.11772,0.28577,0.674,0.128,0.528,0.8807
    ↪ ,0.1478,0.5318'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mP16'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

_cell_length_a 6.29530
_cell_length_b 6.66490
_cell_length_c 5.55800
_cell_angle_alpha 90.00000
_cell_angle_beta 118.13800
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z+1/2
3 -x,-y,-z
4 x,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cl1 Cl 4 e 0.31150 0.09070 0.13340 1.00000
Cu1 Cu 4 e 0.03201 0.11772 0.28577 1.00000
H1 H 4 e 0.67400 0.12800 0.52800 1.00000
O1 O 4 e 0.88070 0.14780 0.53180 1.00000
```

Cu(OH)Cl: ABCD_mP16_14_e_e_e_e - POSCAR

```
ABCD_mP16_14_e_e_e_e & a,b/a,c/a,beta,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,
    ↪ z4 --params=6.2953,1.05871046654,0.882880879386,118.138,0.3115,
    ↪ 0.0907,0.1334,0.03201,0.11772,0.28577,0.674,0.128,0.528,0.8807,
    ↪ 0.1478,0.5318 & P2_{1}/c C_{2h}^{5} #14 (e^4) & mP16 & None &
    ↪ ClCuHO & ClCuHO & Y. Cudennec et al., J. Solid State Chem. 151,
    ↪ 308-312 (2000)
   1.00000000000000
   6.29530000000000    0.00000000000000    0.00000000000000
   0.00000000000000    6.66490000000000    0.00000000000000
  -2.62113516045273    0.00000000000000    4.90112379670606
     Cl    Cu    H     O
     4     4     4     4
Direct
   0.31150000000000    0.09070000000000    0.13340000000000    Cl   (4e)
  -0.31150000000000    0.59070000000000    0.36660000000000    Cl   (4e)
  -0.31150000000000   -0.09070000000000   -0.13340000000000    Cl   (4e)
   0.31150000000000    0.40930000000000    0.63340000000000    Cl   (4e)
   0.03201000000000    0.11772000000000    0.28577000000000    Cu   (4e)
  -0.03201000000000    0.61772000000000    0.21423000000000    Cu   (4e)
  -0.03201000000000   -0.11772000000000   -0.28577000000000    Cu   (4e)
   0.03201000000000    0.38228000000000    0.78577000000000    Cu   (4e)
   0.67400000000000    0.12800000000000    0.52800000000000    H    (4e)
  -0.67400000000000    0.62800000000000   -0.02800000000000    H    (4e)
  -0.67400000000000   -0.12800000000000   -0.52800000000000    H    (4e)
   0.67400000000000    0.37200000000000    1.02800000000000    H    (4e)
   0.88070000000000    0.14780000000000    0.53180000000000    O    (4e)
```

```
  -0.88070000000000   0.64780000000000  -0.03180000000000   O   (4e)
  -0.88070000000000  -0.14780000000000  -0.53180000000000   O   (4e)
   0.88070000000000   0.35220000000000   1.03180000000000   O   (4e)
```

Arsenopyrite (FeAsS, $E0_7$): ABC_mP12_14_e_e_e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Arsenopyrite'
_chemical_formula_sum 'As Fe S'

loop_
_publ_author_name
 'L. Bindi'
 'Y. Mo\"{e}lo'
 'P. L\\'eone'
 'M. Suchaud'
_journal_name_full_name
;
 Canadian Mineralogist
;
_journal_volume 50
_journal_year 2012
_journal_page_first 471
_journal_page_last 479
_publ_Section_title
;
 Stoichiometric Arsenopyrite, FeAsS, from La Roche-Balue Quarry,
      ↪ Loire-Atlantique, France: Crystal Structure And M\"{o}ssbauer
      ↪ Study
;

_aflow_title 'Arsenopyrite (FeAsS, $E0_{7}$) Structure'
_aflow_proto 'ABC_mP12_14_e_e_e'
_aflow_params 'a,b/a,c/a,\beta,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3
      ↪ },y_{3},z_{3}'
_aflow_params_values '5.7612,0.986617371381,1.00107616469,111.721,
      ↪ 0.14746,0.13055,0.86937,0.28353,-0.00643,0.29429,0.6551,0.1311,
      ↪ 0.3211'
_aflow_Strukturbericht '$E0_{7}$'
_aflow_Pearson 'mP12'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

_cell_length_a     5.76120
_cell_length_b     5.68410
_cell_length_c     5.76740
_cell_angle_alpha  90.00000
_cell_angle_beta   111.72100
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z+1/2
3 -x,-y,-z
4 x,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
As1 As 4 e 0.14746 0.13055 0.86937 1.00000
Fe1 Fe 4 e 0.28353 -0.00643 0.29429 1.00000
S1 S 4 e 0.65510 0.13110 0.32110 1.00000
```

Arsenopyrite (FeAsS, $E0_7$): ABC_mP12_14_e_e_e - POSCAR

```
ABC_mP12_14_e_e_e & a,b/a,c/a,beta,x1,y1,z1,x2,y2,z2,x3,y3,z3 --params=
      ↪ 5.7612,0.986617371381,1.00107616469,111.721,0.14746,0.13055,
      ↪ 0.86937,0.28353,-0.00643,0.29429,0.6551,0.1311,0.3211 & P2_{1}/
      ↪ c C_{2h}^{5} #14 (e^3) & mP12 & $E0_{7}$ & AsFeS & Arsenopyrite
      ↪ & L. Bindi et al., Can. Mineral. 50, 471-479 (2012)
   1.00000000000000
   5.76120000000000   0.00000000000000   0.00000000000000
   0.00000000000000   5.68410000000000   0.00000000000000
  -2.13444136312758   0.00000000000000   5.35789723934400
     As     Fe     S
      4      4     4
Direct
   0.14746000000000   0.13055000000000   0.86937000000000   As   (4e)
  -0.14746000000000   0.63055000000000  -0.36937000000000   As   (4e)
  -0.14746000000000  -0.13055000000000  -0.86937000000000   As   (4e)
   0.14746000000000   0.36945000000000   1.36937000000000   As   (4e)
   0.28353000000000  -0.00643000000000   0.29429000000000   Fe   (4e)
  -0.28353000000000   0.49357000000000   0.20571000000000   Fe   (4e)
  -0.28353000000000   0.00643000000000  -0.29429000000000   Fe   (4e)
   0.28353000000000   0.50643000000000   0.79429000000000   Fe   (4e)
   0.65510000000000   0.13110000000000   0.32110000000000    S   (4e)
  -0.65510000000000   0.63110000000000   0.17890000000000    S   (4e)
  -0.65510000000000  -0.13110000000000  -0.32110000000000    S   (4e)
   0.65510000000000   0.36890000000000   0.82110000000000    S   (4e)
```

α-ICl: AB_mP16_14_2e_2e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'ClI'
_chemical_formula_sum 'Cl I'

loop_
_publ_author_name
 'K. H. Boswijk'
 'J. {van der Heide}'
 'A. Vos'
 'E. H. Wiebenga'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 9
_journal_year 1956
_journal_page_first 274
_journal_page_last 277
_publ_Section_title
;
 The crystal structure of $\alpha$-ICl
;

_aflow_title '$\alpha$-ICl Structure'
_aflow_proto 'AB_mP16_14_2e_2e'
_aflow_params 'a,b/a,c/a,\beta,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3
      ↪ },y_{3},z_{3},x_{4},y_{4},z_{4}'
_aflow_params_values '12.6,0.347619047619,0.944444444444,119.5,0.084,
      ↪ 0.152,0.706,0.462,0.858,0.62,0.179,0.366,0.588,0.297,0.632,
      ↪ 0.436'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mP16'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

_cell_length_a     12.60000
_cell_length_b     4.38000
_cell_length_c     11.90000
_cell_angle_alpha  90.00000
_cell_angle_beta   119.50000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z+1/2
3 -x,-y,-z
4 x,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cl1 Cl 4 e 0.08400 0.15200 0.70600 1.00000
Cl2 Cl 4 e 0.46200 0.85800 0.62000 1.00000
I1 I 4 e 0.17900 0.36600 0.58800 1.00000
I2 I 4 e 0.29700 0.63200 0.43600 1.00000
```

α-ICl: AB_mP16_14_2e_2e - POSCAR

```
AB_mP16_14_2e_2e & a,b/a,c/a,beta,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,z4 --
      ↪ params=12.6,0.347619047619,0.944444444444,119.5,0.084,0.152,
      ↪ 0.706,0.462,0.858,0.62,0.179,0.366,0.588,0.297,0.632,0.436 &
      ↪ P2_{1}/c C_{2h}^{5} #14 (e^4) & mP16 & None & ClI & ClI & K. H.
      ↪ Boswijk et al., Acta Cryst. 9, 274-277 (1956)
   1.00000000000000
  12.60000000000000   0.00000000000000   0.00000000000000
   0.00000000000000   4.38000000000000   0.00000000000000
  -5.85984036523126   0.00000000000000  10.35723278168480
     Cl     I
      8     8
Direct
   0.08400000000000   0.15200000000000   0.70600000000000   Cl   (4e)
  -0.08400000000000   0.65200000000000  -0.20600000000000   Cl   (4e)
  -0.08400000000000  -0.15200000000000  -0.70600000000000   Cl   (4e)
   0.08400000000000   0.34800000000000   1.20600000000000   Cl   (4e)
   0.46200000000000   0.85800000000000   0.62000000000000   Cl   (4e)
  -0.46200000000000   1.35800000000000  -0.12000000000000   Cl   (4e)
  -0.46200000000000  -0.85800000000000  -0.62000000000000   Cl   (4e)
   0.46200000000000  -0.35800000000000   1.12000000000000   Cl   (4e)
   0.17900000000000   0.36600000000000   0.58800000000000    I   (4e)
  -0.17900000000000   0.86600000000000  -0.08800000000000    I   (4e)
  -0.17900000000000  -0.36600000000000  -0.58800000000000    I   (4e)
   0.17900000000000   0.13400000000000   1.08800000000000    I   (4e)
   0.29700000000000   0.63200000000000   0.43600000000000    I   (4e)
  -0.29700000000000   1.13200000000000   0.06400000000000    I   (4e)
  -0.29700000000000  -0.63200000000000  -0.43600000000000    I   (4e)
   0.29700000000000  -0.13200000000000   0.93600000000000    I   (4e)
```

LiAs: AB_mP16_14_2e_2e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM
```

```
_chemical_name_mineral 'AsLi'
_chemical_formula_sum 'As Li'

loop_
_publ_author_name
 'D. T. Cromer'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 12
_journal_year 1959
_journal_page_first 36
_journal_page_last 41
_publ_Section_title
;
 The Crystal Structure of LiAs
;

# Found in Pearson's Handbook of Crystallographic Data for Intermetallic
      ↪  Phases, 1991

_aflow_title 'LiAs Structure'
_aflow_proto 'AB_mP16_14_2e_2e'
_aflow_params 'a,b/a,c/a,\beta,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3
      ↪ },y_{3},z_{3},x_{4},y_{4},z_{4}'
_aflow_params_values '5.79,0.905008635579,1.84801381693,117.4,0.3042,
      ↪ 0.9143,0.2992,0.2891,0.1626,0.1011,0.235,0.402,0.329,0.232,
      ↪ 0.669,0.045'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mP16'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

_cell_length_a     5.79000
_cell_length_b     5.24000
_cell_length_c     10.70000
_cell_angle_alpha  90.00000
_cell_angle_beta   117.40000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z+1/2
3 -x,-y,-z
4 x,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
As1 As 4 e 0.30420 0.91430 0.29920 1.00000
As2 As 4 e 0.28910 0.16260 0.10110 1.00000
Li1 Li 4 e 0.23500 0.40200 0.32900 1.00000
Li2 Li 4 e 0.23200 0.66900 0.04500 1.00000
```

LiAs: AB_mP16_14_2e_2e - POSCAR

```
AB_mP16_14_2e_2e & a,b/a,c/a,beta,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,z4 --
      ↪ params=5.79,0.905008635579,1.84801381693,117.4,0.3042,0.9143,
      ↪ 0.2992,0.2891,0.1626,0.1011,0.235,0.402,0.329,0.232,0.669,0.045
      ↪ & P2_{1}/c C_{2h}^{5} #14 (e^4) & mP16 & None & AsLi & AsLi &
      ↪ D. T. Cromer, Acta Cryst. 12, 36-41 (1959)
   1.00000000000000
   5.79000000000000    0.00000000000000    0.00000000000000
   0.00000000000000    5.24000000000000    0.00000000000000
  -4.92413769718721    0.00000000000000    9.49962462095949
     As      Li
      8       8
Direct
   0.30420000000000    0.91430000000000    0.29920000000000    As    (4e)
  -0.30420000000000    1.41430000000000    0.20080000000000    As    (4e)
  -0.30420000000000   -0.91430000000000   -0.29920000000000    As    (4e)
   0.30420000000000   -0.41430000000000    0.79920000000000    As    (4e)
   0.28910000000000    0.16260000000000    0.10110000000000    As    (4e)
  -0.28910000000000    0.66260000000000    0.39890000000000    As    (4e)
  -0.28910000000000   -0.16260000000000   -0.10110000000000    As    (4e)
   0.28910000000000    0.33740000000000    0.60110000000000    As    (4e)
   0.23500000000000    0.40200000000000    0.32900000000000    Li    (4e)
  -0.23500000000000    0.90200000000000    0.17100000000000    Li    (4e)
  -0.23500000000000   -0.40200000000000   -0.32900000000000    Li    (4e)
   0.23500000000000    0.09800000000000    0.82900000000000    Li    (4e)
   0.23200000000000    0.66900000000000    0.04500000000000    Li    (4e)
  -0.23200000000000    1.16900000000000    0.45500000000000    Li    (4e)
  -0.23200000000000   -0.66900000000000   -0.04500000000000    Li    (4e)
   0.23200000000000   -0.16900000000000    0.54500000000000    Li    (4e)
```

ϵ-1,2,3,4,5,6-Hexachlorocyclohexane (C_6Cl_6): AB_mP24_14_3e_3e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral '$\epsilon$-1,2,3,4,5,6-hexachlorocyclohexane'
_chemical_formula_sum 'C Cl'

loop_
_publ_author_name
 'N. Norman'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 4
_journal_year 1950
_journal_page_first 251
_journal_page_last 259
_publ_Section_title
;
 The Crystal Structure of the Epsilon Isomer of 1,2,3,4,5,6-Hexachloro\
      ↪ em{cyclo}hexane
;

_aflow_title '$\epsilon$-1,2,3,4,5,6-Hexachlorocyclohexane (C$_{6}$Cl$_{
      ↪ 6}$) Structure'
_aflow_proto 'AB_mP24_14_3e_3e'
_aflow_params 'a,b/a,c/a,\beta,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3
      ↪ },y_{3},z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},
      ↪ z_{6}'
_aflow_params_values '7.02,1.67948717949,0.968660968661,112.0,0.206,
      ↪ 0.013,0.005,0.089,0.115,0.095,-0.035,0.041,0.177,0.369,-0.071,
      ↪ 0.212,0.27,0.188,0.312,0.81,0.13,0.298'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mP24'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

_cell_length_a     7.02000
_cell_length_b     11.79000
_cell_length_c     6.80000
_cell_angle_alpha  90.00000
_cell_angle_beta   112.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z+1/2
3 -x,-y,-z
4 x,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
C1 C 4 e 0.20600 0.01300 0.00500 1.00000
C2 C 4 e 0.08900 0.11500 0.09500 1.00000
C3 C 4 e -0.03500 0.04100 0.17700 1.00000
Cl1 Cl 4 e 0.36900 -0.07100 0.21200 1.00000
Cl2 Cl 4 e 0.27000 0.18800 0.31200 1.00000
Cl3 Cl 4 e 0.81000 0.13000 0.29800 1.00000
```

ϵ-1,2,3,4,5,6-Hexachlorocyclohexane (C_6Cl_6): AB_mP24_14_3e_3e - POSCAR

```
AB_mP24_14_3e_3e & a,b/a,c/a,beta,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,z4,x5
      ↪ ,y5,z5,x6,y6,z6 --params=7.02,1.67948717949,0.968660968661,
      ↪ 112.0,0.206,0.013,0.005,0.089,0.115,0.095,-0.035,0.041,0.177,
      ↪ 0.369,-0.071,0.212,0.27,0.188,0.312,0.81,0.13,0.298 & P2_{1}/c
      ↪ C_{2h}^{5} #14 (e^6) & mP24 & None & C6Cl6 & $\epsilon$-1,2,3,4
      ↪ ,5,6-hexachlorocyclohexane & N. Norman, Acta Chem. Scand. 4,
      ↪ 251-259 (1950)
   1.00000000000000
   7.02000000000000    0.00000000000000    0.00000000000000
   0.00000000000000   11.79000000000000    0.00000000000000
  -2.54732483522820    0.00000000000000    6.30485021105415
      C      Cl
     12      12
Direct
   0.20600000000000    0.01300000000000    0.00500000000000    C     (4e)
  -0.20600000000000    0.51300000000000    0.49500000000000    C     (4e)
  -0.20600000000000   -0.01300000000000   -0.00500000000000    C     (4e)
   0.20600000000000    0.48700000000000    0.50500000000000    C     (4e)
   0.08900000000000    0.11500000000000    0.09500000000000    C     (4e)
  -0.08900000000000    0.61500000000000    0.40500000000000    C     (4e)
  -0.08900000000000   -0.11500000000000   -0.09500000000000    C     (4e)
   0.08900000000000    0.38500000000000    0.59500000000000    C     (4e)
  -0.03500000000000    0.04100000000000    0.17700000000000    C     (4e)
   0.03500000000000    0.54100000000000    0.32300000000000    C     (4e)
   0.03500000000000   -0.04100000000000   -0.17700000000000    C     (4e)
  -0.03500000000000    0.45900000000000    0.67700000000000    C     (4e)
   0.36900000000000   -0.07100000000000    0.21200000000000    Cl    (4e)
  -0.36900000000000    0.42900000000000    0.28800000000000    Cl    (4e)
  -0.36900000000000    0.07100000000000   -0.21200000000000    Cl    (4e)
   0.36900000000000    0.57100000000000    0.71200000000000    Cl    (4e)
   0.27000000000000    0.18800000000000    0.31200000000000    Cl    (4e)
  -0.27000000000000    0.68800000000000    0.18800000000000    Cl    (4e)
  -0.27000000000000   -0.18800000000000   -0.31200000000000    Cl    (4e)
   0.27000000000000    0.31200000000000    0.81200000000000    Cl    (4e)
   0.81000000000000    0.13000000000000    0.29800000000000    Cl    (4e)
  -0.81000000000000    0.63000000000000    0.20200000000000    Cl    (4e)
  -0.81000000000000   -0.13000000000000   -0.29800000000000    Cl    (4e)
   0.81000000000000    0.37000000000000    0.79800000000000    Cl    (4e)
```

Pararealgar (AsS): AB_mP32_14_4e_4e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Pararealgar'
_chemical_formula_sum 'As S'

loop_
_publ_author_name
 'P. Bonazzi'
 'S. Menchetti'
 'G. Pratesi'
_journal_name_full_name
;
 American Mineralogist
;
_journal_volume 80
_journal_year 1995
_journal_page_first 400
_journal_page_last 403
_publ_Section_title
;
 The crystal structure of pararealgar, As$_{4}$S$_{4}$
;

_aflow_title 'Pararealgar (AsS) Structure'
_aflow_proto 'AB_mP32_14_4e_4e'
_aflow_params 'a,b/a,c/a,\beta,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3
      ↪ },y_{3},z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},
      ↪ z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},z_{8}'
_aflow_params_values '9.909,0.97436673731,0.858007871632,97.29,0.3187,
      ↪ 0.6355,0.0432,0.0819,0.5427,0.3252,0.3698,0.3607,0.3431,0.1455,
      ↪ 0.3439,0.1643,0.1645,0.7187,0.1923,0.2537,0.4782,0.5099,0.4703,
      ↪ 0.5276,0.2192,0.1964,0.4483,-0.0492'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mP32'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

_cell_length_a    9.90900
_cell_length_b    9.65500
_cell_length_c    8.50200
_cell_angle_alpha 90.00000
_cell_angle_beta  97.29000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z+1/2
3 -x,-y,-z
4 x,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
As1 As 4 e 0.31870 0.63550 0.04320 1.00000
As2 As 4 e 0.08190 0.54270 0.32520 1.00000
As3 As 4 e 0.36980 0.36070 0.34310 1.00000
As4 As 4 e 0.14550 0.34390 0.16430 1.00000
S1 S 4 e 0.16450 0.71870 0.19230 1.00000
S2 S 4 e 0.25370 0.47820 0.50990 1.00000
S3 S 4 e 0.47030 0.52760 0.21920 1.00000
S4 S 4 e 0.19640 0.44830 -0.04920 1.00000
```

Pararealgar (AsS): AB_mP32_14_4e_4e - POSCAR

```
AB_mP32_14_4e_4e & a,b/a,c/a,beta,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,z4,x5
      ↪ ,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8 --params=9.909,0.97436673731,
      ↪ 0.858007871632,97.29,0.3187,0.6355,0.0432,0.0819,0.5427,0.3252,
      ↪ 0.3698,0.3607,0.3431,0.1455,0.3439,0.1643,0.1645,0.7187,0.1923,
      ↪ 0.2537,0.4782,0.5099,0.4703,0.5276,0.2192,0.1964,0.4483,-0.0492
      ↪  & P2_{1}/c C_{2h}^{5} #14 (e^8) & mP32 & None & AsS &
      ↪ Pararealgar & P. Bonazzi and S. Menchetti and G. Pratesi, Am.
      ↪ Mineral. 80, 400-403 (1995)
   1.00000000000000
   9.90900000000000    0.00000000000000    0.00000000000000
   0.00000000000000    9.65500000000000    0.00000000000000
  -1.07883143461801    0.00000000000000    8.43327497095168
     As      S
     16     16
Direct
   0.31870000000000    0.63550000000000    0.04320000000000    As    (4e)
  -0.31870000000000    1.13550000000000    0.45680000000000    As    (4e)
  -0.31870000000000   -0.63550000000000   -0.04320000000000    As    (4e)
   0.31870000000000   -0.13550000000000    0.54320000000000    As    (4e)
   0.08190000000000    0.54270000000000    0.32520000000000    As    (4e)
  -0.08190000000000    1.04270000000000    0.17480000000000    As    (4e)
  -0.08190000000000   -0.54270000000000   -0.32520000000000    As    (4e)
   0.08190000000000   -0.04270000000000    0.82520000000000    As    (4e)
   0.36980000000000    0.36070000000000    0.34310000000000    As    (4e)
  -0.36980000000000    0.86070000000000    0.15690000000000    As    (4e)
  -0.36980000000000   -0.36070000000000   -0.34310000000000    As    (4e)
   0.36980000000000    0.13930000000000    0.84310000000000    As    (4e)
   0.14550000000000    0.34390000000000    0.16430000000000    As    (4e)
  -0.14550000000000    0.84390000000000    0.33570000000000    As    (4e)
  -0.14550000000000   -0.34390000000000   -0.16430000000000    As    (4e)
   0.14550000000000    0.15610000000000    0.66430000000000    As    (4e)
   0.16450000000000    0.71870000000000    0.19230000000000     S    (4e)
  -0.16450000000000    1.21870000000000    0.30770000000000     S    (4e)
  -0.16450000000000   -0.71870000000000   -0.19230000000000     S    (4e)
   0.16450000000000   -0.21870000000000    0.69230000000000     S    (4e)
   0.25370000000000    0.47820000000000    0.50990000000000     S    (4e)
  -0.25370000000000    0.97820000000000   -0.00990000000000     S    (4e)
  -0.25370000000000   -0.47820000000000   -0.50990000000000     S    (4e)
   0.25370000000000    0.02180000000000    1.00990000000000     S    (4e)
   0.47030000000000    0.52760000000000    0.21920000000000     S    (4e)
  -0.47030000000000    1.02760000000000    0.28080000000000     S    (4e)
  -0.47030000000000   -0.52760000000000   -0.21920000000000     S    (4e)
   0.47030000000000   -0.02760000000000    0.71920000000000     S    (4e)
   0.19640000000000    0.44830000000000   -0.04920000000000     S    (4e)
  -0.19640000000000    0.94830000000000    0.54920000000000     S    (4e)
  -0.19640000000000   -0.44830000000000    0.04920000000000     S    (4e)
   0.19640000000000    0.05170000000000    0.45080000000000     S    (4e)
```

Realgar (AsS, B_l): AB_mP32_14_4e_4e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Realgar'
_chemical_formula_sum 'As S'

loop_
_publ_author_name
 'T. Ito'
 'N. Morimoto'
 'R. Sadanaga'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 5
_journal_year 1952
_journal_page_first 755
_journal_page_last 782
_publ_Section_title
;
 The Crystal Structure of Realgar
;

_aflow_title 'Realgar (AsS, $B_{1}$) Structure'
_aflow_proto 'AB_mP32_14_4e_4e'
_aflow_params 'a,b/a,c/a,\beta,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3
      ↪ },y_{3},z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},
      ↪ z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},z_{8}'
_aflow_params_values '6.56,2.05792682927,1.47941615854,113.75283,0.359,
      ↪ 0.524,0.118,0.567,0.36,0.425,0.137,0.373,0.318,0.328,0.339,
      ↪ 0.038,0.641,0.508,0.346,0.093,0.524,0.213,0.608,0.275,0.245,
      ↪ 0.067,0.285,0.115'
_aflow_Strukturbericht '$B_{1}$'
_aflow_Pearson 'mP32'

_symmetry_space_group_name_H-M "P 1 21/c 1"
_symmetry_Int_Tables_number 14

_cell_length_a    6.56000
_cell_length_b    13.50000
_cell_length_c    9.70497
_cell_angle_alpha 90.00000
_cell_angle_beta  113.75283
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y+1/2,-z+1/2
3 -x,-y,-z
4 x,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
As1 As 4 e 0.35900 0.52400 0.11800 1.00000
As2 As 4 e 0.56700 0.36000 0.42500 1.00000
As3 As 4 e 0.13700 0.37300 0.31800 1.00000
As4 As 4 e 0.32800 0.33900 0.03800 1.00000
S1 S 4 e 0.64100 0.50800 0.34600 1.00000
S2 S 4 e 0.09300 0.52400 0.21300 1.00000
S3 S 4 e 0.60800 0.27500 0.24500 1.00000
S4 S 4 e 0.06700 0.28500 0.11500 1.00000
```

Realgar (AsS, B_l): AB_mP32_14_4e_4e - POSCAR

```
AB_mP32_14_4e_4e & a,b/a,c/a,beta,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,z4,x5
      ↪ ,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8 --params=6.56,2.05792682927,
      ↪ 1.47941615854,113.75283,0.359,0.524,0.118,0.567,0.36,0.425,
      ↪ 0.137,0.373,0.318,0.328,0.339,0.038,0.641,0.508,0.346,0.093,
      ↪ 0.524,0.213,0.608,0.275,0.245,0.067,0.285,0.115 & P2_{1}/c C_{
      ↪ 2h}^{5} #14 (e^8) & mP32 & $B_{1}$ & AsS & Realgar & T. Ito and
      ↪  N. Morimoto and R. Sadanaga, Acta Cryst. 5, 755-782 (1952)
   1.00000000000000
   6.56000000000000    0.00000000000000    0.00000000000000
   0.00000000000000   13.50000000000000    0.00000000000000
```

```
  -3.90908329779061    0.00000000000000    8.88287737570628
     As      S
     16      16
Direct
   0.35900000000000    0.52400000000000    0.11800000000000    As    (4e)
  -0.35900000000000    1.02400000000000    0.38200000000000    As    (4e)
  -0.35900000000000   -0.52400000000000   -0.11800000000000    As    (4e)
   0.35900000000000   -0.02400000000000    0.61800000000000    As    (4e)
   0.56700000000000    0.36000000000000    0.42500000000000    As    (4e)
  -0.56700000000000    0.86000000000000    0.07500000000000    As    (4e)
  -0.56700000000000   -0.36000000000000   -0.42500000000000    As    (4e)
   0.56700000000000    0.14000000000000    0.92500000000000    As    (4e)
   0.13700000000000    0.37300000000000    0.31800000000000    As    (4e)
  -0.13700000000000    0.87300000000000    0.18200000000000    As    (4e)
  -0.13700000000000   -0.37300000000000   -0.31800000000000    As    (4e)
   0.13700000000000    0.12700000000000    0.81800000000000    As    (4e)
   0.32800000000000    0.33900000000000    0.03800000000000    As    (4e)
  -0.32800000000000    0.83900000000000    0.46200000000000    As    (4e)
  -0.32800000000000   -0.33900000000000   -0.03800000000000    As    (4e)
   0.32800000000000    0.16100000000000    0.53800000000000    As    (4e)
   0.64100000000000    0.50800000000000    0.34600000000000    S     (4e)
  -0.64100000000000    1.00800000000000    0.15400000000000    S     (4e)
  -0.64100000000000   -0.50800000000000   -0.34600000000000    S     (4e)
   0.64100000000000   -0.00800000000000    0.84600000000000    S     (4e)
   0.09300000000000    0.52400000000000    0.21300000000000    S     (4e)
  -0.09300000000000    1.02400000000000    0.28700000000000    S     (4e)
  -0.09300000000000   -0.52400000000000   -0.21300000000000    S     (4e)
   0.09300000000000   -0.02400000000000    0.71300000000000    S     (4e)
   0.60800000000000    0.27500000000000    0.24500000000000    S     (4e)
  -0.60800000000000    0.77500000000000    0.25500000000000    S     (4e)
  -0.60800000000000   -0.27500000000000   -0.24500000000000    S     (4e)
   0.60800000000000    0.22500000000000    0.74500000000000    S     (4e)
   0.06700000000000    0.28500000000000    0.11500000000000    S     (4e)
  -0.06700000000000    0.78500000000000    0.38500000000000    S     (4e)
  -0.06700000000000   -0.28500000000000   -0.11500000000000    S     (4e)
   0.06700000000000    0.21500000000000    0.61500000000000    S     (4e)
```

Ag_2PbO_2: A2B2C_mC20_15_ad_f_e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Ag2O2Pb'
_chemical_formula_sum 'Ag2 O2 Pb'

loop_
_publ_author_name
 'A. Bystr\"{o}m'
 'L. Evers'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 4
_journal_year 1950
_journal_page_first 613
_journal_page_last 627
_publ_Section_title
;
 The Crystal Structures of Ag$_{2}$PbO$_{2}$ and Ag5Pb$_{2}$O$_{6}$
;

_aflow_title 'Ag$_{2}$PbO$_{2}$ Structure'
_aflow_proto 'A2B2C_mC20_15_ad_f_e'
_aflow_params 'a,b/a,c/a,\beta,y_{3},x_{4},y_{4},z_{4}'
_aflow_params_values '8.65,1.00751445087,0.703121387283,130.86424,0.625,
    ↪ 0.804,0.445,0.865'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mC20'

_symmetry_space_group_name_H-M "C 1 2/c 1"
_symmetry_Int_Tables_number 15

_cell_length_a     8.65000
_cell_length_b     8.71500
_cell_length_c     6.08200
_cell_angle_alpha  90.00000
_cell_angle_beta   130.86424
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z+1/2
3 -x,-y,-z
4 x,-y,z+1/2
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z+1/2
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ag1 Ag 4 a 0.00000 0.00000 0.00000 1.00000
Ag2 Ag 4 d 0.25000 0.25000 0.50000 1.00000
Pb1 Pb 4 e 0.00000 0.62500 0.25000 1.00000
O1 O 8 f 0.80400 0.44500 0.86500 1.00000
```

Ag_2PbO_2: A2B2C_mC20_15_ad_f_e - POSCAR

```
A2B2C_mC20_15_ad_f_e & a,b/a,c/a,beta,y3,x4,y4,z4 --params=8.65,
    ↪ 1.00751445087,0.703121387283,130.86424,0.625,0.804,0.445,0.865
    ↪ & C2/c C_{2h}^{6} #15 (adef) & mC20 & None & Ag2O2Pb & Ag2O2Pb
    ↪ & A. Bystr\"{o}m and L. Evers, Acta Chem. Scand. 4, 613-627 (
    ↪ 1950)
   1.00000000000000
   4.32500000000000   -4.35750000000000    0.00000000000000
   4.32500000000000    4.35750000000000    0.00000000000000
  -3.97926366629002    0.00000000000000    4.59958527197226
     Ag      O      Pb
      4      4       2
Direct
   0.00000000000000    0.00000000000000    0.00000000000000    Ag    (4a)
   0.00000000000000    0.00000000000000    0.50000000000000    Ag    (4a)
   0.00000000000000    0.50000000000000    0.50000000000000    Ag    (4d)
   0.50000000000000    0.00000000000000    0.00000000000000    Ag    (4d)
   0.35900000000000    1.24900000000000    0.86500000000000    O     (8f)
  -1.24900000000000   -0.35900000000000   -0.36500000000000    O     (8f)
  -0.35900000000000   -1.24900000000000   -0.86500000000000    O     (8f)
   1.24900000000000    0.35900000000000    1.36500000000000    O     (8f)
  -0.62500000000000    0.62500000000000    0.25000000000000    Pb    (4e)
   0.62500000000000   -0.62500000000000    0.75000000000000    Pb    (4e)
```

Catapleiite ($Na_2ZrSi_3O_9{\cdot}2H_2O$): A2B3C9D3E_mC144_15_2f_bcdef_9f_3f_ae - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Catapleiite'
_chemical_formula_sum '(H2O)2 Na3 O9 Si3 Zr'

loop_
_publ_author_name
 'G. D. Ilyushin'
 'A. A. Voronkov'
 'V. V. Ilyukhin'
 'N. N. Nevskii'
 'N. V. Belov'
_journal_name_full_name
;
 Doklady Akademii Nauk SSSR
;
_journal_volume 260
_journal_year 1981
_journal_page_first 623
_journal_page_last 627
_publ_Section_title
;
 Crystal structure of natural monoclinic catapleiite, Na$_{2}$ZrSi$_{3}
    ↪ $O$_{9}\cdot$ 2H$_{2}$O
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'Catapleiite (Na$_{2}$ZrSi$_{3}$O$_{9}\cdot2H$_{2}$O)
    ↪ Structure'
_aflow_proto 'A2B3C9D3E_mC144_15_2f_bcdef_9f_3f_ae'
_aflow_params 'a,b/a,c/a,\beta,y_{5},y_{6},x_{7},y_{7},z_{7},x_{8},y_{8
    ↪ },z_{8},x_{9},y_{9},z_{9},x_{10},y_{10},z_{10},x_{11},y_{11},z_
    ↪ {11},x_{12},y_{12},z_{12},x_{13},y_{13},z_{13},x_{14},y_{14},z_
    ↪ {14},x_{15},y_{15},z_{15},x_{16},y_{16},z_{16},x_{17},y_{17},z_
    ↪ {17},x_{18},y_{18},z_{18},x_{19},y_{19},z_{19},x_{20},y_{20},z_
    ↪ {20},x_{21},y_{21},z_{21}'
_aflow_params_values '23.8927,0.311057352246,0.843270120162,147.42281,
    ↪ 0.4941,-0.0074,0.8315,0.4922,0.0362,0.3371,0.0092,0.8816,0.2531
    ↪ ,0.2432,0.255,0.8661,0.0088,-0.0767,0.3666,0.49,0.0586,0.0493,
    ↪ 0.4955,0.1747,0.0711,0.1973,0.1285,0.0641,0.1942,0.2558,0.2265,
    ↪ 0.3221,0.3544,0.562,0.298,0.1203,0.569,0.2938,0.2625,0.7239,
    ↪ 0.1738,0.3455,0.5976,0.2004,0.2221,0.0995,0.2923,0.2252,0.1972,
    ↪ 0.0015,0.5723'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mC144'

_symmetry_space_group_name_H-M "C 1 2/c 1"
_symmetry_Int_Tables_number 15

_cell_length_a     23.89270
_cell_length_b     7.43200
_cell_length_c     20.14800
_cell_angle_alpha  90.00000
_cell_angle_beta   147.42281
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z+1/2
3 -x,-y,-z
4 x,-y,z+1/2
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z+1/2
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
```

```
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Zr1  Zr  4  a  0.00000  0.00000  0.00000  1.00000
Na1  Na  4  b  0.00000  0.50000  0.00000  1.00000
Na2  Na  4  c  0.25000  0.25000  0.00000  1.00000
Na3  Na  4  d  0.25000  0.25000  0.50000  1.00000
Na4  Na  4  e  0.00000  0.49410  0.25000  1.00000
Zr2  Zr  4  e  0.00000 -0.00740  0.25000  1.00000
H2O1 H2O 8  f  0.83150  0.49220  0.03620  1.00000
H2O2 H2O 8  f  0.33710  0.00920  0.88160  1.00000
Na5  Na  8  f  0.25310  0.24320  0.25500  1.00000
O1  O  8  f  0.86610  0.00880 -0.07670  1.00000
O2  O  8  f  0.36660  0.49000  0.05860  1.00000
O3  O  8  f  0.04930  0.49550  0.17470  1.00000
O4  O  8  f  0.07110  0.19730  0.12850  1.00000
O5  O  8  f  0.06410  0.19420  0.25580  1.00000
O6  O  8  f  0.22650  0.32210  0.35440  1.00000
O7  O  8  f  0.56200  0.29800  0.12030  1.00000
O8  O  8  f  0.56900  0.29380  0.26250  1.00000
O9  O  8  f  0.72390  0.17380  0.34550  1.00000
Si1  Si  8  f  0.59760  0.20040  0.22210  1.00000
Si2  Si  8  f  0.09950  0.29230  0.22520  1.00000
Si3  Si  8  f  0.19720  0.00150  0.57230  1.00000
```

Catapleiite ($Na_2ZrSi_3O_9{\cdot}2H_2O$): A2B3C9D3E_mC144_15_2f_bcdef_9f_3f_ae - POSCAR

```
A2B3C9D3E_mC144_15_2f_bcdef_9f_3f_ae & a,b/a,c/a,beta,y5,y6,x7,y7,z7,x8,
    ↪ y8,z8,x9,y9,z9,x10,y10,z10,x11,y11,z11,x12,y12,z12,x13,y13,z13,
    ↪ x14,y14,z14,x15,y15,z15,x16,y16,z16,x17,y17,z17,x18,y18,z18,x19
    ↪ ,y19,z19,x20,y20,z20,x21,y21,z21 --params=23.8927,
    ↪ 0.311057352246,0.843270120162,147.42281,0.4941,-0.0074,0.8315,
    ↪ 0.4922,0.0362,0.3371,0.0092,0.8816,0.2531,0.2432,0.255,0.8661,
    ↪ 0.0088,-0.0767,0.3666,0.49,0.0586,0.0493,0.4955,0.1747,0.0711,
    ↪ 0.1973,0.1285,0.0641,0.1942,0.2558,0.2265,0.3221,0.3544,0.562,
    ↪ 0.298,0.1203,0.569,0.2938,0.2625,0.7239,0.1738,0.3455,0.5976,
    ↪ 0.2004,0.2221,0.0995,0.2923,0.2252,0.1972,0.0015,0.5723 & C2/c
    ↪ C_{2h}^{6} #15 (abcde^2f^15) & mC144 & None & (H2O)2Na2O9Si3Zr
    ↪ & Catapleiite & G. D. Ilyushin et al., Doklady Akademii Nauk
    ↪ SSSR 260, 623-627 (1981)
   1.00000000000000
  11.94635000000000  -3.71600000000000   0.00000000000000
  11.94635000000000   3.71600000000000   0.00000000000000
 -16.97805109064560   0.00000000000000  10.84839551101580
    H2O     Na      O     Si     Zr
      8     12     36     12      4
Direct
   0.33930000000000   1.32370000000000   0.03620000000000  H2O   (8f)
  -1.32370000000000  -0.33930000000000   0.46380000000000  H2O   (8f)
  -0.33930000000000  -1.32370000000000  -0.03620000000000  H2O   (8f)
   1.32370000000000   0.33930000000000   0.53620000000000  H2O   (8f)
   0.32790000000000   0.34630000000000   0.88160000000000  H2O   (8f)
  -0.34630000000000  -0.32790000000000  -0.38160000000000  H2O   (8f)
  -0.32790000000000  -0.34630000000000  -0.88160000000000  H2O   (8f)
   0.34630000000000   0.32790000000000   1.38160000000000  H2O   (8f)
   0.50000000000000   0.50000000000000   0.00000000000000  Na    (4b)
   0.50000000000000   0.50000000000000   0.50000000000000  Na    (4b)
   0.00000000000000   0.50000000000000   0.00000000000000  Na    (4c)
   0.50000000000000   0.00000000000000   0.50000000000000  Na    (4c)
   0.00000000000000   0.50000000000000   0.50000000000000  Na    (4d)
   0.50000000000000   0.00000000000000   0.00000000000000  Na    (4d)
  -0.49410000000000   0.49410000000000   0.25000000000000  Na    (4e)
   0.49410000000000  -0.49410000000000   0.75000000000000  Na    (4e)
   0.00990000000000   0.49630000000000   0.25500000000000  Na    (8f)
  -0.49630000000000  -0.00990000000000   0.24500000000000  Na    (8f)
  -0.00990000000000  -0.49630000000000  -0.25500000000000  Na    (8f)
   0.49630000000000   0.00990000000000   0.75500000000000  Na    (8f)
   0.85730000000000   0.87490000000000  -0.07670000000000  O     (8f)
  -0.87490000000000  -0.85730000000000   0.57670000000000  O     (8f)
  -0.85730000000000  -0.87490000000000   0.07670000000000  O     (8f)
   0.87490000000000   0.85730000000000   0.42330000000000  O     (8f)
  -0.12340000000000   0.85660000000000   0.05860000000000  O     (8f)
  -0.85660000000000   0.12340000000000   0.44140000000000  O     (8f)
   0.12340000000000  -0.85660000000000  -0.05860000000000  O     (8f)
   0.85660000000000  -0.12340000000000   0.55860000000000  O     (8f)
  -0.44620000000000   0.54480000000000   0.17470000000000  O     (8f)
  -0.54480000000000   0.44620000000000   0.32530000000000  O     (8f)
   0.44620000000000  -0.54480000000000  -0.17470000000000  O     (8f)
   0.54480000000000  -0.44620000000000   0.67470000000000  O     (8f)
  -0.12620000000000   0.26840000000000   0.12850000000000  O     (8f)
  -0.26840000000000   0.12620000000000   0.37150000000000  O     (8f)
   0.12620000000000  -0.26840000000000  -0.12850000000000  O     (8f)
   0.26840000000000  -0.12620000000000   0.62850000000000  O     (8f)
  -0.13010000000000   0.25830000000000   0.25580000000000  O     (8f)
  -0.25830000000000   0.13010000000000   0.24420000000000  O     (8f)
   0.13010000000000  -0.25830000000000  -0.25580000000000  O     (8f)
   0.25830000000000  -0.13010000000000   0.75580000000000  O     (8f)
  -0.09560000000000   0.54860000000000   0.35440000000000  O     (8f)
  -0.54860000000000   0.09560000000000   0.14560000000000  O     (8f)
   0.09560000000000  -0.54860000000000  -0.35440000000000  O     (8f)
   0.54860000000000  -0.09560000000000   0.85440000000000  O     (8f)
   0.26400000000000   0.86000000000000   0.12030000000000  O     (8f)
  -0.86000000000000  -0.26400000000000   0.37970000000000  O     (8f)
  -0.26400000000000  -0.86000000000000  -0.12030000000000  O     (8f)
   0.86000000000000   0.26400000000000   0.62030000000000  O     (8f)
   0.27520000000000   0.86280000000000   0.26250000000000  O     (8f)
  -0.86280000000000  -0.27520000000000   0.23750000000000  O     (8f)
  -0.27520000000000  -0.86280000000000  -0.26250000000000  O     (8f)
   0.86280000000000   0.27520000000000   0.76250000000000  O     (8f)
   0.55010000000000   0.89770000000000   0.34550000000000  O     (8f)
  -0.89770000000000  -0.55010000000000   0.15450000000000  O     (8f)
  -0.55010000000000  -0.89770000000000  -0.34550000000000  O     (8f)
   0.89770000000000   0.55010000000000   0.84550000000000  O     (8f)
   0.39720000000000   0.79800000000000   0.22210000000000  Si    (8f)
  -0.79800000000000  -0.39720000000000   0.27790000000000  Si    (8f)
  -0.39720000000000  -0.79800000000000  -0.22210000000000  Si    (8f)
   0.79800000000000   0.39720000000000   0.72210000000000  Si    (8f)
  -0.19280000000000   0.39180000000000   0.22520000000000  Si    (8f)
  -0.39180000000000   0.19280000000000   0.27480000000000  Si    (8f)
   0.19280000000000  -0.39180000000000  -0.22520000000000  Si    (8f)
   0.39180000000000  -0.19280000000000   0.72520000000000  Si    (8f)
   0.19570000000000   0.19870000000000   0.57230000000000  Si    (8f)
  -0.19870000000000  -0.19570000000000  -0.07230000000000  Si    (8f)
  -0.19570000000000  -0.19870000000000  -0.57230000000000  Si    (8f)
   0.19870000000000   0.19570000000000   1.07230000000000  Si    (8f)
   0.00000000000000   0.00000000000000   0.00000000000000  Zr    (4a)
   0.00000000000000   0.00000000000000   0.50000000000000  Zr    (4a)
   0.00740000000000  -0.00740000000000   0.25000000000000  Zr    (4e)
  -0.00740000000000   0.00740000000000   0.75000000000000  Zr    (4e)
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Na_2PrO_3: A2B3C_mC48_15_aef_3f_2e - CIF

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_chemical_formula_sum 'Na2 O3 Pr'

loop_
_publ_author_name
 'Y. Hinatsu'
 'Y. Doi'
_journal_name_full_name
;
 Journal of Alloys and Compounds
;
_journal_volume 418
_journal_year 2006
_journal_page_first 155
_journal_page_last 160
_publ_Section_title
;
 Crystal structures and magnetic properties of alkali-metal lanthanide
     ↪ oxides $A_{2}$$Ln$O$_{3}$ ($A$ = Li, Na; $Ln$ = Ce, Pr, Tb)
;

_aflow_title 'Na$_{2}$PrO$_{3}$ Structure'
_aflow_proto 'A2B3C_mC48_15_aef_3f_2e'
_aflow_params 'a,b/a,c/a,\beta,y_{2},y_{3},y_{4},x_{5},y_{5},z_{5},x_{6
    ↪ },y_{6},z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},z_{8}'
_aflow_params_values '5.9649,1.72911532465,1.96968934936,110.09,0.8146,
    ↪ 0.1681,0.4994,-0.04,0.325,-0.004,0.241,0.501,0.141,0.251,0.135,
    ↪ 0.144,0.301,0.861,0.143'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mC48'

_symmetry_space_group_name_H-M "C 1 2/c 1"
_symmetry_Int_Tables_number 15

_cell_length_a     5.96490
_cell_length_b     10.31400
_cell_length_c     11.74900
_cell_angle_alpha  90.00000
_cell_angle_beta   110.09000
_cell_angle_gamma  90.00000

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_space_group_symop_id
_space_group_symop_operation_xyz
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2 -x,y,-z+1/2
3 -x,-y,-z
4 x,-y,z+1/2
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z+1/2
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z+1/2

loop_
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_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Na1 Na 4 a 0.00000 0.00000 0.00000 1.00000
Na2 Na 4 e 0.00000 0.81460 0.25000 1.00000
Pr1 Pr 4 e 0.00000 0.16810 0.25000 1.00000
Pr2 Pr 4 e 0.00000 0.49940 0.25000 1.00000
Na3 Na 8 f -0.04000 0.32500 -0.00400 1.00000
O1 O 8 f 0.24100 0.50100 0.14100 1.00000
O2 O 8 f 0.25100 0.13500 0.14400 1.00000
O3 O 8 f 0.30100 0.86100 0.14300 1.00000
```

Na_2PrO_3: A2B3C_mC48_15_aef_3f_2e - POSCAR

```
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    ↪ y7,z7,x8,y8,z8 --params=5.9649,1.72911532465,1.96968934936,
    ↪ 110.09,0.8146,0.1681,0.4994,-0.04,0.325,-0.004,0.241,0.501,
    ↪ 0.141,0.251,0.135,0.144,0.301,0.861,0.143 & C2/c C_{2h}^{6} #15
    ↪  (ae^3f^4) & mC48 & None & Na2O3Pr & Na2O3Pr & Y. Hinatsu and
    ↪ Y. Doi, J. Alloys Compd. 418, 155-160 (2006)
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   2.98245000000000  -5.15700000000000   0.00000000000000
   2.98245000000000   5.15700000000000   0.00000000000000
  -4.03573199541645   0.00000000000000  11.03412290402690
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    Na      O     Pr
     8     12      4
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   0.00000000000000   0.00000000000000   0.50000000000000   Na   (4a)
  -0.81460000000000   0.81460000000000   0.25000000000000   Na   (4e)
   0.81460000000000  -0.81460000000000   0.75000000000000   Na   (4e)
  -0.36500000000000   0.28500000000000  -0.00400000000000   Na   (8f)
  -0.28500000000000   0.36500000000000   0.50400000000000   Na   (8f)
   0.36500000000000  -0.28500000000000   0.00400000000000   Na   (8f)
   0.28500000000000  -0.36500000000000   0.49600000000000   Na   (8f)
  -0.26000000000000   0.74200000000000   0.14100000000000   O    (8f)
  -0.74200000000000   0.26000000000000   0.35900000000000   O    (8f)
   0.26000000000000  -0.74200000000000  -0.14100000000000   O    (8f)
   0.74200000000000  -0.26000000000000   0.64100000000000   O    (8f)
   0.11600000000000   0.38600000000000   0.14400000000000   O    (8f)
  -0.38600000000000  -0.11600000000000   0.35600000000000   O    (8f)
  -0.11600000000000  -0.38600000000000  -0.14400000000000   O    (8f)
   0.38600000000000   0.11600000000000   0.64400000000000   O    (8f)
  -0.56000000000000   1.16200000000000   0.14300000000000   O    (8f)
  -1.16200000000000   0.56000000000000   0.35700000000000   O    (8f)
   0.56000000000000  -1.16200000000000  -0.14300000000000   O    (8f)
   1.16200000000000  -0.56000000000000   0.64300000000000   O    (8f)
  -0.16810000000000   0.16810000000000   0.25000000000000   Pr   (4e)
   0.16810000000000  -0.16810000000000   0.75000000000000   Pr   (4e)
  -0.49940000000000   0.49940000000000   0.25000000000000   Pr   (4e)
   0.49940000000000  -0.49940000000000   0.75000000000000   Pr   (4e)
```

Eudidymite ($BeHNaO_8Si_3$): A2B4C2D17E6_mC124_15_f_2f_f_e8f_3f - CIF

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_audit_creation_method FINDSYM

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_chemical_formula_sum 'Be2 H4 Na2 O17 Si6'

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_publ_author_name
 'G. {Diego Gatta}'
 'N. Rotiroti'
 'G. J. {McIntyre}'
 'A. Guastoni'
 'F. Nestola'
_journal_name_full_name
;
 American Mineralogist
;
_journal_volume 93
_journal_year 2008
_journal_page_first 1158
_journal_page_last 1165
_publ_Section_title
;
 New insights into the crystal chemistry of epididymite and eudidymite
      ↪ from Malosa, Malawi: A single-crystal neutron diffraction
      ↪ study
;

_aflow_title 'Eudidymite (BeHNaO$_{8}$Si$_{3}$) Structure'
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      ↪ },x_{4},y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{6},x_{7},
      ↪ y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{10},y_{10},
      ↪ z_{10},x_{11},y_{11},z_{11},x_{12},y_{12},z_{12},x_{13},y_{13},
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      ↪ z_{16}'
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      ↪ 0.17321,0.16817,0.32317,0.4971,0.084,0.4645,0.7449,0.4773,
      ↪ 0.0419,0.7486,0.02899,0.34078,-0.07147,0.0727,0.03247,0.62614,
      ↪ 0.24036,0.1803,0.1154,0.24045,0.16471,0.57054,0.23677,0.43859,
      ↪ 0.24484,0.1548,0.49864,0.05623,0.12854,0.05189,0.87541,0.05237,
      ↪ 0.25694,0.43729,0.01209,0.43098,0.74466,0.20415,0.01965,0.64127
      ↪ ,0.25166,0.1025,0.86648,0.02964,0.09397,0.36195'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mC124'

_symmetry_space_group_name_H-M "C 1 2/c 1"
_symmetry_Int_Tables_number 15

_cell_length_a    12.61880
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_cell_length_c    13.99400
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_cell_angle_beta  103.76200
_cell_angle_gamma 90.00000

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_space_group_symop_id
_space_group_symop_operation_xyz
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2 -x,y,-z+1/2
3 -x,-y,-z
4 x,-y,z+1/2
5 x+1/2,y+1/2,z
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7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
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Be1 Be 8 f 0.16817 0.32317 0.49710 1.00000
H1 H 8 f 0.08400 0.46450 0.74490 1.00000
H2 H 8 f 0.47730 0.04190 0.74860 1.00000
Na1 Na 8 f 0.02899 0.34078 -0.07147 1.00000
O2 O 8 f 0.07270 0.03247 0.62614 1.00000
O3 O 8 f 0.24036 0.18030 0.11540 1.00000
O4 O 8 f 0.24045 0.16471 0.57054 1.00000
O5 O 8 f 0.23677 0.43859 0.24484 1.00000
O6 O 8 f 0.15480 0.49864 0.05623 1.00000
O7 O 8 f 0.12854 0.05189 0.87541 1.00000
O8 O 8 f 0.05237 0.25694 0.43729 1.00000
O9 O 8 f 0.01209 0.43098 0.74466 1.00000
Si1 Si 8 f 0.20415 0.01965 0.64127 1.00000
Si2 Si 8 f 0.25166 0.10250 0.86648 1.00000
Si3 Si 8 f 0.02964 0.09397 0.36195 1.00000
```

Eudidymite ($BeHNaO_8Si_3$): A2B4C2D17E6_mC124_15_f_2f_f_e8f_3f - POSCAR

```
A2B4C2D17E6_mC124_15_f_2f_f_e8f_3f & a,b/a,c/a,beta,y1,x2,y2,z2,x3,y3,z3
      ↪ ,x4,y4,z4,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,
      ↪ z10,x11,y11,z11,x12,y12,z12,x13,y13,z13,x14,y14,z14,x15,y15,z15
      ↪ ,x16,y16,z16 --params=12.6188,0.584691095825,1.10898025169,
      ↪ 103.762,0.17321,0.16817,0.32317,0.4971,0.084,0.4645,0.7449,
      ↪ 0.4773,0.0419,0.7486,0.02899,0.34078,-0.07147,0.0727,0.03247,
      ↪ 0.62614,0.24036,0.1803,0.1154,0.24045,0.16471,0.57054,0.23677,
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      ↪ 0.05237,0.25694,0.43729,0.01209,0.43098,0.74466,0.20415,0.01965
      ↪ ,0.64127,0.25166,0.1025,0.86648,0.02964,0.09397,0.36195 & C2/c
      ↪ C_{2h}^{6} #15 (ef^15) & mC124 & None & BeHNaO8Si3 & Eudidymite
      ↪ & G. {Diego Gatta} et al., Am. Mineral. 93, 1158-1165 (2008)
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  6.30940000000000   3.68905000000000   0.00000000000000
 -3.32902320731927   0.00000000000000  13.59226399409350
    Be      H     Na      O     Si
     4      8      4     34     12
Direct
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   0.15500000000000  -0.49134000000000  -0.49710000000000   Be   (8f)
   0.49134000000000  -0.15500000000000   0.99710000000000   Be   (8f)
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   0.38050000000000  -0.54850000000000  -0.74490000000000   H    (8f)
   0.54850000000000  -0.38050000000000   1.24490000000000   H    (8f)
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  -0.43540000000000  -0.51920000000000  -0.74860000000000   H    (8f)
   0.51920000000000   0.43540000000000   1.24860000000000   H    (8f)
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  -0.36977000000000   0.31179000000000   0.57147000000000   Na   (8f)
   0.31179000000000  -0.36977000000000   0.07147000000000   Na   (8f)
   0.36977000000000  -0.31179000000000   0.42853000000000   Na   (8f)
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   0.17321000000000  -0.17321000000000   0.75000000000000   O    (4e)
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  -0.04023000000000  -0.10517000000000  -0.62614000000000   O    (8f)
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   0.07574000000000   0.40516000000000   0.57054000000000   O    (8f)
  -0.40516000000000  -0.07574000000000  -0.07054000000000   O    (8f)
  -0.07574000000000  -0.40516000000000  -0.57054000000000   O    (8f)
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   0.20182000000000  -0.67536000000000  -0.24484000000000   O    (8f)
   0.67536000000000  -0.20182000000000   0.74484000000000   O    (8f)
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  -0.07665000000000  -0.18043000000000  -0.87541000000000   O    (8f)
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   0.20457000000000  -0.30931000000000  -0.43729000000000   O    (8f)
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   0.41889000000000  -0.44307000000000  -0.74466000000000   O    (8f)
   0.44307000000000  -0.41889000000000   1.24466000000000   O    (8f)
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  -0.22380000000000  -0.18450000000000  -0.14127000000000   Si   (8f)
  -0.18450000000000  -0.22380000000000  -0.64127000000000   Si   (8f)
   0.22380000000000   0.18450000000000   1.14127000000000   Si   (8f)
   0.14916000000000   0.35416000000000   0.86648000000000   Si   (8f)
  -0.35416000000000  -0.14916000000000  -0.36648000000000   Si   (8f)
  -0.14916000000000  -0.35416000000000  -0.86648000000000   Si   (8f)
   0.35416000000000   0.14916000000000   1.36648000000000   Si   (8f)
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  -0.12361000000000   0.06433000000000   0.13805000000000   Si   (8f)
   0.06433000000000  -0.12361000000000  -0.36195000000000   Si   (8f)
   0.12361000000000  -0.06433000000000   0.86195000000000   Si   (8f)
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ζ-Nb_2O_5 (B-Nb_2O_5): A2B5_mC28_15_f_e2f - CIF

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 'T. S. Ercit'
_journal_name_full_name
;
 Mineralogy and Petrology
;
_journal_volume 43
_journal_year 1991
_journal_page_first 217
_journal_page_last 223
_publ_Section_title
;
 Refinement of the structure of $\zeta$-Nb$_{2}$O$_{5}$ and its
      ↪ relationship to the rutile and thoreaulite structures
;

# Found in ResearchGate, {~},

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      ↪ 0.3598,0.25874,0.2512,0.1088,0.4421,0.4709,0.2055,0.0723,0.1256
      ↪ '
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mC28'

_symmetry_space_group_name_H-M "C 1 2/c 1"
_symmetry_Int_Tables_number 15

_cell_length_a      12.74000
_cell_length_b      4.88300
_cell_length_c      5.56090
_cell_angle_alpha   90.00000
_cell_angle_beta    105.02000
_cell_angle_gamma   90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z+1/2
3 -x,-y,-z
4 x,-y,z+1/2
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z+1/2
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 4 e 0.00000 0.64650 0.25000 1.00000
Nb1 Nb 8 f 0.35980 0.25874 0.25120 1.00000
O2 O 8 f 0.10880 0.44210 0.47090 1.00000
O3 O 8 f 0.20550 0.07230 0.12560 1.00000
```

ζ-Nb_2O_5 (B-Nb_2O_5): A2B5_mC28_15_f_e2f - POSCAR

```
A2B5_mC28_15_f_e2f & a,b/a,c/a,beta,y1,x2,y2,z2,x3,y3,z3,x4,y4,z4 --
      ↪ params=12.74,0.38328100471,0.436491365777,105.02,0.6465,0.3598,
      ↪ 0.25874,0.2512,0.1088,0.4421,0.4709,0.2055,0.0723,0.1256 & C2/c
      ↪ C_{2h}^{6} #15 (ef^3) & mC28 & None & Nb2O5 & Nb2O5 & T. S.
      ↪ Ercit, Mineral. Petrol. 43, 217-223 (1991)
   1.00000000000000
   6.37000000000000   -2.44150000000000    0.00000000000000
   6.37000000000000    2.44150000000000    0.00000000000000
  -1.44114171840500    0.00000000000000    5.37091420127642
     Nb      O
      4     10
Direct
   0.10106000000000    0.61854000000000    0.25120000000000   Nb   (8f)
  -0.61854000000000   -0.10106000000000    0.24880000000000   Nb   (8f)
  -0.10106000000000   -0.61854000000000   -0.25120000000000   Nb   (8f)
   0.61854000000000    0.10106000000000    0.75120000000000   Nb   (8f)
  -0.64650000000000    0.64650000000000    0.25000000000000   O    (4e)
   0.64650000000000   -0.64650000000000    0.75000000000000   O    (4e)
  -0.33330000000000    0.55090000000000    0.47090000000000   O    (8f)
  -0.55090000000000    0.33330000000000    0.02910000000000   O    (8f)
   0.33330000000000   -0.55090000000000   -0.47090000000000   O    (8f)
   0.55090000000000   -0.33330000000000    0.97090000000000   O    (8f)
   0.13320000000000    0.27780000000000    0.12560000000000   O    (8f)
  -0.27780000000000   -0.13320000000000    0.37440000000000   O    (8f)
  -0.13320000000000   -0.27780000000000   -0.12560000000000   O    (8f)
   0.27780000000000    0.13320000000000    0.62560000000000   O    (8f)
```

Muscovite ($KH_2Al_3Si_3O_{12}$, $S5_1$): A2BC10D2E4_mC76_15_f_e_5f_f_2f - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Muscovite'
_chemical_formula_sum 'Al2 K O10 (OH)2 Si4'

loop_
_publ_author_name
 'S. M. Richardson'
 'J. W. {Richardson, Jr.}'
_journal_name_full_name
;
 American Mineralogist
;
_journal_volume 67
_journal_year 1982
_journal_page_first 69
_journal_page_last 75
_publ_Section_title
;
 Crystal structure of a pink muscovite from Archer\'s Post, Kenya:
      ↪ Implications for reverse pleochroism in dioctahedral micas
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'Muscovite (KH$_{2}$Al$_{3}$Si$_{3}$O$_{12}$, $S5_{1}$)
      ↪ Structure'
_aflow_proto 'A2BC10D2E4_mC76_15_f_e_5f_f_2f'
_aflow_params 'a,b/a,c/a,\beta,y_{1},x_{2},y_{2},z_{2},x_{3},y_{3},z_{3
      ↪ },x_{4},y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{6},x_{7},
      ↪ y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{10},y_{10},
      ↪ z_{10}'
_aflow_params_values '5.1988,1.73628529661,3.86739247519,95.782,0.0992,
      ↪ 0.2506,0.0838,0.0002,0.3872,0.2525,0.0543,0.0366,0.4431,0.4459,
      ↪ 0.4178,0.0931,0.1685,0.2475,0.3712,0.1685,0.2509,0.3132,0.3424,
      ↪ 0.0422,0.0622,0.4492,0.451,0.2587,0.1355,0.0354,0.4298,0.3646'
_aflow_Strukturbericht '$S5_{1}$'
_aflow_Pearson 'mC76'

_symmetry_space_group_name_H-M "C 1 2/c 1"
_symmetry_Int_Tables_number 15

_cell_length_a      5.19880
_cell_length_b      9.02660
_cell_length_c      20.10580
_cell_angle_alpha   90.00000
_cell_angle_beta    95.78200
_cell_angle_gamma   90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z+1/2
3 -x,-y,-z
4 x,-y,z+1/2
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z+1/2
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
K1 K 4 e 0.00000 0.09920 0.25000 1.00000
Al1 Al 8 f 0.25060 0.08380 0.00020 1.00000
O1 O 8 f 0.38720 0.25250 0.05430 1.00000
O2 O 8 f 0.03660 0.44310 0.44590 1.00000
O3 O 8 f 0.41780 0.09310 0.16850 1.00000
O4 O 8 f 0.24750 0.37120 0.16850 1.00000
O5 O 8 f 0.25090 0.31320 0.34240 1.00000
OH1 OH 8 f 0.04220 0.06220 0.44920 1.00000
Si1 Si 8 f 0.45100 0.25870 0.13550 1.00000
Si2 Si 8 f 0.03540 0.42980 0.36460 1.00000
```

Muscovite ($KH_2Al_3Si_3O_{12}$, $S5_1$): A2BC10D2E4_mC76_15_f_e_5f_f_2f - POSCAR

```
A2BC10D2E4_mC76_15_f_e_5f_f_2f & a,b/a,c/a,beta,y1,x2,y2,z2,x3,y3,z3,x4,
      ↪ y4,z4,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10
      ↪ --params=5.1988,1.73628529661,3.86739247519,95.782,0.0992,
      ↪ 0.2506,0.0838,0.0002,0.3872,0.2525,0.0543,0.0366,0.4431,0.4459,
      ↪ 0.4178,0.0931,0.1685,0.2475,0.3712,0.1685,0.2509,0.3132,0.3424,
      ↪ 0.0422,0.0622,0.4492,0.451,0.2587,0.1355,0.0354,0.4298,0.3646 &
      ↪  C2/c C_{2h}^{6} #15 (ef^9) & mC76 & $S5_{1}$ & Al3KO10(OH)2Si3
      ↪ & Muscovite & S. M. Richardson and J. W. {Richardson, Jr.},
      ↪ Am. Mineral. 67, 69-75 (1982)
   1.00000000000000
   2.59940000000000   -4.51330000000000    0.00000000000000
   2.59940000000000    4.51330000000000    0.00000000000000
  -2.02553351199094    0.00000000000000   20.00350988281310
     Al      K      O     OH     Si
      4      2     20      4      8
Direct
   0.16680000000000    0.33440000000000    0.00020000000000   Al   (8f)
  -0.33440000000000   -0.16680000000000    0.49980000000000   Al   (8f)
  -0.16680000000000   -0.33440000000000   -0.00020000000000   Al   (8f)
   0.33440000000000    0.16680000000000    0.50020000000000   Al   (8f)
  -0.09920000000000    0.09920000000000    0.25000000000000    K   (4e)
```

```
   0.09920000000000  −0.09920000000000   0.75000000000000   K   (4e)
   0.13470000000000   0.63970000000000   0.05430000000000   O   (8f)
  −0.63970000000000  −0.13470000000000   0.44570000000000   O   (8f)
  −0.13470000000000  −0.63970000000000  −0.05430000000000   O   (8f)
   0.63970000000000   0.13470000000000   0.55430000000000   O   (8f)
  −0.40650000000000   0.47970000000000   0.44590000000000   O   (8f)
  −0.47970000000000   0.40650000000000   0.05410000000000   O   (8f)
   0.40650000000000  −0.47970000000000  −0.44590000000000   O   (8f)
   0.47970000000000  −0.40650000000000   0.94590000000000   O   (8f)
   0.32470000000000   0.51090000000000   0.16850000000000   O   (8f)
  −0.51090000000000  −0.32470000000000   0.33150000000000   O   (8f)
  −0.32470000000000  −0.51090000000000  −0.16850000000000   O   (8f)
   0.51090000000000   0.32470000000000   0.66850000000000   O   (8f)
  −0.12370000000000   0.61870000000000   0.16850000000000   O   (8f)
  −0.61870000000000   0.12370000000000   0.33150000000000   O   (8f)
   0.12370000000000  −0.61870000000000  −0.16850000000000   O   (8f)
   0.61870000000000  −0.12370000000000   0.66850000000000   O   (8f)
  −0.06230000000000   0.56410000000000   0.34240000000000   O   (8f)
  −0.56410000000000   0.06230000000000   0.15760000000000   O   (8f)
   0.06230000000000  −0.56410000000000  −0.34240000000000   O   (8f)
   0.56410000000000  −0.06230000000000   0.84240000000000   O   (8f)
  −0.02000000000000   0.10440000000000   0.44920000000000   OH  (8f)
  −0.10440000000000   0.02000000000000   0.05080000000000   OH  (8f)
   0.02000000000000  −0.10440000000000  −0.44920000000000   OH  (8f)
   0.10440000000000  −0.02000000000000   0.94920000000000   OH  (8f)
   0.19230000000000   0.70970000000000   0.13550000000000   Si  (8f)
  −0.70970000000000  −0.19230000000000   0.36450000000000   Si  (8f)
  −0.19230000000000  −0.70970000000000  −0.13550000000000   Si  (8f)
   0.70970000000000   0.19230000000000   0.63550000000000   Si  (8f)
  −0.39440000000000   0.46520000000000   0.36460000000000   Si  (8f)
  −0.46520000000000   0.39440000000000   0.13540000000000   Si  (8f)
   0.39440000000000  −0.46520000000000  −0.36460000000000   Si  (8f)
   0.46520000000000  −0.39440000000000   0.86460000000000   Si  (8f)
```

$Rb_2C_2O_4{\cdot}H_2O$: A2BC4D2_mC36_15_f_e_2f_f - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'C2(H2O)O4Rb2'
_chemical_formula_sum 'C2 (H2O) O4 Rb2'

loop_
_publ_author_name
 'B. F. Pedersen'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 19
_journal_year 1965
_journal_page_first 1815
_journal_page_last 1818
_publ_Section_title
;
 The Crystal Structure of Rubidium. Oxalate Monohydrate, Rb$_{2}$C$_{2}
    ↪ $O$_{4}\cdot$H$_{2}$O
;

_aflow_title 'Rb$_{2}$C$_{2}$O$_{4}\cdot$H$_{2}$O Structure'
_aflow_proto 'A2BC4D2_mC36_15_f_e_2f_f'
_aflow_params 'a,b/a,c/a,\beta,y_{1},x_{2},y_{2},z_{2},x_{3},y_{3},z_{3
    ↪ },x_{4},y_{4},z_{4},x_{5},y_{5},z_{5}'
_aflow_params_values '9.662,0.657213827365,1.147588491,109.4,0.4733,
    ↪ 0.2396,0.3216,0.0546,0.1355,0.2748,0.0933,0.3262,0.4768,0.094,
    ↪ 0.129,0.8156,0.1297'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mC36'

_symmetry_space_group_name_H-M "C 1 2/c 1"
_symmetry_Int_Tables_number 15

_cell_length_a      9.66200
_cell_length_b      6.35000
_cell_length_c      11.08800
_cell_angle_alpha  90.00000
_cell_angle_beta   109.40000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z+1/2
3 -x,-y,-z
4 x,-y,z+1/2
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z+1/2
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
H2O1 H2O 4 e 0.00000 0.47330 0.25000 1.00000
C1 C 8 f 0.23960 0.32160 0.05460 1.00000
O1 O 8 f 0.13550 0.27480 0.09330 1.00000
O2 O 8 f 0.32620 0.47680 0.09400 1.00000
Rb1 Rb 8 f 0.12900 0.81560 0.12970 1.00000
```

$Rb_2C_2O_4{\cdot}H_2O$: A2BC4D2_mC36_15_f_e_2f_f - POSCAR

```
A2BC4D2_mC36_15_f_e_2f_f & a,b/a,c/a,beta,y1,x2,y2,z2,x3,y3,z3,x4,y4,z4,
    ↪ x5,y5,z5 --params=9.662,0.657213827365,1.147588491,109.4,0.4733
    ↪ ,0.2396,0.3216,0.0546,0.1355,0.2748,0.0933,0.3262,0.4768,0.094,
    ↪ 0.129,0.8156,0.1297 & C2/c C_{2h}^{6} #15 (ef^4) & mC36 & None
    ↪ & C2(H2O)O4Rb2 & C2(H2O)O4Rb2 & B. F. Pedersen, Acta Chem.
    ↪ Scand. 19, 1815-1818 (1965)
   1.00000000000000
   4.83100000000000  -3.17500000000000   0.00000000000000
   4.83100000000000   3.17500000000000   0.00000000000000
  -3.68300263032650   0.00000000000000  10.45845283132300
     C    H2O      O     Rb
     4      2      8      4
Direct
  -0.08200000000000   0.56120000000000   0.05460000000000   C   (8f)
  -0.56120000000000   0.08200000000000   0.44540000000000   C   (8f)
   0.08200000000000  -0.56120000000000  -0.05460000000000   C   (8f)
   0.56120000000000  -0.08200000000000   0.55460000000000   C   (8f)
  -0.47330000000000   0.47330000000000   0.25000000000000   H2O (4e)
   0.47330000000000  -0.47330000000000   0.75000000000000   H2O (4e)
  -0.13930000000000   0.41030000000000   0.09330000000000   O   (8f)
  -0.41030000000000   0.13930000000000   0.40670000000000   O   (8f)
   0.13930000000000  -0.41030000000000  -0.09330000000000   O   (8f)
   0.41030000000000  -0.13930000000000   0.59330000000000   O   (8f)
  -0.15060000000000   0.80300000000000   0.09400000000000   O   (8f)
  -0.80300000000000   0.15060000000000   0.40600000000000   O   (8f)
   0.15060000000000  -0.80300000000000  -0.09400000000000   O   (8f)
   0.80300000000000  -0.15060000000000   0.59400000000000   O   (8f)
  -0.68660000000000   0.94460000000000   0.12970000000000   Rb  (8f)
  -0.94460000000000   0.68660000000000   0.37030000000000   Rb  (8f)
   0.68660000000000  -0.94460000000000  -0.12970000000000   Rb  (8f)
   0.94460000000000  -0.68660000000000   0.62970000000000   Rb  (8f)
```

Alluaudite $[NaMnFe_2(PO_4)_3]$: A2BCD12E3_mC76_15_f_e_b_6f_ef - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Alluaudite'
_chemical_formula_sum 'Fe2 Mn Na O12 P3'

loop_
_publ_author_name
 'P. B. Moore'
_journal_name_full_name
;
 American Mineralogist
;
_journal_volume 56
_journal_year 1971
_journal_page_first 1955
_journal_page_last 1975
_publ_Section_title
;
 Crystal Chemistry of the Alluaudite Structure Type: Contribution to the
    ↪ Paragenesis of Pegmatite Phosphate Giant Crystals
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'Alluaudite [NaMnFe$_{2}$(PO$_{4}$)$_{3}$] Structure'
_aflow_proto 'A2BCD12E3_mC76_15_f_e_b_6f_ef'
_aflow_params 'a,b/a,c/a,\beta,y_{2},y_{3},x_{4},y_{4},z_{4},x_{5},y_{5
    ↪ },z_{5},x_{6},y_{6},z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},z_{8},
    ↪ x_{9},y_{9},z_{9},x_{10},y_{10},z_{10},x_{11},y_{11},z_{11}'
_aflow_params_values '12.004,1.04406864379,0.533488837054,114.4,0.2599,
    ↪ 0.7145,0.2812,0.6525,0.3713,0.4533,0.7152,0.5342,0.0988,0.6375,
    ↪ 0.2401,0.3272,0.6633,0.102,0.1213,0.3974,0.3119,0.2251,0.822,
    ↪ 0.3172,0.3102,0.5021,0.3735,0.2424,0.8911,0.1325'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mC76'

_symmetry_space_group_name_H-M "C 1 2/c 1"
_symmetry_Int_Tables_number 15

_cell_length_a      12.00400
_cell_length_b      12.53300
_cell_length_c      6.40400
_cell_angle_alpha  90.00000
_cell_angle_beta   114.40000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z+1/2
3 -x,-y,-z
4 x,-y,z+1/2
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z+1/2
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
```

```
_atom_site_fract_z
_atom_site_occupancy
Na1 Na 4 b 0.00000 0.50000 0.00000 1.00000
Mn1 Mn 4 e 0.00000 0.25990 0.25000 1.00000
P1 P 4 e 0.00000 0.71450 0.25000 1.00000
Fe1 Fe 8 f 0.28120 0.65250 0.37130 1.00000
O1 O 8 f 0.45330 0.71520 0.53420 1.00000
O2 O 8 f 0.09880 0.63750 0.24010 1.00000
O3 O 8 f 0.32720 0.66330 0.10200 1.00000
O4 O 8 f 0.12130 0.39740 0.31190 1.00000
O5 O 8 f 0.22510 0.82200 0.31720 1.00000
O6 O 8 f 0.31020 0.50210 0.37350 1.00000
P2 P 8 f 0.24240 0.89110 0.13250 1.00000
```

Alluaudite [$NaMnFe_2(PO_4)_3$]: A2BCD12E3_mC76_15_f_e_b_6f_ef - POSCAR

```
A2BCD12E3_mC76_15_f_e_b_6f_ef & a,b/a,c/a,beta,y2,y3,x4,y4,z4,x5,y5,z5,
    ↪ x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10,x11,y11,z11 --
    ↪ params=12.004,1.04406864379,0.533488837054,114.4,0.2599,0.7145,
    ↪ 0.2812,0.6525,0.3713,0.4533,0.7152,0.5342,0.0988,0.6375,0.2401,
    ↪ 0.3272,0.6633,0.102,0.1213,0.3974,0.3119,0.2251,0.822,0.3172,
    ↪ 0.3102,0.5021,0.3735,0.2424,0.8911,0.1325 & C2/c C_{2h}^{6} #15
    ↪ (be^2f^8) & mC76 & None & Fe2MnNaO12P3 & Alluaudite & P. B.
    ↪ Moore, Am. Mineral. 56, 1955–1975 (1971)
  1.00000000000000
  6.00200000000000  -6.26650000000000   0.00000000000000
  6.00200000000000   6.26650000000000   0.00000000000000
 -2.64552076859637   0.00000000000000   5.83201816380276
   Fe  Mn  Na  O  P
    4   2   2 24  6
Direct
 -0.37130000000000   0.93370000000000   0.37130000000000  Fe  (8f)
 -0.93370000000000   0.37130000000000   0.12870000000000  Fe  (8f)
  0.37130000000000  -0.93370000000000  -0.37130000000000  Fe  (8f)
  0.93370000000000  -0.37130000000000   0.87130000000000  Fe  (8f)
 -0.25990000000000   0.25990000000000   0.25000000000000  Mn  (4e)
  0.25990000000000  -0.25990000000000   0.75000000000000  Mn  (4e)
  0.50000000000000   0.50000000000000   0.00000000000000  Na  (4b)
  0.50000000000000   0.50000000000000   0.50000000000000  Na  (4b)
 -0.26190000000000   1.16850000000000   0.53420000000000  O  (8f)
 -1.16850000000000   0.26190000000000  -0.03420000000000  O  (8f)
  0.26190000000000  -1.16850000000000  -0.53420000000000  O  (8f)
  1.16850000000000  -0.26190000000000   1.03420000000000  O  (8f)
 -0.53870000000000   0.73630000000000   0.24010000000000  O  (8f)
 -0.73630000000000   0.53870000000000   0.25990000000000  O  (8f)
  0.53870000000000  -0.73630000000000  -0.24010000000000  O  (8f)
  0.73630000000000  -0.53870000000000   0.74010000000000  O  (8f)
 -0.33610000000000   0.99050000000000   0.10200000000000  O  (8f)
 -0.99050000000000   0.33610000000000   0.39800000000000  O  (8f)
  0.33610000000000  -0.99050000000000  -0.10200000000000  O  (8f)
  0.99050000000000  -0.33610000000000   0.60200000000000  O  (8f)
 -0.27610000000000   0.51870000000000   0.31190000000000  O  (8f)
 -0.51870000000000   0.27610000000000   0.18810000000000  O  (8f)
  0.27610000000000  -0.51870000000000  -0.31190000000000  O  (8f)
  0.51870000000000  -0.27610000000000   0.81190000000000  O  (8f)
 -0.59690000000000   1.04710000000000   0.31720000000000  O  (8f)
 -1.04710000000000   0.59690000000000   0.18280000000000  O  (8f)
  0.59690000000000  -1.04710000000000  -0.31720000000000  O  (8f)
  1.04710000000000  -0.59690000000000   0.81720000000000  O  (8f)
 -0.19190000000000   0.81230000000000   0.37350000000000  O  (8f)
 -0.81230000000000   0.19190000000000   0.12650000000000  O  (8f)
  0.19190000000000  -0.81230000000000  -0.37350000000000  O  (8f)
  0.81230000000000  -0.19190000000000   0.87350000000000  O  (8f)
 -0.71450000000000   0.71450000000000   0.25000000000000  P  (4e)
  0.71450000000000  -0.71450000000000   0.75000000000000  P  (4e)
 -0.64870000000000   1.13350000000000   0.13250000000000  P  (8f)
 -1.13350000000000   0.64870000000000   0.36750000000000  P  (8f)
  0.64870000000000  -1.13350000000000  -0.13250000000000  P  (8f)
  1.13350000000000  -0.64870000000000   0.63250000000000  P  (8f)
```

ThC_2 (C_g): A2B_mC12_15_f_e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'C2Th'
_chemical_formula_sum 'C2 Th'

loop_
_publ_author_name
 'A. L. Bowman'
 'N. H. Krikorian'
 'G. P. Arnold'
 'T. C. Wallace'
 'N. G. Nereson'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 24
_journal_year 1968
_journal_page_first 1121
_journal_page_last 1123
_publ_Section_title
;
 The Crystal Structure of ThC$_{2}$
;

_aflow_title 'ThC$_{2}$ (C$_{g}$) Structure'
_aflow_proto 'A2B_mC12_15_f_e'
_aflow_params 'a,b/a,c/a,\beta,y_{1},x_{2},y_{2},z_{2}'
_aflow_params_values '6.692,0.631052002391,1.00777047221,103.12,0.2074,
    ↪ 0.2992,0.1326,0.054'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mC12'

_symmetry_space_group_name_H-M "C 1 2/c 1"
_symmetry_Int_Tables_number 15

_cell_length_a    6.69200
_cell_length_b    4.22300
_cell_length_c    6.74400
_cell_angle_alpha 90.00000
_cell_angle_beta  103.12000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z+1/2
3 -x,-y,-z
4 x,-y,z+1/2
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z+1/2
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Th1 Th 4 e 0.00000 0.20740 0.25000 1.00000
C1 C 8 f 0.29920 0.13260 0.05400 1.00000
```

ThC_2 (C_g): A2B_mC12_15_f_e - POSCAR

```
A2B_mC12_15_f_e & a,b/a,c/a,beta,y1,x2,y2,z2 --params=6.692,
    ↪ 0.631052002391,1.00777047221,103.12,0.2074,0.2992,0.1326,0.054
    ↪ & C2/c C_{2h}^{6} #15 (ef) & mC12 & None & C2Th & C2Th & A. L.
    ↪ Bowman et al., Acta Crystallogr. Sect. B Struct. Sci. 24,
    ↪ 1121–1123 (1968)
  1.00000000000000
  3.34600000000000  -2.11150000000000   0.00000000000000
  3.34600000000000   2.11150000000000   0.00000000000000
 -1.53082916110101   0.00000000000000   6.56795996330084
   C  Th
   4   2
Direct
  0.16660000000000   0.43180000000000   0.05400000000000  C  (8f)
 -0.43180000000000  -0.16660000000000   0.44600000000000  C  (8f)
 -0.16660000000000  -0.43180000000000  -0.05400000000000  C  (8f)
  0.43180000000000   0.16660000000000   0.55400000000000  C  (8f)
 -0.20740000000000   0.20740000000000   0.25000000000000  Th  (4e)
  0.20740000000000  -0.20740000000000   0.75000000000000  Th  (4e)
```

Clinocervantite (β-Sb_2O_4): A2B_mC24_15_2f_ce - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Clinocervantite'
_chemical_formula_sum 'O2 Sb'

loop_
_publ_author_name
 'R. Basso'
 'G. Lucchetti'
 'L. Zefiro'
 'A. Palenzona'
_journal_name_full_name
;
 European Journal of Mineralogy
;
_journal_volume 11
_journal_year 1999
_journal_page_first 95
_journal_page_last 100
_publ_Section_title
;
 Clinocervantite, $\beta$-Sb$_{2}$O$_{4}$, the natural monoclinic
    ↪ polymorph of cervantite from the Cetine mine, Siena, Italy
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'Clinocervantite ($\beta$-Sb$_{2}$O$_{4}$) Structure'
_aflow_proto 'A2B_mC24_15_2f_ce'
_aflow_params 'a,b/a,c/a,\beta,y_{2},x_{3},y_{3},z_{3},x_{4},y_{4},z_{4
    ↪ }'
_aflow_params_values '12.061,0.40096177763,0.446314567615,103.12,0.2851,
    ↪ 0.1918,0.0517,0.6746,0.0939,0.4122,-0.0351'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mC24'

_symmetry_space_group_name_H-M "C 1 2/c 1"
_symmetry_Int_Tables_number 15

_cell_length_a    12.06100
_cell_length_b    4.83600
_cell_length_c    5.38300
_cell_angle_alpha 90.00000
_cell_angle_beta  103.12000
```

```
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z+1/2
3 -x,-y,-z
4 x,-y,z+1/2
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z+1/2
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Sb1 Sb 4 c 0.25000 0.25000 0.00000 1.00000
Sb2 Sb 4 e 0.00000 0.28510 0.25000 1.00000
O1 O 8 f 0.19180 0.05170 0.67460 1.00000
O2 O 8 f 0.09390 0.41220 -0.03510 1.00000
```

Clinocervantite (β-Sb_2O_4): A2B_mC24_15_2f_ce - POSCAR

```
A2B_mC24_15_2f_ce & a,b/a,c/a,beta,y2,x3,y3,z3,x4,y4,z4 --params=12.061,
    ↪ 0.40096177763,0.446314567615,103.12,0.2851,0.1918,0.0517,0.6746
    ↪ ,0.0939,0.4122,-0.0351 & C2/c C_{2h}^{6} #15 (cef^2) & mC24 &
    ↪ None & O2Sb & Clinocervantite & R. Basso et al., Eur. J.
    ↪ Mineral. 11, 95-100 (1999)
   1.00000000000000
   6.03050000000000  -2.41800000000000   0.00000000000000
   6.03050000000000   2.41800000000000   0.00000000000000
  -1.22189403532129   0.00000000000000   5.24248642978179
     O     Sb
     8      4
Direct
   0.14010000000000   0.24350000000000   0.67460000000000   O   (8f)
  -0.24350000000000  -0.14010000000000  -0.17460000000000   O   (8f)
  -0.14010000000000  -0.24350000000000  -0.67460000000000   O   (8f)
   0.24350000000000   0.14010000000000   1.17460000000000   O   (8f)
  -0.31830000000000   0.50610000000000  -0.03510000000000   O   (8f)
  -0.50610000000000   0.31830000000000   0.53510000000000   O   (8f)
   0.31830000000000  -0.50610000000000   0.03510000000000   O   (8f)
   0.50610000000000  -0.31830000000000   0.46490000000000   O   (8f)
   0.00000000000000   0.50000000000000   0.00000000000000  Sb   (4c)
   0.50000000000000   0.00000000000000   0.50000000000000  Sb   (4c)
  -0.28510000000000   0.28510000000000   0.25000000000000  Sb   (4e)
   0.28510000000000  -0.28510000000000   0.75000000000000  Sb   (4e)
```

($CdSO_4$)$_3$·8H_2O ($H4_{20}$): A3B16C20D3_mC168_15_ef_8f_10f_ef - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Cd3H16O20S3'
_chemical_formula_sum 'Cd3 H16 O20 S3'

loop_
_publ_author_name
 'R. Caminiti'
 'G. Johansson'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 35a
_journal_year 1981
_journal_page_first 451
_journal_page_last 455
_publ_Section_title
;
 A refinement of the Crystal Structure of the Cadmium Sulfate 3CdSO$_{4}
    ↪ $$\cdot$8H$_{2}$O
;

_aflow_title '(CdSO$_{4}$)$_{3}$$\cdot$8H$_{2}$O ($H4_{20}$) Structure'
_aflow_proto 'A3B16C20D3_mC168_15_ef_8f_10f_ef'
_aflow_params 'a,b/a,c/a,\beta,y_{1},y_{2},x_{3},y_{3},z_{3},x_{4},y_{4
    ↪ },z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{6},x_{7},y_{7},z_{7},
    ↪ x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{10},y_{10},z_{10},x_{11
    ↪ },y_{11},z_{11},x_{12},y_{12},z_{12},x_{13},y_{13},z_{13},x_{14
    ↪ },y_{14},z_{14},x_{15},y_{15},z_{15},x_{16},y_{16},z_{16},x_{17
    ↪ },y_{17},z_{17},x_{18},y_{18},z_{18},x_{19},y_{19},z_{19},x_{20
    ↪ },y_{20},z_{20},x_{21},y_{21},z_{21},x_{22},y_{22},z_{22}'
_aflow_params_values '14.818,0.803279794844,0.638952625186,97.39,-
    ↪ 0.05363,0.50828,0.15443,0.40511,0.035,0.083,0.252,0.716,0.095,
    ↪ 0.209,0.601,0.251,0.079,0.406,0.271,0.075,0.309,0.195,0.326,
    ↪ 0.288,0.273,0.347,0.273,0.094,0.203,0.346,0.06,0.3,0.402,0.0258
    ↪ ,0.4358,0.1357,0.4224,0.0799,0.1936,0.3962,0.4709,0.0521,0.1008
    ↪ ,0.2269,0.0033,0.0848,0.0936,0.1832,0.2296,0.1152,0.0982,0.0936
    ↪ ,0.1928,0.6827,0.2796,0.0746,0.4019,0.2214,0.3584,0.2573,0.0955
    ↪ ,0.2506,0.4077,0.12978,0.11705,0.05632'
_aflow_Strukturbericht '$H4_{20}$'
_aflow_Pearson 'mC168'

_symmetry_space_group_name_H-M "C 1 2/c 1"
_symmetry_Int_Tables_number 15
_cell_length_a     14.81800
_cell_length_b     11.90300
_cell_length_c     9.46800
_cell_angle_alpha  90.00000
_cell_angle_beta   97.39000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z+1/2
3 -x,-y,-z
4 x,-y,z+1/2
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z+1/2
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cd1 Cd 4 e 0.00000 -0.05363 0.25000 1.00000
S1 S 4 e 0.00000 0.50828 0.25000 1.00000
Cd2 Cd 8 f 0.15443 0.40511 0.03500 1.00000
H1 H 8 f 0.08300 0.25200 0.71600 1.00000
H2 H 8 f 0.09500 0.20900 0.60100 1.00000
H3 H 8 f 0.25100 0.07900 0.40600 1.00000
H4 H 8 f 0.27100 0.07500 0.30900 1.00000
H5 H 8 f 0.19500 0.32600 0.28800 1.00000
H6 H 8 f 0.27300 0.34700 0.27300 1.00000
H7 H 8 f 0.09400 0.20300 0.34600 1.00000
H8 H 8 f 0.06000 0.30000 0.40200 1.00000
O1 O 8 f 0.02580 0.43580 0.13570 1.00000
O2 O 8 f 0.42240 0.07990 0.19360 1.00000
O3 O 8 f 0.39620 0.47090 0.05210 1.00000
O4 O 8 f 0.10080 0.22690 0.00330 1.00000
O5 O 8 f 0.08480 0.09360 0.18320 1.00000
O6 O 8 f 0.22960 0.11520 0.09820 1.00000
O7 O 8 f 0.09360 0.19280 0.68270 1.00000
O8 O 8 f 0.27960 0.07460 0.40190 1.00000
O9 O 8 f 0.22140 0.35840 0.25730 1.00000
O10 O 8 f 0.09550 0.25060 0.40770 1.00000
S2 S 8 f 0.12978 0.11705 0.05632 1.00000
```

($CdSO_4$)$_3$·8H_2O ($H4_{20}$): A3B16C20D3_mC168_15_ef_8f_10f_ef - POSCAR

```
A3B16C20D3_mC168_15_ef_8f_10f_ef & a,b/a,c/a,beta,y1,y2,x3,y3,z3,x4,y4,
    ↪ z4,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10,x11
    ↪ ,y11,z11,x12,y12,z12,x13,y13,z13,x14,y14,z14,x15,y15,z15,x16,
    ↪ y16,z16,x17,y17,z17,x18,y18,z18,x19,y19,z19,x20,y20,z20,x21,y21
    ↪ ,z21,x22,y22,z22 --params=14.818,0.803279794844,0.638952625186,
    ↪ 97.39,-0.05363,0.50828,0.15443,0.40511,0.035,0.083,0.252,0.716,
    ↪ 0.095,0.209,0.601,0.251,0.079,0.406,0.271,0.075,0.309,0.195,
    ↪ 0.326,0.288,0.273,0.347,0.273,0.094,0.203,0.346,0.06,0.3,0.402,
    ↪ 0.0258,0.4358,0.1357,0.4224,0.0799,0.1936,0.3962,0.4709,0.0521,
    ↪ 0.1008,0.2269,0.0033,0.0848,0.0936,0.1832,0.2296,0.1152,0.0982,
    ↪ 0.0936,0.1928,0.6827,0.2796,0.0746,0.4019,0.2214,0.3584,0.2573,
    ↪ 0.0955,0.2506,0.4077,0.12978,0.11705,0.05632 & C2/c C_{2h}^{6}
    ↪ #15 (e^2f^20) & mC168 & $H4_{20}$ & Cd3H16O20S3 & Cd3H16O20S3 &
    ↪  R. Caminiti and G. Johansson, Acta Chem. Scand. 35a, 451-455 (
    ↪ 1981)
   1.00000000000000
   7.40900000000000  -5.95150000000000   0.00000000000000
   7.40900000000000   5.95150000000000   0.00000000000000
  -1.21779797531454   0.00000000000000   9.38935525429301
    Cd    H    O    S
     6   32   40    6
Direct
   0.05363000000000  -0.05363000000000   0.25000000000000  Cd   (4e)
  -0.05363000000000   0.05363000000000   0.75000000000000  Cd   (4e)
  -0.25068000000000   0.55954000000000   0.03500000000000  Cd   (8f)
  -0.55954000000000   0.25068000000000   0.46500000000000  Cd   (8f)
   0.25068000000000  -0.55954000000000  -0.03500000000000  Cd   (8f)
   0.55954000000000  -0.25068000000000   0.53500000000000  Cd   (8f)
  -0.16900000000000   0.33500000000000   0.71600000000000   H   (8f)
  -0.33500000000000   0.16900000000000  -0.21600000000000   H   (8f)
   0.16900000000000  -0.33500000000000  -0.71600000000000   H   (8f)
   0.33500000000000  -0.16900000000000   1.21600000000000   H   (8f)
  -0.11400000000000   0.30400000000000   0.60100000000000   H   (8f)
  -0.30400000000000   0.11400000000000  -0.10100000000000   H   (8f)
   0.11400000000000  -0.30400000000000  -0.60100000000000   H   (8f)
   0.30400000000000  -0.11400000000000   1.10100000000000   H   (8f)
   0.17200000000000   0.33000000000000   0.40600000000000   H   (8f)
  -0.33000000000000  -0.17200000000000   0.09400000000000   H   (8f)
  -0.17200000000000  -0.33000000000000  -0.40600000000000   H   (8f)
   0.33000000000000   0.17200000000000   0.90600000000000   H   (8f)
   0.19600000000000   0.34600000000000   0.30900000000000   H   (8f)
  -0.34600000000000  -0.19600000000000   0.19100000000000   H   (8f)
  -0.19600000000000  -0.34600000000000  -0.30900000000000   H   (8f)
   0.34600000000000   0.19600000000000   0.80900000000000   H   (8f)
  -0.13100000000000   0.52100000000000   0.28800000000000   H   (8f)
  -0.52100000000000   0.13100000000000   0.21200000000000   H   (8f)
   0.13100000000000  -0.52100000000000  -0.28800000000000   H   (8f)
   0.52100000000000  -0.13100000000000   0.78800000000000   H   (8f)
  -0.07400000000000   0.62000000000000   0.27300000000000   H   (8f)
  -0.62000000000000   0.07400000000000   0.22700000000000   H   (8f)
   0.07400000000000  -0.62000000000000  -0.27300000000000   H   (8f)
   0.62000000000000  -0.07400000000000   0.77300000000000   H   (8f)
```

−0.10900000000000	0.29700000000000	0.34600000000000	H	(8f)
−0.29700000000000	0.10900000000000	0.15400000000000	H	(8f)
0.10900000000000	−0.29700000000000	−0.34600000000000	H	(8f)
0.29700000000000	−0.10900000000000	0.84600000000000	H	(8f)
−0.24000000000000	0.36000000000000	0.40200000000000	H	(8f)
−0.36000000000000	0.24000000000000	0.09800000000000	H	(8f)
0.24000000000000	−0.36000000000000	−0.40200000000000	H	(8f)
0.36000000000000	−0.24000000000000	0.90200000000000	H	(8f)
−0.41000000000000	0.46160000000000	0.13570000000000	O	(8f)
−0.46160000000000	0.41000000000000	0.36430000000000	O	(8f)
0.41000000000000	−0.46160000000000	−0.13570000000000	O	(8f)
0.46160000000000	−0.41000000000000	0.63570000000000	O	(8f)
0.34250000000000	0.50230000000000	0.19360000000000	O	(8f)
−0.50230000000000	−0.34250000000000	0.30640000000000	O	(8f)
−0.34250000000000	−0.50230000000000	−0.19360000000000	O	(8f)
0.50230000000000	0.34250000000000	0.69360000000000	O	(8f)
−0.07470000000000	0.86710000000000	0.05210000000000	O	(8f)
−0.86710000000000	0.07470000000000	0.44790000000000	O	(8f)
0.07470000000000	−0.86710000000000	−0.05210000000000	O	(8f)
0.86710000000000	−0.07470000000000	0.55210000000000	O	(8f)
−0.12610000000000	0.32770000000000	0.00330000000000	O	(8f)
−0.32770000000000	0.12610000000000	0.49670000000000	O	(8f)
0.12610000000000	−0.32770000000000	−0.00330000000000	O	(8f)
0.32770000000000	−0.12610000000000	0.50330000000000	O	(8f)
−0.00880000000000	0.17840000000000	0.18320000000000	O	(8f)
−0.17840000000000	0.00880000000000	0.31680000000000	O	(8f)
0.00880000000000	−0.17840000000000	−0.18320000000000	O	(8f)
0.17840000000000	−0.00880000000000	0.68320000000000	O	(8f)
0.11440000000000	0.34480000000000	0.09820000000000	O	(8f)
−0.34480000000000	−0.11440000000000	0.40180000000000	O	(8f)
−0.11440000000000	−0.34480000000000	−0.09820000000000	O	(8f)
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−0.09920000000000	0.28640000000000	0.68270000000000	O	(8f)
−0.28640000000000	0.09920000000000	−0.18270000000000	O	(8f)
0.09920000000000	−0.28640000000000	−0.68270000000000	O	(8f)
0.28640000000000	−0.09920000000000	1.18270000000000	O	(8f)
0.20500000000000	0.35420000000000	0.40190000000000	O	(8f)
−0.35420000000000	−0.20500000000000	0.09810000000000	O	(8f)
−0.20500000000000	−0.35420000000000	−0.40190000000000	O	(8f)
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−0.13700000000000	0.57980000000000	0.25730000000000	O	(8f)
−0.57980000000000	0.13700000000000	0.24270000000000	O	(8f)
0.13700000000000	−0.57980000000000	−0.25730000000000	O	(8f)
0.57980000000000	−0.13700000000000	0.75730000000000	O	(8f)
−0.15510000000000	0.34610000000000	0.40770000000000	O	(8f)
−0.34610000000000	0.15510000000000	0.09230000000000	O	(8f)
0.15510000000000	−0.34610000000000	−0.40770000000000	O	(8f)
0.34610000000000	−0.15510000000000	0.90770000000000	O	(8f)
−0.50828000000000	0.50828000000000	0.25000000000000	S	(4e)
0.50828000000000	−0.50828000000000	0.75000000000000	S	(4e)
0.01273000000000	0.24683000000000	0.05632000000000	S	(8f)
−0.24683000000000	−0.01273000000000	0.44368000000000	S	(8f)
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Chlorite [$Mg_3(Mg_2Al)(Si_3Al)O_{10}(OH)_8$, $S5_5$]: A3B5C4D2_mC112_15_a3ef_5f_4f_2f - CIF

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_chemical_formula_sum 'Mg3 O5 (OH)4 Si2'

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_publ_author_name
 'R. C. {McMurchy}'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 88
_journal_year 1934
_journal_page_first 420
_journal_page_last 432
_publ_Section_title
;
 The Crystal Structure of the Chlorite Minerals
;

# Found in The American Mineralogist Crystal Structure Database, 2003

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      ↪ z_{9},x_{10},y_{10},z_{10},x_{11},y_{11},z_{11},x_{12},y_{12},
      ↪ z_{12},x_{13},y_{13},z_{13},x_{14},y_{14},z_{14},x_{15},y_{15},
      ↪ z_{15},x_{16},y_{16},z_{16}'
_aflow_params_values '5.305,1.73213949105,5.37229029218,97.147,0.16667,-
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      ↪ 0.039,-0.006,0.08333,0.114,-0.006,-0.417,0.114,-0.256,-0.16667,
      ↪ 0.114,-0.308,0.33333,0.039,0.142,0.0,0.211,0.142,0.33333,0.211,
      ↪ 0.142,-0.33333,0.211,-0.269,0.0,0.094,-0.269,-0.33333,0.094'
_aflow_Strukturbericht '$S5_{5}$'
_aflow_Pearson 'mC112'

_symmetry_space_group_name_H-M "C 1 2/c 1"
_symmetry_Int_Tables_number 15

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_cell_length_c    28.50000
_cell_angle_alpha 90.00000
_cell_angle_beta  97.14700
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2 -x,y,-z+1/2
3 -x,-y,-z
4 x,-y,z+1/2
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z+1/2
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z+1/2

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_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
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Mg2 Mg 4 e 0.00000 0.16667 0.25000 1.00000
Mg3 Mg 4 e 0.00000 -0.16667 0.25000 1.00000
Mg4 Mg 4 e 0.00000 0.50000 0.25000 1.00000
Mg5 Mg 8 f 0.00000 0.33333 0.00000 1.00000
O1 O 8 f -0.30800 -0.33333 0.03900 1.00000
O2 O 8 f -0.30800 0.00000 0.03900 1.00000
O3 O 8 f -0.00600 0.08333 0.11400 1.00000
O4 O 8 f -0.00600 -0.41700 0.11400 1.00000
O5 O 8 f -0.25600 -0.16667 0.11400 1.00000
OH1 OH 8 f -0.30800 0.33333 0.03900 1.00000
OH2 OH 8 f 0.14200 0.00000 0.21100 1.00000
OH3 OH 8 f 0.14200 0.33333 0.21100 1.00000
OH4 OH 8 f 0.14200 -0.33333 0.21100 1.00000
Si1 Si 8 f -0.26900 0.00000 0.09400 1.00000
Si2 Si 8 f -0.26900 -0.33333 0.09400 1.00000
```

Chlorite [$Mg_3(Mg_2Al)(Si_3Al)O_{10}(OH)_8$, $S5_5$]: A3B5C4D2_mC112_15_a3ef_5f_4f_2f - POSCAR

```
A3B5C4D2_mC112_15_a3ef_5f_4f_2f & a,b/a,c/a,beta,y2,y3,y4,x5,y5,z5,x6,y6
      ↪ ,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10,x11,y11,z11,x12,y12,
      ↪ z12,x13,y13,z13,x14,y14,z14,x15,y15,z15,x16,y16,z16 --params=
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      ↪ 0.0,0.33333,0.0,-0.308,-0.33333,0.039,-0.308,0.0,0.039,-0.006,
      ↪ 0.08333,0.114,-0.006,-0.417,0.114,-0.256,-0.16667,0.114,-0.308,
      ↪ 0.33333,0.039,0.142,0.0,0.211,0.142,0.33333,0.211,0.142,-
      ↪ 0.33333,0.211,-0.269,0.0,0.094,-0.269,-0.33333,0.094 & C2/c C_{
      ↪ 2h}^{6} #15 (ae^3f^12) & mC112 & $S5_{5}$ & Mg3O5(OH)4Si2 &
      ↪ Chlorite & R. C. McMurchy, Zeitschrift f{"u}r Kristallographie
      ↪ - Crystalline Materials 88, 420-432 (1934)
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   2.65250000000000   4.59450000000000   0.00000000000000
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     Mg     O     OH     Si
     12    20     16      8
Direct
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   0.00000000000000   0.00000000000000   0.50000000000000  Mg  (4a)
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   0.16667000000000  -0.16667000000000   0.75000000000000  Mg  (4e)
   0.16667000000000  -0.16667000000000   0.25000000000000  Mg  (4e)
  -0.16667000000000   0.16667000000000   0.75000000000000  Mg  (4e)
  -0.50000000000000   0.50000000000000   0.25000000000000  Mg  (4e)
   0.50000000000000  -0.50000000000000   0.75000000000000  Mg  (4e)
  -0.33333000000000   0.33333000000000   0.00000000000000  Mg  (8f)
  -0.33333000000000   0.33333000000000   0.50000000000000  Mg  (8f)
   0.33333000000000  -0.33333000000000   0.00000000000000  Mg  (8f)
   0.33333000000000  -0.33333000000000   0.50000000000000  Mg  (8f)
   0.02533000000000  -0.64133000000000   0.03900000000000  O  (8f)
   0.64133000000000  -0.02533000000000   0.46100000000000  O  (8f)
  -0.02533000000000   0.64133000000000  -0.03900000000000  O  (8f)
  -0.64133000000000   0.02533000000000   0.53900000000000  O  (8f)
  -0.30800000000000  -0.30800000000000   0.03900000000000  O  (8f)
   0.30800000000000   0.30800000000000   0.46100000000000  O  (8f)
   0.30800000000000   0.30800000000000  -0.03900000000000  O  (8f)
  -0.30800000000000  -0.30800000000000   0.53900000000000  O  (8f)
  -0.08933000000000   0.07733000000000   0.11400000000000  O  (8f)
  -0.07733000000000   0.08933000000000   0.38600000000000  O  (8f)
   0.08933000000000  -0.07733000000000  -0.11400000000000  O  (8f)
   0.07733000000000  -0.08933000000000   0.61400000000000  O  (8f)
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   0.42300000000000  -0.41100000000000   0.38600000000000  O  (8f)
  -0.41100000000000   0.42300000000000  -0.11400000000000  O  (8f)
  -0.42300000000000   0.41100000000000   0.61400000000000  O  (8f)
  -0.08933000000000  -0.42267000000000   0.11400000000000  O  (8f)
   0.42267000000000   0.08933000000000   0.38600000000000  O  (8f)
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  -0.42267000000000  -0.08933000000000   0.61400000000000  O  (8f)
  -0.64133000000000   0.02533000000000   0.03900000000000  OH  (8f)
  -0.02533000000000   0.64133000000000   0.46100000000000  OH  (8f)
   0.64133000000000  -0.02533000000000  -0.03900000000000  OH  (8f)
   0.02533000000000  -0.64133000000000   0.53900000000000  OH  (8f)
   0.14200000000000   0.14200000000000   0.21100000000000  OH  (8f)
  -0.14200000000000  -0.14200000000000   0.28900000000000  OH  (8f)
  -0.14200000000000  -0.14200000000000  -0.21100000000000  OH  (8f)
   0.14200000000000   0.14200000000000   0.71100000000000  OH  (8f)
  -0.19133000000000   0.47533000000000   0.21100000000000  OH  (8f)
  -0.47533000000000   0.19133000000000   0.28900000000000  OH  (8f)
   0.19133000000000  -0.47533000000000  -0.21100000000000  OH  (8f)
   0.47533000000000  -0.19133000000000   0.71100000000000  OH  (8f)
```

```
   0.47533000000000  -0.19133000000000   0.21100000000000   OH   (8f)
   0.19133000000000  -0.47533000000000   0.28900000000000   OH   (8f)
  -0.47533000000000   0.19133000000000  -0.21100000000000   OH   (8f)
  -0.19133000000000   0.47533000000000   0.71100000000000   OH   (8f)
  -0.26900000000000  -0.26900000000000   0.09400000000000   Si   (8f)
   0.26900000000000   0.26900000000000   0.40600000000000   Si   (8f)
   0.26900000000000   0.26900000000000  -0.09400000000000   Si   (8f)
  -0.26900000000000  -0.26900000000000   0.59400000000000   Si   (8f)
   0.06433000000000  -0.60233000000000   0.09400000000000   Si   (8f)
   0.60233000000000  -0.06433000000000   0.40600000000000   Si   (8f)
  -0.06433000000000   0.60233000000000  -0.09400000000000   Si   (8f)
  -0.60233000000000   0.06433000000000   0.59400000000000   Si   (8f)
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Y_2SiO_5 (RE_2SiO_5 X2): A5BC2_mC64_15_5f_f_2f - CIF

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 'A. M. Korovkin'
 'T. I. Merkulyaeva'
 'A. M. Morozova'
 'M. V. Petrov'
 'I. R. Savinova'
 'V. R. Startsev'
 'P. P. Feofilov'
_journal_name_full_name
;
 Inorganic Materials
;
_journal_volume 17
_journal_year 1981
_journal_page_first 754
_journal_page_last 758
_publ_Section_title
;
 Growth of lanthanide oxyorthosilicate single crystals, and their
      ↪ structural and optical characteristics
;

# Found in Y$_{2}$SiO$_{5}$ (Y$_{2}$[SiO$_{4}$]O ht) Crystal Structure,
      ↪ {PAULING FILE in: Inorganic Solid Phases, SpringerMaterials (
      ↪ online database), Springer, Heidelberg (ed.) SpringerMaterials
      ↪ }, 2016

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      ↪ z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},z_{8}'
_aflow_params_values '14.43,0.466597366597,0.721413721414,122.13,0.089,
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      ↪ ,0.485,0.102,0.103,0.181,0.093,0.308,0.037,0.257,0.466,0.359,
      ↪ 0.122,0.165'
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_aflow_Pearson 'mC64'

_symmetry_space_group_name_H-M "C 1 2/c 1"
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_cell_angle_gamma  90.00000

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_space_group_symop_operation_xyz
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3 -x,-y,-z
4 x,-y,z+1/2
5 x+1/2,y+1/2,z
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8 x+1/2,-y+1/2,z+1/2

loop_
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_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
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O2 O 8 f 0.11800 0.28700 0.31800 1.00000
O3 O 8 f 0.29700 0.42900 0.06000 1.00000
O4 O 8 f 0.29800 0.15700 0.33000 1.00000
O5 O 8 f 0.48500 0.10200 0.10300 1.00000
Si1 Si 8 f 0.18100 0.09300 0.30800 1.00000
Y1 Y 8 f 0.03700 0.25700 0.46600 1.00000
Y2 Y 8 f 0.35900 0.12200 0.16500 1.00000
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Y_2SiO_5 (RE_2SiO_5 X2): A5BC2_mC64_15_5f_f_2f - POSCAR

```
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      ↪ z4,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8 --params=14.43,
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      ↪ 0.287,0.318,0.297,0.429,0.06,0.298,0.157,0.33,0.485,0.102,0.103
      ↪ ,0.181,0.093,0.308,0.037,0.257,0.466,0.359,0.122,0.165 & C2/c
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      ↪ Anan'eva et al., Inorg. Mat. 17, 754-758 (1981)
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    20     4     8
Direct
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   0.72600000000000  -0.13200000000000   0.56000000000000   O   (8f)
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  -0.45500000000000  -0.14100000000000   0.17000000000000   O   (8f)
  -0.14100000000000  -0.45500000000000  -0.33000000000000   O   (8f)
   0.45500000000000   0.14100000000000   0.83000000000000   O   (8f)
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  -0.58700000000000  -0.38300000000000   0.39700000000000   O   (8f)
  -0.38300000000000  -0.58700000000000  -0.10300000000000   O   (8f)
   0.58700000000000   0.38300000000000   0.60300000000000   O   (8f)
   0.08800000000000   0.27400000000000   0.30800000000000   Si   (8f)
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  -0.08800000000000  -0.27400000000000  -0.30800000000000   Si   (8f)
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   0.29400000000000  -0.22000000000000   0.96600000000000   Y   (8f)
   0.23700000000000   0.48100000000000   0.16500000000000   Y   (8f)
  -0.48100000000000  -0.23700000000000   0.33500000000000   Y   (8f)
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α-$Zn_2V_2O_7$: A7B2C2_mC44_15_e3f_f_f - CIF

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 'R. Gopal'
 'C. Calvo'
_journal_name_full_name
;
 Canadian Journal of Chemistry
;
_journal_volume 51
_journal_year 1973
_journal_page_first 1004
_journal_page_last 1009
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;
 Crystal Structure of $\alpha$-Zn$_{2}$V$_{2}$O$_{7}$
;

# Found in $\alpha$ Cupric Divanadate, 1975

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_aflow_Pearson 'mC44'

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_cell_angle_beta   111.37000
_cell_angle_gamma  90.00000

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_space_group_symop_operation_xyz
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2 -x,y,-z+1/2
3 -x,-y,-z
4 x,-y,z+1/2
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z+1/2
7 -x+1/2,-y+1/2,-z
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8 x+1/2,-y+1/2,z+1/2

loop_
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_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
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_atom_site_occupancy
O1 O 4 e 0.00000 0.06120 0.25000 1.00000
O2 O 8 f 0.39840 -0.01890 0.36200 1.00000
O3 O 8 f 0.24400 0.15410 0.10560 1.00000
O4 O 8 f 0.15310 0.83530 0.11380 1.00000
V1 V 8 f 0.20160 0.00490 0.20580 1.00000
Zn1 Zn 8 f -0.04958 0.32390 0.51955 1.00000
```

α-$Zn_2V_2O_7$: A7B2C2_mC44_15_e3f_f_f - POSCAR

```
A7B2C2_mC44_15_e3f_f_f & a,b/a,c/a,beta,y1,x2,y2,z2,x3,y3,z3,x4,y4,z4,x5
    ↪ ,y5,z5,x6,y6,z6 --params=7.429,1.12262754072,1.35926773455,
    ↪ 111.37,0.0612,0.3984,-0.0189,0.362,0.244,0.1541,0.1056,0.1531,
    ↪ 0.8353,0.1138,0.2016,0.0049,0.2058,-0.04958,0.3239,0.51955 & C2
    ↪ /c C_{2h}^{6} #15 (ef^5) & mC44 & None & O7V2Zn2 & O7V2Zn2 & R.
    ↪  Gopal and C. Calvo, Can. J. Chem. 51, 1004-1009 (1973)
   1.00000000000000
   3.71450000000000  -4.17000000000000   0.00000000000000
   3.71450000000000   4.17000000000000   0.00000000000000
  -3.67960249157826   0.00000000000000   9.40372955289395
     O     V     Zn
    14     4     4
Direct
  -0.06120000000000   0.06120000000000   0.25000000000000   O   (4e)
   0.06120000000000  -0.06120000000000   0.75000000000000   O   (4e)
   0.41730000000000   0.37950000000000   0.36200000000000   O   (8f)
  -0.37950000000000  -0.41730000000000   0.13800000000000   O   (8f)
  -0.41730000000000  -0.37950000000000  -0.36200000000000   O   (8f)
   0.37950000000000   0.41730000000000   0.86200000000000   O   (8f)
   0.08990000000000   0.39810000000000   0.10560000000000   O   (8f)
  -0.39810000000000  -0.08990000000000   0.39440000000000   O   (8f)
  -0.08990000000000  -0.39810000000000  -0.10560000000000   O   (8f)
   0.39810000000000   0.08990000000000   0.60560000000000   O   (8f)
  -0.68220000000000   0.98840000000000   0.11380000000000   O   (8f)
  -0.98840000000000   0.68220000000000   0.38620000000000   O   (8f)
   0.68220000000000  -0.98840000000000  -0.11380000000000   O   (8f)
   0.98840000000000  -0.68220000000000   0.61380000000000   O   (8f)
   0.19670000000000   0.20650000000000   0.20580000000000   V   (8f)
  -0.20650000000000  -0.19670000000000   0.29420000000000   V   (8f)
  -0.19670000000000  -0.20650000000000  -0.20580000000000   V   (8f)
   0.20650000000000   0.19670000000000   0.70580000000000   V   (8f)
  -0.37348000000000   0.27432000000000   0.51955000000000   Zn   (8f)
  -0.27432000000000   0.37348000000000  -0.01955000000000   Zn   (8f)
   0.37348000000000  -0.27432000000000  -0.51955000000000   Zn   (8f)
   0.27432000000000  -0.37348000000000   1.01955000000000   Zn   (8f)
```

Manganese-leonite 185 K $[K_2Mn(SO_4)_2{\cdot}4H_2O]$: A8B2CD12E2_mC200_15_8f_2f_ce_2e11f_2f - CIF

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_chemical_formula_sum 'H8 K2 Mn O12 S2'

loop_
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 'B. Hertweck'
 'G. Giester'
 'E. Libowitzky'
_journal_name_full_name
;
 American Mineralogist
;
_journal_volume 86
_journal_year 2001
_journal_page_first 1282
_journal_page_last 1292
_publ_Section_title
;
 The crystal structures of the low-temperature phases of leonite-type
    ↪ compounds, K$_{2}$$Me$(SO$_{4}$)$_{2}$$\cdot$4H$_{2}$O ($Me^{2
    ↪ +}$ = Mg, Mn, Fe)
;

_aflow_title 'Manganese-leonite 185~K [K$_{2}$Mn(SO$_{4}$)$_{2}$$\
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    ↪ z_{9},x_{10},y_{10},z_{10},x_{11},y_{11},z_{11},x_{12},y_{12},
    ↪ z_{12},x_{13},y_{13},z_{13},x_{14},y_{14},z_{14},x_{15},y_{15},
    ↪ z_{15},x_{16},y_{16},z_{16},x_{17},y_{17},z_{17},x_{18},y_{18},
    ↪ z_{18},x_{19},y_{19},z_{19},x_{20},y_{20},z_{20},x_{21},y_{21},
    ↪ z_{21},x_{22},y_{22},z_{22},x_{23},y_{23},z_{23},x_{24},y_{24},
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    ↪ z_{27}'
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    ↪ 0.1105,0.0332,0.3651,0.3925,0.0298,0.4808,0.036,0.1753,0.481,
    ↪ 0.4628,0.1753,0.1379,0.2306,0.4589,0.1408,0.258,0.334,0.3794,
    ↪ 0.017,0.2991,0.372,0.4906,0.292,0.4444,0.1248,0.036,0.4419,
    ↪ 0.3781,0.0326,0.2991,0.3418,0.3983,0.2731,0.0985,0.3462,0.0301,
    ↪ 0.2498,0.1074,0.5483,0.257,0.1144,0.2827,0.2554,0.2006,0.1872,
    ↪ 0.2697,0.2437,0.3248,0.0844,0.0375,0.3231,0.417,0.0324,0.1102,
    ↪ 0.2518,0.3703,0.4779,0.2528,0.0207,0.2598,0.2429,0.2977'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mC200'

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_cell_angle_beta    117.54800
_cell_angle_gamma   90.00000

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2 -x,y,-z+1/2
3 -x,-y,-z
4 x,-y,z+1/2
5 x+1/2,y+1/2,z
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8 x+1/2,-y+1/2,z+1/2

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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
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Mn2 Mn 4 e 0.00000 0.24960 0.25000 1.00000
O1 O 4 e 0.00000 0.47800 0.25000 1.00000
O2 O 4 e 0.00000 0.02150 0.25000 1.00000
H1 H 8 f 0.19420 0.49160 0.02300 1.00000
H2 H 8 f 0.19390 0.00040 0.52050 1.00000
H3 H 8 f 0.36640 0.11050 0.03320 1.00000
H4 H 8 f 0.36510 0.39250 0.02980 1.00000
H5 H 8 f 0.48080 0.03600 0.17530 1.00000
H6 H 8 f 0.48100 0.46280 0.17530 1.00000
H7 H 8 f 0.13790 0.23060 0.45890 1.00000
H8 H 8 f 0.14080 0.25800 0.33400 1.00000
K1 K 8 f 0.37940 0.01700 0.29910 1.00000
K2 K 8 f 0.37200 0.49060 0.29200 1.00000
O3 O 8 f 0.44440 0.12480 0.03600 1.00000
O4 O 8 f 0.44190 0.37810 0.03260 1.00000
O5 O 8 f 0.29910 0.34180 0.39830 1.00000
O6 O 8 f 0.27310 0.09850 0.34620 1.00000
O7 O 8 f 0.03010 0.24980 0.10740 1.00000
O8 O 8 f 0.54830 0.25700 0.11440 1.00000
O9 O 8 f 0.28270 0.25540 0.20060 1.00000
O10 O 8 f 0.18720 0.26970 0.24370 1.00000
O11 O 8 f 0.32480 0.08440 0.03750 1.00000
O12 O 8 f 0.32310 0.41700 0.03240 1.00000
O13 O 8 f 0.11020 0.25180 0.37030 1.00000
S1 S 8 f 0.47790 0.25280 0.02070 1.00000
S2 S 8 f 0.25980 0.24290 0.29770 1.00000
```

Manganese-leonite 185 K $[K_2Mn(SO_4)_2{\cdot}4H_2O]$: A8B2CD12E2_mC200_15_8f_2f_ce_2e11f_2f - POSCAR

```
A8B2CD12E2_mC200_15_8f_2f_ce_2e11f_2f & a,b/a,c/a,beta,y2,y3,y4,x5,y5,z5
    ↪ ,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10,x11,y11,z11,
    ↪ x12,y12,z12,x13,y13,z13,x14,y14,z14,x15,y15,z15,x16,y16,z16,x17
    ↪ ,y17,z17,x18,y18,z18,x19,y19,z19,x20,y20,z20,x21,y21,z21,x22,
    ↪ y22,z22,x23,y23,z23,x24,y24,z24,x25,y25,z25,x26,y26,z26,x27,y27
    ↪ ,z27 --params=22.291,0.42837916648,0.539903997129,117.548,
    ↪ 0.2496,0.478,0.0215,0.1942,0.4916,0.023,0.1939,0.0004,0.5205,
    ↪ 0.3664,0.1105,0.0332,0.3651,0.3925,0.0298,0.4808,0.036,0.1753,
    ↪ 0.481,0.4628,0.1753,0.1379,0.2306,0.4589,0.1408,0.258,0.334,
    ↪ 0.3794,0.017,0.2991,0.372,0.4906,0.292,0.4444,0.1248,0.036,
    ↪ 0.4419,0.3781,0.0326,0.2991,0.3418,0.3983,0.2731,0.0985,0.3462,
    ↪ 0.0301,0.2498,0.1074,0.5483,0.257,0.1144,0.2827,0.2554,0.2006,
    ↪ 0.1872,0.2697,0.2437,0.3248,0.0844,0.0375,0.3231,0.417,0.0324,
    ↪ 0.1102,0.2518,0.3703,0.4779,0.2528,0.0207,0.2598,0.2429,0.2977
    ↪ & C2/c C_{2h}^{6} #15 (ce^3f^23) & mC200 & None & H8K2MnO12S2 &
    ↪  Manganese-leonite & B. Hertweck and G. Giester and E.
    ↪ Libowitzky, Am. Mineral. 86, 1282-1292 (2001)
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  11.14550000000000   4.77450000000000   0.00000000000000
  -5.56608582313949   0.00000000000000  10.67051608917980
     H     K     Mn     O     S
    32     8     4    48     8
Direct
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  -0.68580000000000   0.29740000000000   0.47700000000000   H   (8f)
   0.29740000000000  -0.68580000000000  -0.02300000000000   H   (8f)
   0.68580000000000  -0.29740000000000   0.52300000000000   H   (8f)
   0.19350000000000   0.19430000000000   0.52050000000000   H   (8f)
  -0.19430000000000  -0.19350000000000  -0.02050000000000   H   (8f)
  -0.19350000000000  -0.19430000000000  -0.52050000000000   H   (8f)
   0.19430000000000   0.19350000000000   1.02050000000000   H   (8f)
   0.25590000000000   0.47690000000000   0.03320000000000   H   (8f)
  -0.47690000000000  -0.25590000000000   0.46680000000000   H   (8f)
  -0.25590000000000  -0.47690000000000  -0.03320000000000   H   (8f)
   0.47690000000000   0.25590000000000   0.53320000000000   H   (8f)
  -0.02740000000000   0.75760000000000   0.02980000000000   H   (8f)
  -0.75760000000000   0.02740000000000   0.47020000000000   H   (8f)
```

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   0.02740000000000  −0.75760000000000  −0.02980000000000   H   (8f)
   0.75760000000000  −0.02740000000000   0.52980000000000   H   (8f)
   0.44480000000000   0.51680000000000   0.17530000000000   H   (8f)
  −0.51680000000000  −0.44480000000000   0.32470000000000   H   (8f)
  −0.44480000000000  −0.51680000000000  −0.17530000000000   H   (8f)
   0.51680000000000   0.44480000000000   0.67530000000000   H   (8f)
   0.01820000000000   0.94380000000000   0.17530000000000   H   (8f)
  −0.94380000000000  −0.01820000000000   0.32470000000000   H   (8f)
  −0.01820000000000  −0.94380000000000  −0.17530000000000   H   (8f)
   0.94380000000000   0.01820000000000   0.67530000000000   H   (8f)
  −0.09270000000000   0.36850000000000   0.45890000000000   H   (8f)
  −0.36850000000000   0.09270000000000   0.04110000000000   H   (8f)
   0.09270000000000  −0.36850000000000  −0.45890000000000   H   (8f)
   0.36850000000000  −0.09270000000000   0.95890000000000   H   (8f)
  −0.11720000000000   0.39880000000000   0.33400000000000   H   (8f)
  −0.39880000000000   0.11720000000000   0.16600000000000   H   (8f)
   0.11720000000000  −0.39880000000000  −0.33400000000000   H   (8f)
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  −0.39640000000000  −0.36240000000000   0.20090000000000   K   (8f)
  −0.36240000000000  −0.39640000000000  −0.29910000000000   K   (8f)
   0.39640000000000   0.36240000000000   0.79910000000000   K   (8f)
  −0.11860000000000   0.86260000000000   0.29200000000000   K   (8f)
  −0.86260000000000   0.11860000000000   0.20800000000000   K   (8f)
   0.11860000000000  −0.86260000000000  −0.29200000000000   K   (8f)
   0.86260000000000  −0.11860000000000   0.79200000000000   K   (8f)
   0.00000000000000   0.50000000000000   0.00000000000000   Mn  (4c)
   0.50000000000000   0.00000000000000   0.50000000000000   Mn  (4c)
  −0.24960000000000   0.24960000000000   0.25000000000000   Mn  (4e)
   0.24960000000000  −0.24960000000000   0.75000000000000   Mn  (4e)
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   0.47800000000000  −0.47800000000000   0.75000000000000   O   (4e)
  −0.02150000000000   0.02150000000000   0.25000000000000   O   (4e)
   0.02150000000000  −0.02150000000000   0.75000000000000   O   (4e)
   0.31960000000000   0.56920000000000   0.03600000000000   O   (8f)
  −0.56920000000000  −0.31960000000000   0.46400000000000   O   (8f)
  −0.31960000000000  −0.56920000000000  −0.03600000000000   O   (8f)
   0.56920000000000   0.31960000000000   0.53600000000000   O   (8f)
   0.06380000000000   0.82000000000000   0.03260000000000   O   (8f)
  −0.82000000000000  −0.06380000000000   0.46740000000000   O   (8f)
  −0.06380000000000  −0.82000000000000  −0.03260000000000   O   (8f)
   0.82000000000000   0.06380000000000   0.53260000000000   O   (8f)
  −0.04270000000000   0.64090000000000   0.39830000000000   O   (8f)
  −0.64090000000000   0.04270000000000   0.10170000000000   O   (8f)
   0.04270000000000  −0.64090000000000  −0.39830000000000   O   (8f)
   0.64090000000000  −0.04270000000000   0.89830000000000   O   (8f)
   0.17460000000000   0.37160000000000   0.34620000000000   O   (8f)
  −0.37160000000000  −0.17460000000000   0.15380000000000   O   (8f)
  −0.17460000000000  −0.37160000000000  −0.34620000000000   O   (8f)
   0.37160000000000   0.17460000000000   0.84620000000000   O   (8f)
  −0.21970000000000   0.27990000000000   0.10740000000000   O   (8f)
  −0.27990000000000   0.21970000000000   0.39260000000000   O   (8f)
   0.21970000000000  −0.27990000000000  −0.10740000000000   O   (8f)
   0.27990000000000  −0.21970000000000   0.60740000000000   O   (8f)
   0.29130000000000   0.80530000000000   0.11440000000000   O   (8f)
  −0.80530000000000  −0.29130000000000   0.38560000000000   O   (8f)
  −0.29130000000000  −0.80530000000000  −0.11440000000000   O   (8f)
   0.80530000000000   0.29130000000000   0.61440000000000   O   (8f)
   0.02730000000000   0.53810000000000   0.20060000000000   O   (8f)
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  −0.45690000000000   0.08250000000000   0.25630000000000   O   (8f)
   0.08250000000000  −0.45690000000000  −0.24370000000000   O   (8f)
   0.45690000000000  −0.08250000000000   0.74370000000000   O   (8f)
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  −0.24040000000000  −0.40920000000000  −0.03750000000000   O   (8f)
   0.40920000000000   0.24040000000000   0.53750000000000   O   (8f)
  −0.09390000000000   0.74010000000000   0.03240000000000   O   (8f)
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   0.09390000000000  −0.74010000000000  −0.03240000000000   O   (8f)
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   0.22510000000000   0.73070000000000   0.02070000000000   S   (8f)
  −0.73070000000000  −0.22510000000000   0.47930000000000   S   (8f)
  −0.22510000000000  −0.73070000000000  −0.02070000000000   S   (8f)
   0.73070000000000   0.22510000000000   0.52070000000000   S   (8f)
   0.01690000000000   0.50270000000000   0.29770000000000   S   (8f)
  −0.50270000000000  −0.01690000000000   0.20230000000000   S   (8f)
  −0.01690000000000  −0.50270000000000  −0.29770000000000   S   (8f)
   0.50270000000000   0.01690000000000   0.79770000000000   S   (8f)
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BaNi(CN)$_4$·4H$_2$O ($H4_{22}$): AB4C4D4E_mC56_15_e_2f_2f_2f_a - CIF

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 'R. G. Hazell'
 'S. E. Rasmussen'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 23
_journal_year 1969
_journal_page_first 61
_journal_page_last 69
_publ_Section_title
;
 The Crystal Structure of Barium Tetracyanonickelate(II) Tetrahydrate
;

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_aflow_proto 'AB4C4D4E_mC56_15_e_2f_2f_2f_a'
_aflow_params 'a,b/a,c/a,\beta,y_{2},x_{3},y_{3},z_{3},x_{4},y_{4},z_{4
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      ↪ y_{8},z_{8}'
_aflow_params_values '12.07,1.12758906379,0.556752278376,107.54,0.37049,
      ↪ 0.0641,0.125,0.0557,0.6475,0.4437,0.0905,0.3588,0.1896,0.3431,
      ↪ 0.4018,0.0551,0.0562,0.0997,0.2032,0.0914,0.7393,0.4076,0.1439'
_aflow_Strukturbericht '$H4_{22}$'
_aflow_Pearson 'mC56'

_symmetry_space_group_name_H-M "C 1 2/c 1"
_symmetry_Int_Tables_number 15

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_cell_length_b      13.61000
_cell_length_c      6.72000
_cell_angle_alpha  90.00000
_cell_angle_beta   107.54000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z+1/2
3 -x,-y,-z
4 x,-y,z+1/2
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z+1/2
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ni1 Ni 4 a 0.00000 0.00000 0.00000 1.00000
Ba1 Ba 4 e 0.00000 0.37049 0.25000 1.00000
C1 C 8 f 0.06410 0.12500 0.05570 1.00000
C2 C 8 f 0.64750 0.44370 0.09050 1.00000
H2O1 H2O 8 f 0.35880 0.18960 0.34310 1.00000
H2O2 H2O 8 f 0.40180 0.05510 0.05620 1.00000
N1 N 8 f 0.09970 0.20320 0.09140 1.00000
N2 N 8 f 0.73930 0.40760 0.14390 1.00000
```

BaNi(CN)$_4$·4H$_2$O ($H4_{22}$): AB4C4D4E_mC56_15_e_2f_2f_2f_a - POSCAR

```
AB4C4D4E_mC56_15_e_2f_2f_2f_a & a,b/a,c/a,beta,y2,x3,y3,z3,x4,y4,z4,x5,
      ↪ y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8 --params=12.07,1.12758906379,
      ↪ 0.556752278376,107.54,0.37049,0.0641,0.125,0.0557,0.6475,0.4437
      ↪ ,0.0905,0.3588,0.1896,0.3431,0.4018,0.0551,0.0562,0.0997,0.2032
      ↪ ,0.0914,0.7393,0.4076,0.1439 & C2/c C_{2h}^{6} #15 (aef^6) &
      ↪ mC56 & $H4_{22}$ & BaC4(H2O)4N4Ni & BaC4(H2O)4N4Ni & F. K.
      ↪ Larsen and R. G. Hazell and S. E. Rasmussen, Acta Chem. Scand.
      ↪ 23, 61-69 (1969)
   1.00000000000000
   6.03500000000000  -6.80500000000000   0.00000000000000
   6.03500000000000   6.80500000000000   0.00000000000000
  -2.02521679051048   0.00000000000000   6.40756560258531
     Ba     C    H2O     N     Ni
      2     8     8     8     2
Direct
  -0.37049000000000   0.37049000000000   0.25000000000000   Ba   (4e)
   0.37049000000000  -0.37049000000000   0.75000000000000   Ba   (4e)
  -0.06090000000000   0.18910000000000   0.05570000000000   C    (8f)
  -0.18910000000000   0.06090000000000   0.44430000000000   C    (8f)
   0.06090000000000  -0.18910000000000  -0.05570000000000   C    (8f)
   0.18910000000000  -0.06090000000000   0.55570000000000   C    (8f)
   0.20380000000000   1.09120000000000   0.09050000000000   C    (8f)
  -1.09120000000000  -0.20380000000000   0.40950000000000   C    (8f)
  -0.20380000000000  -1.09120000000000  -0.09050000000000   C    (8f)
   1.09120000000000   0.20380000000000   0.59050000000000   C    (8f)
   0.16920000000000   0.54840000000000   0.34310000000000   H2O  (8f)
  -0.54840000000000  -0.16920000000000   0.15690000000000   H2O  (8f)
  -0.16920000000000  -0.54840000000000  -0.34310000000000   H2O  (8f)
   0.54840000000000   0.16920000000000   0.84310000000000   H2O  (8f)
   0.34670000000000   0.45690000000000   0.05620000000000   H2O  (8f)
  -0.45690000000000  -0.34670000000000   0.44380000000000   H2O  (8f)
  -0.34670000000000  -0.45690000000000  -0.05620000000000   H2O  (8f)
   0.45690000000000   0.34670000000000   0.55620000000000   H2O  (8f)
  -0.10350000000000   0.30290000000000   0.09140000000000   N    (8f)
  -0.30290000000000   0.10350000000000   0.40860000000000   N    (8f)
   0.10350000000000  -0.30290000000000  -0.09140000000000   N    (8f)
   0.30290000000000  -0.10350000000000   0.59140000000000   N    (8f)
   0.33170000000000   1.14690000000000   0.14390000000000   N    (8f)
  -1.14690000000000  -0.33170000000000   0.35610000000000   N    (8f)
  -0.33170000000000  -1.14690000000000  -0.14390000000000   N    (8f)
   1.14690000000000   0.33170000000000   0.64390000000000   N    (8f)
   0.00000000000000   0.00000000000000   0.00000000000000   Ni   (4a)
   0.00000000000000   0.00000000000000   0.50000000000000   Ni   (4a)
```

Gypsum ($CaSO_4{\cdot}2H_2O$, $H4_6$): AB4C6D_mC48_15_e_2f_3f_e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Gypsum'
_chemical_formula_sum 'Ca H4 O6 S'

loop_
_publ_author_name
 'P. Comodi'
 'S. Nazzareni'
 'P. F. Zanazzi'
 'S. Speziale'
_journal_name_full_name
;
 American Mineralogist
;
_journal_volume 93
_journal_year 2008
_journal_page_first 1530
_journal_page_last 1537
_publ_Section_title
;
 High-pressure behavior of gypsum: A single-crystal X-ray study
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'Gypsum (CaSO$_{4}$\cdot2H$_{2}$O, $H4_{6}$) Structure'
_aflow_proto 'AB4C6D_mC48_15_e_2f_3f_e'
_aflow_params 'a,b/a,c/a,\beta,y_{1},y_{2},x_{3},y_{3},z_{3},x_{4},y_{4
      ↪ },z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{6},x_{7},y_{7},z_{7}'
_aflow_params_values '6.277,2.41851202804,0.90361637725,114.11,0.1705,
      ↪ 0.67273,0.742,0.087,0.766,0.756,0.02,-0.077,0.08319,0.27218,
      ↪ 0.59103,0.19997,0.38195,-0.08702,0.79177,0.06826,-0.07831'
_aflow_Strukturbericht '$H4_{6}$'
_aflow_Pearson 'mC48'

_symmetry_space_group_name_H-M "C 1 2/c 1"
_symmetry_Int_Tables_number 15

_cell_length_a    6.27700
_cell_length_b    15.18100
_cell_length_c    5.67200
_cell_angle_alpha 90.00000
_cell_angle_beta  114.11000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z+1/2
3 -x,-y,-z
4 x,-y,z+1/2
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z+1/2
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ca1 Ca 4 e 0.00000 0.17050 0.25000 1.00000
S1 S 4 e 0.00000 0.67273 0.25000 1.00000
H1 H 8 f 0.74200 0.08700 0.76600 1.00000
H2 H 8 f 0.75600 0.02000 -0.07700 1.00000
O1 O 8 f 0.08319 0.27218 0.59103 1.00000
O2 O 8 f 0.19997 0.38195 -0.08702 1.00000
O3 O 8 f 0.79177 0.06826 -0.07831 1.00000
```

Gypsum ($CaSO_4{\cdot}2H_2O$, $H4_6$): AB4C6D_mC48_15_e_2f_3f_e - POSCAR

```
AB4C6D_mC48_15_e_2f_3f_e & a,b/a,c/a,beta,y1,y2,x3,y3,z3,x4,y4,z4,x5,y5,
      ↪ z5,x6,y6,z6,x7,y7,z7 --params=6.277,2.41851202804,0.90361637725
      ↪ ,114.11,0.1705,0.67273,0.742,0.087,0.766,0.756,0.02,-0.077,
      ↪ 0.08319,0.27218,0.59103,0.19997,0.38195,-0.08702,0.79177,
      ↪ 0.06826,-0.07831 & C2/c C_{2h}^{6} #15 (e^2f^5) & mC48 & $H4_{6
      ↪ }$ & CaH4O6S & Gypsum & P. Comodi et al., Am. Mineral. 93,
      ↪ 1530-1537 (2008)
   1.00000000000000
   3.13850000000000  -7.59050000000000   0.00000000000000
   3.13850000000000   7.59050000000000   0.00000000000000
  -2.31695399688312   0.00000000000000   5.17719114736237
     Ca     H     O     S
      2     8    12     2
Direct
  -0.17050000000000   0.17050000000000   0.25000000000000   Ca   (4e)
   0.17050000000000  -0.17050000000000   0.75000000000000   Ca   (4e)
   0.65550000000000   0.82900000000000   0.76600000000000   H    (8f)
  -0.82900000000000  -0.65550000000000  -0.26600000000000   H    (8f)
  -0.65550000000000  -0.82900000000000  -0.76600000000000   H    (8f)
   0.82900000000000   0.65550000000000   1.26600000000000   H    (8f)
   0.73600000000000   0.77600000000000  -0.07700000000000   H    (8f)
  -0.77600000000000  -0.73600000000000   0.57700000000000   H    (8f)
  -0.73600000000000  -0.77600000000000   0.07700000000000   H    (8f)
   0.77600000000000   0.73600000000000   0.42300000000000   H    (8f)
  -0.18899000000000   0.35537000000000   0.59103000000000   O    (8f)
  -0.35537000000000   0.18899000000000  -0.09103000000000   O    (8f)
   0.18899000000000  -0.35537000000000  -0.59103000000000   O    (8f)
   0.35537000000000  -0.18899000000000   1.09103000000000   O    (8f)
  -0.18198000000000   0.58192000000000  -0.08702000000000   O    (8f)
  -0.58192000000000   0.18198000000000   0.58702000000000   O    (8f)
   0.18198000000000  -0.58192000000000   0.08702000000000   O    (8f)
   0.58192000000000  -0.18198000000000   0.41298000000000   O    (8f)
   0.72351000000000   0.86003000000000  -0.07831000000000   O    (8f)
  -0.86003000000000  -0.72351000000000   0.57831000000000   O    (8f)
  -0.72351000000000  -0.86003000000000   0.07831000000000   O    (8f)
   0.86003000000000   0.72351000000000   0.42169000000000   O    (8f)
  -0.67273000000000   0.67273000000000   0.25000000000000   S    (4e)
   0.67273000000000  -0.67273000000000   0.75000000000000   S    (4e)
```

Ta_2NiSe_5: AB5C2_mC32_15_e_e2f_f - CIF

```
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_audit_creation_method FINDSYM

_chemical_name_mineral 'NiSe5Ta2'
_chemical_formula_sum 'Ni Se5 Ta2'

loop_
_publ_author_name
 'S. A. Sunshine'
 'J. A. Ibers'
_journal_name_full_name
;
 Inorganic Chemistry
;
_journal_volume 24
_journal_year 1985
_journal_page_first 3611
_journal_page_last 3614
_publ_Section_title
;
 Structure and physical properties of the new layered ternary
      ↪ chalcogenides tantalum nickel sulfide (Ta$_{2}$NiS$_{5}$) and
      ↪ tantalum nickel selenide (Ta$_{2}$NiSe$_{5}$)
;

# Found in Physical and structural properties of the new layered
      ↪ compounds Ta$_{2}$NiS$_{5}$ and Ta$_{2}$NiSe$_{5}$, 1986

_aflow_title 'Ta$_{2}$NiSe$_{5}$ Structure'
_aflow_proto 'AB5C2_mC32_15_e_e2f_f'
_aflow_params 'a,b/a,c/a,\beta,y_{1},y_{2},x_{3},y_{3},z_{3},x_{4},y_{4
      ↪ },z_{4},x_{5},y_{5},z_{5}'
_aflow_params_values '3.496,3.66962242563,4.47397025172,90.53,0.70113,
      ↪ 0.32714,0.5053,0.08039,0.13798,-0.00513,0.14565,-0.04913,-
      ↪ 0.00793,0.22135,0.11044'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mC32'

_symmetry_space_group_name_H-M "C 1 2/c 1"
_symmetry_Int_Tables_number 15

_cell_length_a    3.49600
_cell_length_b    12.82900
_cell_length_c    15.64100
_cell_angle_alpha 90.00000
_cell_angle_beta  90.53000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
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2 -x,y,-z+1/2
3 -x,-y,-z
4 x,-y,z+1/2
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z+1/2
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ni1 Ni 4 e 0.00000 0.70113 0.25000 1.00000
Se1 Se 4 e 0.00000 0.32714 0.25000 1.00000
Se2 Se 8 f 0.50530 0.08039 0.13798 1.00000
Se3 Se 8 f -0.00513 0.14565 -0.04913 1.00000
Ta1 Ta 8 f -0.00793 0.22135 0.11044 1.00000
```

Ta_2NiSe_5: AB5C2_mC32_15_e_e2f_f - POSCAR

```
AB5C2_mC32_15_e_e2f_f & a,b/a,c/a,beta,y1,y2,x3,y3,z3,x4,y4,z4,x5,y5,z5
      ↪ --params=3.496,3.66962242563,4.47397025172,90.53,0.70113,
      ↪ 0.32714,0.5053,0.08039,0.13798,-0.00513,0.14565,-0.04913,-
      ↪ 0.00793,0.22135,0.11044 & C2/c C_{2h}^{6} #15 (e^2f^3) & mC32 &
      ↪  None & NiSe5Ta2 & NiSe5Ta2 & S. A. Sunshine and J. A. Ibers,
      ↪ Inorg. Chem. 24, 3611-3614 (1985)
   1.00000000000000
   1.74800000000000  -6.41450000000000   0.00000000000000
   1.74800000000000   6.41450000000000   0.00000000000000
  -0.14468101925982   0.00000000000000  15.64033082778830
```

```
    Ni    Se    Ta
     2    10     4
Direct
  -0.70113000000000   0.70113000000000   0.25000000000000   Ni   (4e)
   0.70113000000000  -0.70113000000000   0.75000000000000   Ni   (4e)
  -0.32714000000000   0.32714000000000   0.25000000000000   Se   (4e)
   0.32714000000000  -0.32714000000000   0.75000000000000   Se   (4e)
   0.42491000000000   0.58569000000000   0.13798000000000   Se   (8f)
  -0.58569000000000  -0.42491000000000   0.36202000000000   Se   (8f)
  -0.42491000000000  -0.58569000000000  -0.13798000000000   Se   (8f)
   0.58569000000000   0.42491000000000   0.63798000000000   Se   (8f)
  -0.15078000000000   0.14052000000000  -0.04913000000000   Se   (8f)
  -0.14052000000000   0.15078000000000   0.54913000000000   Se   (8f)
   0.15078000000000  -0.14052000000000   0.04913000000000   Se   (8f)
   0.14052000000000  -0.15078000000000   0.45087000000000   Se   (8f)
  -0.22928000000000   0.21342000000000   0.11044000000000   Ta   (8f)
  -0.21342000000000   0.22928000000000   0.38956000000000   Ta   (8f)
   0.22928000000000  -0.21342000000000  -0.11044000000000   Ta   (8f)
   0.21342000000000  -0.22928000000000   0.61044000000000   Ta   (8f)
```

Pyrophyllite [$AlSi_2O_5(OH)$, $S5_6$]: AB5CD2_mC72_15_f_5f_f_2f - CIF

```
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_chemical_name_mineral 'Pyrophyllite'
_chemical_formula_sum 'Al O5 (OH) Si2'

loop_
_publ_author_name
 'J. W. Gruner'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 88
_journal_year 1934
_journal_page_first 412
_journal_page_last 419
_publ_Section_title
;
 The Crystal Structures of Talc and Pyrophyllite
;

# Found in Strukturbericht Band III 1933-1935, 1937

_aflow_title 'Pyrophyllite [AlSi$_{2}$O$_{5}$(OH), $S5_{6}$] Structure'
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_aflow_params 'a,b/a,c/a,\beta,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3
      ↪ },y_{3},z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},
      ↪ z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9}'
_aflow_params_values '5.14,1.73151750973,3.60894941634,99.91667,0.0,
      ↪ 0.33333,0.0,0.20278,0.5,0.05833,0.20278,0.16667,0.05833,0.025,
      ↪ 0.08333,0.17639,0.525,0.08333,0.17639,0.275,0.33333,0.17639,
      ↪ 0.20278,0.83333,0.05833,0.76111,0.0,0.14306,0.26111,0.16667,
      ↪ 0.14306'
_aflow_Strukturbericht '$S5_{6}$'
_aflow_Pearson 'mC72'

_symmetry_space_group_name_H-M "C 1 2/c 1"
_symmetry_Int_Tables_number 15

_cell_length_a    5.14000
_cell_length_b    8.90000
_cell_length_c    18.55000
_cell_angle_alpha 90.00000
_cell_angle_beta  99.91667
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
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2 -x,y,-z+1/2
3 -x,-y,-z
4 x,-y,z+1/2
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z+1/2
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
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O1 O 8 f 0.20278 0.50000 0.05833 1.00000
O2 O 8 f 0.20278 0.16667 0.05833 1.00000
O3 O 8 f 0.02500 0.08333 0.17639 1.00000
O4 O 8 f 0.52500 0.08333 0.17639 1.00000
O5 O 8 f 0.27500 0.33333 0.17639 1.00000
OH1 OH 8 f 0.20278 0.83333 0.05833 1.00000
Si1 Si 8 f 0.76111 0.00000 0.14306 1.00000
Si2 Si 8 f 0.26111 0.16667 0.14306 1.00000
```

Pyrophyllite [$AlSi_2O_5(OH)$, $S5_6$]: AB5CD2_mC72_15_f_5f_f_2f - POSCAR

```
AB5CD2_mC72_15_f_5f_f_2f & a,b/a,c/a,beta,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,
      ↪ y4,z4,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9 --params=
      ↪ 5.14,1.73151750973,3.60894941634,99.91667,0.0,0.33333,0.0,
      ↪ 0.20278,0.5,0.05833,0.20278,0.16667,0.05833,0.025,0.08333,
      ↪ 0.17639,0.525,0.08333,0.17639,0.275,0.33333,0.17639,0.20278,
      ↪ 0.83333,0.05833,0.76111,0.0,0.14306,0.26111,0.16667,0.14306 &
      ↪ C2/c C_{2h}^{6} #15 (f^9) & mC72 & $S5_{6}$ & AlO5(OH)Si2 &
      ↪ Pyrophyllite & J. W. Gruner, Zeitschrift f{"u}r
      ↪ Kristallographie - Crystalline Materials 88, 412-419 (1934)
  1.00000000000000
  2.57000000000000  -4.45000000000000   0.00000000000000
  2.57000000000000   4.45000000000000   0.00000000000000
 -3.19460136479521   0.00000000000000  18.27284931585790
    Al    O   OH   Si
     4   20    4    8
Direct
  -0.33333000000000   0.33333000000000   0.00000000000000   Al   (8f)
  -0.33333000000000   0.33333000000000   0.50000000000000   Al   (8f)
   0.33333000000000  -0.33333000000000   0.00000000000000   Al   (8f)
   0.33333000000000  -0.33333000000000   0.50000000000000   Al   (8f)
  -0.29722000000000   0.70278000000000   0.05833000000000   O    (8f)
  -0.70278000000000   0.29722000000000   0.44167000000000   O    (8f)
   0.29722000000000  -0.70278000000000  -0.05833000000000   O    (8f)
   0.70278000000000  -0.29722000000000   0.55833000000000   O    (8f)
   0.03611000000000   0.36945000000000   0.05833000000000   O    (8f)
  -0.36945000000000  -0.03611000000000   0.44167000000000   O    (8f)
  -0.03611000000000  -0.36945000000000  -0.05833000000000   O    (8f)
   0.36945000000000   0.03611000000000   0.55833000000000   O    (8f)
  -0.05833000000000   0.10833000000000   0.17639000000000   O    (8f)
  -0.10833000000000   0.05833000000000   0.32361000000000   O    (8f)
   0.05833000000000  -0.10833000000000  -0.17639000000000   O    (8f)
   0.10833000000000  -0.05833000000000   0.67639000000000   O    (8f)
   0.44167000000000   0.60833000000000   0.17639000000000   O    (8f)
  -0.60833000000000  -0.44167000000000   0.32361000000000   O    (8f)
  -0.44167000000000  -0.60833000000000  -0.17639000000000   O    (8f)
   0.60833000000000   0.44167000000000   0.67639000000000   O    (8f)
  -0.05833000000000   0.60833000000000   0.17639000000000   O    (8f)
  -0.60833000000000   0.05833000000000   0.32361000000000   O    (8f)
   0.05833000000000  -0.60833000000000  -0.17639000000000   O    (8f)
   0.60833000000000  -0.05833000000000   0.67639000000000   O    (8f)
  -0.63055000000000   1.03611000000000   0.05833000000000   OH   (8f)
  -1.03611000000000   0.63055000000000   0.44167000000000   OH   (8f)
   0.63055000000000  -1.03611000000000  -0.05833000000000   OH   (8f)
   1.03611000000000  -0.63055000000000   0.55833000000000   OH   (8f)
   0.76111000000000   0.76111000000000   0.14306000000000   Si   (8f)
  -0.76111000000000  -0.76111000000000   0.35694000000000   Si   (8f)
  -0.76111000000000  -0.76111000000000  -0.14306000000000   Si   (8f)
   0.76111000000000   0.76111000000000   0.64306000000000   Si   (8f)
   0.09444000000000   0.42778000000000   0.14306000000000   Si   (8f)
  -0.42778000000000  -0.09444000000000   0.35694000000000   Si   (8f)
  -0.09444000000000  -0.42778000000000  -0.14306000000000   Si   (8f)
   0.42778000000000   0.09444000000000   0.64306000000000   Si   (8f)
```

Titanite ($CaTiSiO_5$, $S0_6$): AB5CD_mC32_15_e_e2f_e_b - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Titanite'
_chemical_formula_sum 'Ca O5 Si Ti'

loop_
_publ_author_name
 'F. C. Hawthorne'
 'L. A. Groat'
 'M. Raudsepp'
 'N. A. Ball'
 'M. Kimata'
 'F. D. Spike'
 'R. Gaba'
 'N. M. Halden'
 'G. R. Lumpkin'
 'R. C. Ewing'
 'R. B. Greegor'
 'F. W. Lytle'
 'T. {Scott Ercit}'
 'G. R. Rossman'
 'F. J. Wicks'
 'R. A. Ramik'
 'B. L. Sherriff'
 'M. E. Fleet'
 'C. {McCammon}'
_journal_name_full_name
;
 American Mineralogist
;
_journal_volume 76
_journal_year 1991
_journal_page_first 370
_journal_page_last 396
_publ_Section_title
;
 Alpha-decay damage in titanite
;

_aflow_title 'Titanite (CaTiSiO$_{5}$, $S0_{6}$) Structure'
_aflow_proto 'AB5CD_mC32_15_e_e2f_e_b'
_aflow_params 'a,b/a,c/a,\beta,y_{2},y_{3},y_{4},x_{5},y_{5},z_{5},x_{6
      ↪ },y_{6},z_{6}'
_aflow_params_values '6.549,1.32768361582,1.07802717972,113.87,0.8323,
      ↪ 0.5714,0.1828,0.1855,0.0663,0.4102,0.1025,0.2893,0.1185'
_aflow_Strukturbericht '$S0_{6}$'
_aflow_Pearson 'mC32'

_symmetry_space_group_name_H-M "C 1 2/c 1"
_symmetry_Int_Tables_number 15
```

```
_cell_length_a      6.54900
_cell_length_b      8.69500
_cell_length_c      7.06000
_cell_angle_alpha  90.00000
_cell_angle_beta   113.87000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z+1/2
3 -x,-y,-z
4 x,-y,z+1/2
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z+1/2
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ti1 Ti 4 b 0.00000 0.50000 0.00000 1.00000
Ca1 Ca 4 e 0.00000 0.83230 0.25000 1.00000
O1 O 4 e 0.00000 0.57140 0.25000 1.00000
Si1 Si 4 e 0.00000 0.18280 0.25000 1.00000
O2 O 8 f 0.18550 0.06630 0.41020 1.00000
O3 O 8 f 0.10250 0.28930 0.11850 1.00000
```

Titanite ($CaTiSiO_5$, $S0_6$): AB5CD_mC32_15_e_e2f_e_b - POSCAR

```
AB5CD_mC32_15_e_e2f_e_b & a,b/a,c/a,beta,y2,y3,y4,x5,y5,z5,x6,y6,z6 --
    ↪ params=6.549,1.32768361582,1.07802717972,113.87,0.8323,0.5714,
    ↪ 0.1828,0.1855,0.0663,0.4102,0.1025,0.2893,0.1185 & C2/c C_{2h
    ↪ }^{6} #15 (be^3f^2) & mC32 & $S0_{6}$ & CaO5SiTi & Titanite &
    ↪ F. C. Hawthorne et al., Am. Mineral. 76, 370-396 (1991)
   1.00000000000000
   3.27450000000000  -4.34750000000000   0.00000000000000
   3.27450000000000   4.34750000000000   0.00000000000000
  -2.85691957284066   0.00000000000000   6.45612968846816
     Ca      O      Si      Ti
      2     10       2       2
Direct
  -0.83230000000000   0.83230000000000   0.25000000000000   Ca    (4e)
   0.83230000000000  -0.83230000000000   0.75000000000000   Ca    (4e)
  -0.57140000000000   0.57140000000000   0.25000000000000   O     (4e)
   0.57140000000000  -0.57140000000000   0.75000000000000   O     (4e)
   0.11920000000000   0.25180000000000   0.41020000000000   O     (8f)
  -0.25180000000000  -0.11920000000000   0.08980000000000   O     (8f)
  -0.11920000000000  -0.25180000000000  -0.41020000000000   O     (8f)
   0.25180000000000   0.11920000000000   0.91020000000000   O     (8f)
  -0.18680000000000   0.39180000000000   0.11850000000000   O     (8f)
  -0.39180000000000   0.18680000000000   0.38150000000000   O     (8f)
   0.18680000000000  -0.39180000000000  -0.11850000000000   O     (8f)
   0.39180000000000  -0.18680000000000   0.61850000000000   O     (8f)
  -0.18280000000000   0.18280000000000   0.25000000000000   Si    (4e)
   0.18280000000000  -0.18280000000000   0.75000000000000   Si    (4e)
   0.50000000000000   0.50000000000000   0.00000000000000   Ti    (4b)
   0.50000000000000   0.50000000000000   0.50000000000000   Ti    (4b)
```

$KFeS_2$ ($F5_a$): ABC2_mC16_15_e_e_f - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'FeKS2'
_chemical_formula_sum 'Fe K S2'

loop_
_publ_author_name
 'W. Bronger'
 'A. Kyas'
 'P. M\"{u}ller'
_journal_name_full_name
;
 Journal of Solid State Chemistry
;
_journal_volume 70
_journal_year 1987
_journal_page_first 262
_journal_page_last 270
_publ_Section_title
;
 The antiferromagnetic structures of KFeS$_{2}$, RbFeS$_{2}$, KFeSe$_{2}
    ↪ $, and RbFeSe$_{2}$ and the correlation between magnetic
    ↪ moments and crystal field calculations
;

_aflow_title 'KFeS$_{2}$ ($F5_{a}$) Structure'
_aflow_proto 'ABC2_mC16_15_e_e_f'
_aflow_params 'a,b/a,c/a,\beta,y_{1},y_{2},x_{3},y_{3},z_{3}'
_aflow_params_values '7.084,1.595567476,0.761434217956,113.2,-0.00332,
    ↪ 0.3572,0.196,0.1098,0.1068'
_aflow_Strukturbericht '$F5_{a}$'
_aflow_Pearson 'mC16'

_symmetry_space_group_name_H-M "C 1 2/c 1"
_symmetry_Int_Tables_number 15

_cell_length_a      7.08400
_cell_length_b     11.30300
_cell_length_c      5.39400
_cell_angle_alpha  90.00000
_cell_angle_beta   113.20000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z+1/2
3 -x,-y,-z
4 x,-y,z+1/2
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z+1/2
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Fe1 Fe 4 e 0.00000 -0.00332 0.25000 1.00000
K1 K 4 e 0.00000 0.35720 0.25000 1.00000
S1 S 8 f 0.19600 0.10980 0.10680 1.00000
```

$KFeS_2$ ($F5_a$): ABC2_mC16_15_e_e_f - POSCAR

```
ABC2_mC16_15_e_e_f & a,b/a,c/a,beta,y1,y2,x3,y3,z3 --params=7.084,
    ↪ 1.595567476,0.761434217956,113.2,-0.00332,0.3572,0.196,0.1098,
    ↪ 0.1068 & C2/c C_{2h}^{6} #15 (e^2f) & mC16 & $F5_{a}$ & FeKS2 &
    ↪  FeKS2 & W. Bronger and A. Kyas and P. M\"{u}ller, J. Solid
    ↪ State Chem. 70, 262-270 (1987)
   1.00000000000000
   3.54200000000000  -5.65150000000000   0.00000000000000
   3.54200000000000   5.65150000000000   0.00000000000000
  -2.12492266033359   0.00000000000000   4.95781601994273
     Fe      K      S
      2      2      4
Direct
   0.00332000000000  -0.00332000000000   0.25000000000000   Fe    (4e)
  -0.00332000000000   0.00332000000000   0.75000000000000   Fe    (4e)
  -0.35720000000000   0.35720000000000   0.25000000000000   K     (4e)
   0.35720000000000  -0.35720000000000   0.75000000000000   K     (4e)
   0.08620000000000   0.30580000000000   0.10680000000000   S     (8f)
  -0.30580000000000  -0.08620000000000   0.39320000000000   S     (8f)
  -0.08620000000000  -0.30580000000000  -0.10680000000000   S     (8f)
   0.30580000000000   0.08620000000000   0.60680000000000   S     (8f)
```

Diopside [$CaMg(SiO_3)_2$, $S4_1$]: ABC6D2_mC40_15_e_e_3f_f - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Diopside'
_chemical_formula_sum 'Ca Mg O6 Si2'

loop_
_publ_author_name
 'L. W. Finger'
 'Y. Ohashi'
_journal_name_full_name
;
 American Mineralogist
;
_journal_volume 61
_journal_year 1976
_journal_page_first 303
_journal_page_last 310
_publ_Section_title
;
 The thermal expansion of diopside to 800$^{\circ}$C and a refinement of
    ↪  the crystal structure at 700$^{\circ}$C
;

_aflow_title 'Diopside [CaMg(SiO$_{3}$)$_{2}$, $S4_{1}$] Structure'
_aflow_proto 'ABC6D2_mC40_15_e_e_3f_f'
_aflow_params 'a,b/a,c/a,\beta,y_{1},y_{2},x_{3},y_{3},z_{3},x_{4},y_{4
    ↪ },z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{6}'
_aflow_params_values '9.804,0.921052631579,0.538249694002,106.02,-0.0933
    ↪ ,0.3003,0.1167,0.0872,0.1417,0.3617,0.246,0.3166,0.3493,0.0155
    ↪ ,-0.0022,0.2864,0.0923,0.2299'
_aflow_Strukturbericht '$S4_{1}$'
_aflow_Pearson 'mC40'

_symmetry_space_group_name_H-M "C 1 2/c 1"
_symmetry_Int_Tables_number 15

_cell_length_a      9.80400
_cell_length_b      9.03000
_cell_length_c      5.27700
_cell_angle_alpha  90.00000
_cell_angle_beta   106.02000
_cell_angle_gamma  90.00000

loop_
```

```
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z+1/2
3 -x,-y,-z
4 x,-y,z+1/2
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z+1/2
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ca1 Ca 4 e 0.00000 -0.09330 0.25000 1.00000
Mg1 Mg 4 e 0.00000 0.30030 0.25000 1.00000
O1 O 8 f 0.11670 0.08720 0.14170 1.00000
O2 O 8 f 0.36170 0.24600 0.31660 1.00000
O3 O 8 f 0.34930 0.01550 -0.00220 1.00000
Si1 Si 8 f 0.28640 0.09230 0.22990 1.00000
```

Diopside [$CaMg(SiO_3)_2$, $S4_1$]: ABC6D2_mC40_15_e_e_3f_f - POSCAR

```
ABC6D2_mC40_15_e_e_3f_f & a,b/a,c/a,beta,y1,y2,x3,y3,z3,x4,y4,z4,x5,y5,
    ↪ z5,x6,y6,z6 --params=9.804,0.921052631579,0.538249694002,106.02
    ↪ ,-0.0933,0.3003,0.1167,0.0872,0.1417,0.3617,0.246,0.3166,0.3493
    ↪ ,0.0155,-0.0022,0.2864,0.0923,0.2299 & C2/c C_{2h}^{6} #15 (e^
    ↪ 2f^4) & mC40 & $S4_{1}$ & CaMgO6Si2 & Diopside & L. W. Finger
    ↪ and Y. Ohashi, Am. Mineral. 61, 303-310 (1976)
   1.00000000000000
   4.90200000000000  -4.51500000000000   0.00000000000000
   4.90200000000000   4.51500000000000   0.00000000000000
  -1.45630890173743   0.00000000000000   5.07206993077977
     Ca     Mg      O     Si
      2      2     12      4
Direct
   0.09330000000000  -0.09330000000000   0.25000000000000   Ca    (4e)
  -0.09330000000000   0.09330000000000   0.75000000000000   Ca    (4e)
  -0.30030000000000   0.30030000000000   0.25000000000000   Mg    (4e)
   0.30030000000000  -0.30030000000000   0.75000000000000   Mg    (4e)
   0.02950000000000   0.20390000000000   0.14170000000000   O     (8f)
  -0.20390000000000  -0.02950000000000   0.35830000000000   O     (8f)
  -0.02950000000000  -0.20390000000000  -0.14170000000000   O     (8f)
   0.20390000000000   0.02950000000000   0.64170000000000   O     (8f)
   0.11570000000000   0.60770000000000   0.31660000000000   O     (8f)
  -0.60770000000000  -0.11570000000000   0.18340000000000   O     (8f)
  -0.11570000000000  -0.60770000000000  -0.31660000000000   O     (8f)
   0.60770000000000   0.11570000000000   0.81660000000000   O     (8f)
   0.33380000000000   0.36480000000000  -0.00220000000000   O     (8f)
  -0.36480000000000  -0.33380000000000   0.50220000000000   O     (8f)
  -0.33380000000000  -0.36480000000000   0.00220000000000   O     (8f)
   0.36480000000000   0.33380000000000   0.49780000000000   O     (8f)
   0.19410000000000   0.37870000000000   0.22990000000000   Si    (8f)
  -0.37870000000000  -0.19410000000000   0.27010000000000   Si    (8f)
  -0.19410000000000  -0.37870000000000  -0.22990000000000   Si    (8f)
   0.37870000000000   0.19410000000000   0.72990000000000   Si    (8f)
```

β-Ga (*obsolete*): A_mC4_15_e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Ga'
_chemical_formula_sum 'Ga'

loop_
_publ_author_name
 'L. Bosio'
 'A. Defrain'
 'H. Curien'
 'A. Rimsky'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 25
_journal_year 1969
_journal_page_first 995
_journal_page_last 995
_publ_Section_title
;
 Structure cristalline du gallium $\beta$
;

# Found in The Structures of the Elements, {Reprint of the 1974 John
    ↪ Wilely \& Sons edition}, 1982

_aflow_title '$\beta$-Ga ({\em{obsolete}}) Structure'
_aflow_proto 'A_mC4_15_e'
_aflow_params 'a,b/a,c/a,\beta,y_{1}'
_aflow_params_values '2.766,2.91142443962,1.20462762111,92.03333,0.131'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'mC4'

_symmetry_space_group_name_H-M "C 1 2/c 1"
_symmetry_Int_Tables_number 15

_cell_length_a      2.76600
_cell_length_b      8.05300
_cell_length_c      3.33200
_cell_angle_alpha   90.00000
_cell_angle_beta    92.03333
_cell_angle_gamma   90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,y,-z+1/2
3 -x,-y,-z
4 x,-y,z+1/2
5 x+1/2,y+1/2,z
6 -x+1/2,y+1/2,-z+1/2
7 -x+1/2,-y+1/2,-z
8 x+1/2,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ga1 Ga 4 e 0.00000 0.13100 0.25000 1.00000
```

β-Ga (*obsolete*): A_mC4_15_e - POSCAR

```
A_mC4_15_e & a,b/a,c/a,beta,y1 --params=2.766,2.91142443962,
    ↪ 1.20462762111,92.03333,0.131 & C2/c C_{2h}^{6} #15 (e) & mC4 &
    ↪ None & Ga & Ga & L. Bosio et al., Acta Crystallogr. Sect. B
    ↪ Struct. Sci. 25, 995(1969)
   1.00000000000000
   1.38300000000000  -4.02650000000000   0.00000000000000
   1.38300000000000   4.02650000000000   0.00000000000000
  -0.11822220765196   0.00000000000000   3.32990202702991
     Ga
      2
Direct
  -0.13100000000000   0.13100000000000   0.25000000000000   Ga    (4e)
   0.13100000000000  -0.13100000000000   0.75000000000000   Ga    (4e)
```

$NaNbO_3$: ABC3_oP40_17_abcd_2e_abcd4e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'NaNbO3'
_chemical_formula_sum 'Na Nb O3'

loop_
_publ_author_name
 'P. Vousden'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 4
_journal_year 1951
_journal_page_first 545
_journal_page_last 551
_publ_Section_title
;
 The Structure of Ferroelectric Sodium Niobate at Room Temperature
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'NaNbO$_{3}$ Structure'
_aflow_proto 'ABC3_oP40_17_abcd_2e_abcd4e'
_aflow_params 'a,b/a,c/a,x_{1},x_{2},x_{3},x_{4},y_{5},y_{6},y_{7},y_{8
    ↪ },x_{9},y_{9},z_{9},x_{10},y_{10},z_{10},x_{11},y_{11},z_{11},
    ↪ x_{12},y_{12},z_{12},x_{13},y_{13},z_{13},x_{14},y_{14},z_{14}'
_aflow_params_values '5.504,1.01162790698,2.81976744186,0.481,-0.025,
    ↪ 0.019,0.525,0.519,0.025,-0.019,0.475,0.014,0.005,0.375,0.514,
    ↪ 0.505,0.375,0.25,0.25,0.375,0.75,0.25,0.375,0.25,0.75,0.375,
    ↪ 0.75,0.75,0.375'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP40'

_symmetry_space_group_name_H-M "P 2 2 21"
_symmetry_Int_Tables_number 17

_cell_length_a      5.50400
_cell_length_b      5.56800
_cell_length_c      15.52000
_cell_angle_alpha   90.00000
_cell_angle_beta    90.00000
_cell_angle_gamma   90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z+1/2
4 -x,-y,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
```

```
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Na1 Na 2 a 0.48100 0.00000 0.00000 1.00000
O1 O 2 a -0.02500 0.00000 0.00000 1.00000
Na2 Na 2 b 0.01900 0.50000 0.00000 1.00000
O2 O 2 b 0.52500 0.50000 0.00000 1.00000
Na3 Na 2 c 0.00000 0.51900 0.25000 1.00000
O3 O 2 c 0.00000 0.02500 0.25000 1.00000
Na4 Na 2 d 0.50000 -0.01900 0.25000 1.00000
O4 O 2 d 0.50000 0.47500 0.25000 1.00000
Nb1 Nb 4 e 0.01400 0.00500 0.37500 1.00000
Nb2 Nb 4 e 0.51400 0.50500 0.37500 1.00000
O5 O 4 e 0.25000 0.25000 0.37500 1.00000
O6 O 4 e 0.75000 0.25000 0.37500 1.00000
O7 O 4 e 0.25000 0.75000 0.37500 1.00000
O8 O 4 e 0.75000 0.75000 0.37500 1.00000
```

$NaNbO_3$: ABC3_oP40_17_abcd_2e_abcd4e - POSCAR

```
ABC3_oP40_17_abcd_2e_abcd4e & a,b/a,c/a,x1,x2,x3,x4,y5,y6,y7,y8,x9,y9,z9
      ↪ ,x10,y10,z10,x11,y11,z11,x12,y12,z12,x13,y13,z13,x14,y14,z14 --
      ↪ params=5.504,1.01162790698,2.81976744186,0.481,-0.025,0.019,
      ↪ 0.525,0.519,0.025,-0.019,0.475,0.014,0.005,0.375,0.514,0.505,
      ↪ 0.375,0.25,0.25,0.375,0.75,0.25,0.375,0.25,0.75,0.375,0.75,0.75
      ↪ ,0.375 & P222_{1} D_{2}^{2} #17 (a^2b^2c^2d^2e^6) & oP40 & None
      ↪ & NaNbO3 & NaNbO3 & P. Vousden, Acta Cryst. 4, 545-551 (1951)
   1.00000000000000
   5.50400000000000    0.00000000000000    0.00000000000000
   0.00000000000000    5.56800000000000    0.00000000000000
   0.00000000000000    0.00000000000000   15.52000000000000
     Na     Nb      O
      8      8     24
Direct
   0.48100000000000    0.00000000000000    0.00000000000000   Na   (2a)
  -0.48100000000000    0.00000000000000    0.50000000000000   Na   (2a)
   0.01900000000000    0.50000000000000    0.00000000000000   Na   (2b)
  -0.01900000000000    0.50000000000000    0.50000000000000   Na   (2b)
   0.00000000000000    0.51900000000000    0.25000000000000   Na   (2c)
   0.00000000000000   -0.51900000000000    0.75000000000000   Na   (2c)
   0.50000000000000   -0.01900000000000    0.25000000000000   Na   (2d)
   0.50000000000000    0.01900000000000    0.75000000000000   Na   (2d)
   0.01400000000000    0.00500000000000    0.37500000000000   Nb   (4e)
  -0.01400000000000   -0.00500000000000    0.87500000000000   Nb   (4e)
  -0.01400000000000    0.00500000000000    0.12500000000000   Nb   (4e)
   0.01400000000000   -0.00500000000000   -0.37500000000000   Nb   (4e)
   0.51400000000000    0.50500000000000    0.37500000000000   Nb   (4e)
  -0.51400000000000   -0.50500000000000    0.87500000000000   Nb   (4e)
  -0.51400000000000    0.50500000000000    0.12500000000000   Nb   (4e)
   0.51400000000000   -0.50500000000000   -0.37500000000000   Nb   (4e)
  -0.02500000000000    0.00000000000000    0.00000000000000   O    (2a)
   0.02500000000000    0.00000000000000    0.50000000000000   O    (2a)
   0.52500000000000    0.50000000000000    0.00000000000000   O    (2b)
  -0.52500000000000    0.50000000000000    0.50000000000000   O    (2b)
   0.00000000000000    0.02500000000000    0.25000000000000   O    (2c)
   0.00000000000000   -0.02500000000000    0.75000000000000   O    (2c)
   0.50000000000000    0.47500000000000    0.25000000000000   O    (2d)
   0.50000000000000   -0.47500000000000    0.75000000000000   O    (2d)
   0.25000000000000    0.25000000000000    0.37500000000000   O    (4e)
  -0.25000000000000   -0.25000000000000    0.87500000000000   O    (4e)
  -0.25000000000000    0.25000000000000    0.12500000000000   O    (4e)
   0.25000000000000   -0.25000000000000   -0.37500000000000   O    (4e)
   0.75000000000000    0.25000000000000    0.37500000000000   O    (4e)
  -0.75000000000000   -0.25000000000000    0.87500000000000   O    (4e)
  -0.75000000000000    0.25000000000000    0.12500000000000   O    (4e)
   0.75000000000000   -0.25000000000000   -0.37500000000000   O    (4e)
   0.25000000000000    0.75000000000000    0.37500000000000   O    (4e)
  -0.25000000000000   -0.75000000000000    0.87500000000000   O    (4e)
  -0.25000000000000    0.75000000000000    0.12500000000000   O    (4e)
   0.25000000000000   -0.75000000000000   -0.37500000000000   O    (4e)
   0.75000000000000    0.75000000000000    0.37500000000000   O    (4e)
  -0.75000000000000   -0.75000000000000    0.87500000000000   O    (4e)
  -0.75000000000000    0.75000000000000    0.12500000000000   O    (4e)
   0.75000000000000   -0.75000000000000   -0.37500000000000   O    (4e)
```

γ-TeO_2: A2B_oP12_18_2c_c - CIF

```
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_audit_creation_method FINDSYM

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_chemical_formula_sum 'O2 Te'

loop_
_publ_author_name
 'J. C. {Champarnaud-Mesjard}'
 'S. Blanchandin'
 'P. Thomas'
 'A. Mirgorodsky'
 'T. {Merle-M\\'{e}jean}'
 'B. Frit'
_journal_name_full_name
;
 Journal of Physics and Chemistry of Solids
;
_journal_volume 61
_journal_year 2000
_journal_page_first 1499
_journal_page_last 1507
_publ_Section_title
;
 Crystal structure, Raman spectrum and lattice dynamics of a new
      ↪ metastable form of tellurium dioxide: $\gamma$-TeO$_{2}$
;

# Found in {\em Ab initio} study of the vibrational properties of
      ↪ crystalline TeO$_{2}$: The $\alpha$, $\beta$, and $\gamma$
      ↪ phases, 2006

_aflow_title '$\gamma$-TeO$_{2}$ Structure'
_aflow_proto 'A2B_oP12_18_2c_c'
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_aflow_params_values '4.898,1.75091874234,0.888321763985,0.759,0.281,
      ↪ 0.173,0.855,0.036,0.727,-0.0304,0.1016,0.1358'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP12'

_symmetry_space_group_name_H-M "P 21 21 2"
_symmetry_Int_Tables_number 18

_cell_length_a     4.89800
_cell_length_b     8.57600
_cell_length_c     4.35100
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
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2 x+1/2,-y+1/2,-z
3 -x+1/2,y+1/2,-z
4 -x,-y,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 4 c 0.75900 0.28100 0.17300 1.00000
O2 O 4 c 0.85500 0.03600 0.72700 1.00000
Te1 Te 4 c -0.03040 0.10160 0.13580 1.00000
```

γ-TeO_2: A2B_oP12_18_2c_c - POSCAR

```
A2B_oP12_18_2c_c & a,b/a,c/a,x1,y1,z1,x2,y2,z2,x3,y3,z3 --params=4.898,
      ↪ 1.75091874234,0.888321763985,0.759,0.281,0.173,0.855,0.036,
      ↪ 0.727,-0.0304,0.1016,0.1358 & P2_{1}2_{1}2 D_{2}^{3} #18 (c^3)
      ↪ & oP12 & None & O2Te & O2Te & J. C. {Champarnaud-Mesjard} et
      ↪ al., J. Phys. Chem. Solids 61, 1499-1507 (2000)
   1.00000000000000
   4.89800000000000    0.00000000000000    0.00000000000000
   0.00000000000000    8.57600000000000    0.00000000000000
   0.00000000000000    0.00000000000000    4.35100000000000
     O     Te
     8      4
Direct
   0.75900000000000    0.28100000000000    0.17300000000000   O    (4c)
  -0.75900000000000   -0.28100000000000    0.17300000000000   O    (4c)
  -0.25900000000000    0.78100000000000   -0.17300000000000   O    (4c)
   1.25900000000000    0.21900000000000   -0.17300000000000   O    (4c)
   0.85500000000000    0.03600000000000    0.72700000000000   O    (4c)
  -0.85500000000000   -0.03600000000000    0.72700000000000   O    (4c)
  -0.35500000000000    0.53600000000000   -0.72700000000000   O    (4c)
   1.35500000000000    0.46400000000000   -0.72700000000000   O    (4c)
  -0.03040000000000    0.10160000000000    0.13580000000000   Te   (4c)
   0.03040000000000   -0.10160000000000    0.13580000000000   Te   (4c)
   0.53040000000000    0.60160000000000   -0.13580000000000   Te   (4c)
   0.46960000000000    0.39840000000000   -0.13580000000000   Te   (4c)
```

Diamminetriamidodizinc Chloride ($[Zn_2(NH_3)_2(NH_2)_3]Cl$): AB12C5D2_oP40_18_a_6c_b2c_c - CIF

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_chemical_formula_sum 'Cl H12 N5 Zn2'

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_publ_author_name
 'T. M. M. Richter'
 'S. Strobel'
 'N. S. A. Alt'
 'E. Schl\"{u}cker'
 'R. Niewa'
_journal_name_full_name
;
 Inorganics
;
_journal_volume 4
_journal_year 2016
_journal_page_first 41
_journal_page_last 41
_publ_Section_title
;
 Ammonothermal Synthesis and Crystal Structures of
      ↪ Diamminetriamidodizinc Chloride [Zn$_{2}$(NH$_{3}$)$_{2}$(NH$_
      ↪ {2}$)$_{3}$]Cl and Diamminemonoamidozinc Bromide [Zn(NH$_{3}$)
      ↪ $_{2}$(NH$_{2}$)]Br
```

```
;

_aflow_title 'Diamminetriamidodizinc Chloride ([Zn$_{2}$(NH$_{3}$)$_{2}$
      ↪ (NH$_{2}$)$_{3}$]Cl) Structure'
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      ↪ },x_{5},y_{5},z_{5},x_{6},y_{6},z_{6},x_{7},y_{7},z_{7},x_{8},
      ↪ y_{8},z_{8},x_{9},y_{9},z_{9},x_{10},y_{10},z_{10},x_{11},y_{11
      ↪ },z_{11}'
_aflow_params_values '5.7715,1.77352508014,1.13412457767,-0.0434,0.395,
      ↪ 0.35,0.248,0.67,0.5,0.354,0.67,0.0,0.259,-0.09,0.09,0.36,-0.04,
      ↪ 0.41,0.118,0.08,0.354,0.01,0.65,0.449,0.3291,0.553,0.007,0.3377
      ↪ ,0.853,0.1009,0.3376,0.5374'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP40'

_symmetry_space_group_name_H-M "P 21 21 2"
_symmetry_Int_Tables_number 18

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_cell_length_c    6.54560
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
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_space_group_symop_operation_xyz
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2 x+1/2,-y+1/2,-z
3 -x+1/2,y+1/2,-z
4 -x,-y,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cl1 Cl 2 a 0.00000 0.00000 -0.04340 1.00000
N1 N 2 b 0.00000 0.50000 0.39500 1.00000
H1 H 4 c 0.35000 0.24800 0.67000 1.00000
H2 H 4 c 0.50000 0.35400 0.67000 1.00000
H3 H 4 c 0.00000 0.25900 -0.09000 1.00000
H4 H 4 c 0.09000 0.36000 -0.04000 1.00000
H5 H 4 c 0.41000 0.11800 0.08000 1.00000
H6 H 4 c 0.35400 0.01000 0.65000 1.00000
N2 N 4 c 0.44900 0.32910 0.55300 1.00000
N3 N 4 c 0.00700 0.33770 0.85300 1.00000
Zn1 Zn 4 c 0.10090 0.33760 0.53740 1.00000
```

Diamminetriamidodizinc Chloride ($[Zn_2(NH_3)_2(NH_2)_3]Cl$): AB12C5D2_oP40_18_a_6c_b2c_c - POSCAR

```
AB12C5D2_oP40_18_a_6c_b2c_c & a,b/a,c/a,z1,z2,x3,y3,z3,x4,y4,z4,x5,y5,z5
      ↪ ,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10,x11,y11,z11 --
      ↪ params=5.7715,1.77352508014,1.13412457767,-0.0434,0.395,0.35,
      ↪ 0.248,0.67,0.5,0.354,0.67,0.0,0.259,-0.09,0.09,0.36,-0.04,0.41,
      ↪ 0.118,0.08,0.354,0.01,0.65,0.449,0.3291,0.553,0.007,0.3377,
      ↪ 0.853,0.1009,0.3376,0.5374 & P2_{1}2_{1}2 D_{2}^{3} #18 (abc^9)
      ↪ & oP40 & None & ClH12N5Zn2 & Diamminetriamidodizinc chloride &
      ↪ T. M. M. Richter et al., Inorganics 4, 41(2016)
   1.00000000000000
   5.77150000000000    0.00000000000000    0.00000000000000
   0.00000000000000   10.23590000000000    0.00000000000000
   0.00000000000000    0.00000000000000    6.54560000000000
     Cl   H   N   Zn
      2  24  10   4
Direct
   0.00000000000000    0.00000000000000   -0.04340000000000   Cl   (2a)
   0.50000000000000    0.50000000000000    0.04340000000000   Cl   (2a)
   0.35000000000000    0.24800000000000    0.67000000000000   H    (4c)
  -0.35000000000000   -0.24800000000000    0.67000000000000   H    (4c)
   0.15000000000000    0.74800000000000   -0.67000000000000   H    (4c)
   0.85000000000000    0.25200000000000   -0.67000000000000   H    (4c)
   0.50000000000000    0.35400000000000    0.67000000000000   H    (4c)
  -0.50000000000000   -0.35400000000000    0.67000000000000   H    (4c)
   0.00000000000000    0.85400000000000   -0.67000000000000   H    (4c)
   1.00000000000000    0.14600000000000   -0.67000000000000   H    (4c)
   0.00000000000000    0.25900000000000   -0.09000000000000   H    (4c)
   0.00000000000000   -0.25900000000000   -0.09000000000000   H    (4c)
   0.50000000000000    0.75900000000000    0.09000000000000   H    (4c)
   0.50000000000000    0.24100000000000    0.09000000000000   H    (4c)
   0.09000000000000    0.36000000000000   -0.04000000000000   H    (4c)
  -0.09000000000000   -0.36000000000000   -0.04000000000000   H    (4c)
   0.41000000000000    0.86000000000000    0.04000000000000   H    (4c)
   0.59000000000000    0.14000000000000    0.04000000000000   H    (4c)
   0.41000000000000    0.11800000000000    0.08000000000000   H    (4c)
  -0.41000000000000   -0.11800000000000    0.08000000000000   H    (4c)
   0.09000000000000    0.61800000000000   -0.08000000000000   H    (4c)
   0.91000000000000    0.38200000000000   -0.08000000000000   H    (4c)
   0.35400000000000    0.01000000000000    0.65000000000000   H    (4c)
  -0.35400000000000   -0.01000000000000    0.65000000000000   H    (4c)
   0.14600000000000    0.51000000000000   -0.65000000000000   H    (4c)
   0.85400000000000    0.49000000000000   -0.65000000000000   H    (4c)
   0.00000000000000    0.50000000000000    0.39500000000000   N    (2b)
   0.50000000000000    0.00000000000000   -0.39500000000000   N    (2b)
   0.44900000000000    0.32910000000000    0.55300000000000   N    (4c)
  -0.44900000000000   -0.32910000000000    0.55300000000000   N    (4c)
   0.05100000000000    0.82910000000000   -0.55300000000000   N    (4c)
   0.94900000000000    0.17090000000000   -0.55300000000000   N    (4c)
   0.00700000000000    0.33770000000000    0.85300000000000   N    (4c)
  -0.00700000000000   -0.33770000000000    0.85300000000000   N    (4c)
   0.49300000000000    0.83770000000000   -0.85300000000000   N    (4c)
   0.50700000000000    0.16230000000000   -0.85300000000000   N    (4c)
   0.10090000000000    0.33760000000000    0.53740000000000   Zn   (4c)
  -0.10090000000000   -0.33760000000000    0.53740000000000   Zn   (4c)
   0.39910000000000    0.83760000000000   -0.53740000000000   Zn   (4c)
   0.60090000000000    0.16240000000000   -0.53740000000000   Zn   (4c)
```

Morenosite ($NiSO_4{\cdot}7H_2O$, $H4_{12}$): A14BC11D_oP108_19_14a_a_11a_a - CIF

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_chemical_formula_sum 'H14 Ni O11 S'

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 'H. Ptasiewicz-Bak'
 'I. Olovsson'
 'G. J. McIntyre'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 53
_journal_year 1997
_journal_page_first 325
_journal_page_last 336
_publ_Section_title
;
 Charge Density in Orthorhombic NiSO$_{4}$$\cdot$7H$_{2}$O at Room
      ↪ Temperature and 25 K
;

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      ↪ Structure'
_aflow_proto 'A14BC11D_oP108_19_14a_a_11a_a'
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      ↪ x_{7},y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{10},y_
      ↪ {10},z_{10},x_{11},y_{11},z_{11},x_{12},y_{12},z_{12},x_{13},y_
      ↪ {13},z_{13},x_{14},y_{14},z_{14},x_{15},y_{15},z_{15},x_{16},y_
      ↪ {16},z_{16},x_{17},y_{17},z_{17},x_{18},y_{18},z_{18},x_{19},y_
      ↪ {19},z_{19},x_{20},y_{20},z_{20},x_{21},y_{21},z_{21},x_{22},y_
      ↪ {22},z_{22},x_{23},y_{23},z_{23},x_{24},y_{24},z_{24},x_{25},y_
      ↪ {25},z_{25},x_{26},y_{26},z_{26},x_{27},y_{27},z_{27}'
_aflow_params_values '6.706,1.75902177155,1.78183716075,0.109,0.7344,
      ↪ 0.7161,0.875,0.74,0.6932,0.298,-0.0819,0.7675,0.251,0.0439,
      ↪ 0.7498,0.689,-0.0769,0.7175,0.74,0.0473,0.6778,0.197,0.1238,
      ↪ 0.5537,0.007,0.1119,0.48,-0.023,0.8655,0.3955,0.779,-0.0807,
      ↪ 0.4437,0.319,0.7843,0.5251,0.396,-0.0866,0.4995,-0.05,-0.0782,-
      ↪ 0.0187,-0.002,-0.0288,0.8664,0.04106,-0.07911,0.60442,0.42704,
      ↪ 0.18381,0.57465,0.48169,0.35142,0.68668,0.69542,0.18889,0.70501
      ↪ ,0.36068,0.17853,0.77191,0.00529,0.7607,0.66624,0.18986,-
      ↪ 0.03077,0.74842,0.7752,-0.03247,0.66931,0.07373,0.08217,0.54556
      ↪ ,0.8931,0.87229,0.46088,0.29961,0.8655,0.537,-0.0694,-0.01048,-
      ↪ 0.0641,0.49094,0.2263,0.68354'
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_aflow_Pearson 'oP108'

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_symmetry_Int_Tables_number 19

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_cell_length_c    11.94900
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

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_space_group_symop_operation_xyz
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2 x+1/2,-y+1/2,-z
3 -x,y+1/2,-z+1/2
4 -x+1/2,-y,z+1/2

loop_
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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
H1 H 4 a 0.10900 0.73440 0.71610 1.00000
H2 H 4 a 0.87500 0.74000 0.69320 1.00000
H3 H 4 a 0.29800 -0.08190 0.76750 1.00000
H4 H 4 a 0.25100 0.04390 0.74980 1.00000
H5 H 4 a 0.68900 -0.07690 0.71750 1.00000
H6 H 4 a 0.74000 0.04730 0.67780 1.00000
H7 H 4 a 0.19700 0.12380 0.55370 1.00000
H8 H 4 a 0.00700 0.11190 0.48000 1.00000
H9 H 4 a -0.02300 0.86550 0.39550 1.00000
H10 H 4 a 0.77900 -0.08070 0.44370 1.00000
H11 H 4 a 0.31900 0.78430 0.52510 1.00000
H12 H 4 a 0.39600 -0.08660 0.49950 1.00000
H13 H 4 a -0.05000 -0.07820 -0.01870 1.00000
H14 H 4 a -0.00200 -0.02880 0.86640 1.00000
Ni1 Ni 4 a 0.04106 -0.07911 0.60442 1.00000
```

```
O1 O 4 a 0.42704 0.18381 0.57465 1.00000
O2 O 4 a 0.48169 0.35142 0.68668 1.00000
O3 O 4 a 0.69542 0.18889 0.70501 1.00000
O4 O 4 a 0.36068 0.17853 0.77191 1.00000
O5 O 4 a 0.00529 0.76070 0.66624 1.00000
O6 O 4 a 0.18986 −0.03077 0.74842 1.00000
O7 O 4 a 0.77520 −0.03247 0.66931 1.00000
O8 O 4 a 0.07373 0.08217 0.54556 1.00000
O9 O 4 a 0.89310 0.87229 0.46088 1.00000
O10 O 4 a 0.29961 0.86550 0.53700 1.00000
O11 O 4 a −0.06940 −0.01048 −0.06410 1.00000
S1 S 4 a 0.49094 0.22630 0.68354 1.00000
```

Morenosite ($NiSO_4{\cdot}7H_2O$, $H4_{12}$): A14BC11D_oP108_19_14a_a_11a_a - POSCAR

```
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    ↪ y4,z4,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10,
    ↪ x11,y11,z11,x12,y12,z12,x13,y13,z13,x14,y14,z14,x15,y15,z15,x16
    ↪ ,y16,z16,x17,y17,z17,x18,y18,z18,x19,y19,z19,x20,y20,z20,x21,
    ↪ y21,z21,x22,y22,z22,x23,y23,z23,x24,y24,z24,x25,y25,z25,x26,y26
    ↪ ,z26,x27,y27,z27 --params=6.706,1.75902177155,1.78183716075,
    ↪ 0.109,0.7344,0.7161,0.875,0.74,0.6932,0.298,-0.0819,0.7675,
    ↪ 0.251,0.0439,0.7498,0.689,-0.0769,0.7175,0.74,0.0473,0.6778,
    ↪ 0.197,0.1238,0.5537,0.007,0.1119,0.48,-0.023,0.8655,0.3955,
    ↪ 0.779,-0.0807,0.4437,0.319,0.7843,0.5251,0.396,-0.0866,0.4995,-
    ↪ 0.05,-0.0782,-0.0187,-0.002,-0.0288,0.8664,0.04106,-0.07911,
    ↪ 0.60442,0.42704,0.18381,0.57465,0.48169,0.35142,0.68668,0.69542
    ↪ ,0.18889,0.70501,0.36068,0.17853,0.77191,0.00529,0.7607,0.66624
    ↪ ,0.18986,-0.03077,0.74842,0.7752,-0.03247,0.66931,0.07373,
    ↪ 0.08217,0.54556,0.8931,0.87229,0.46088,0.29961,0.8655,0.537,-
    ↪ 0.0694,-0.01048,-0.0641,0.49094,0.2263,0.68354 & P2_{1}2_{1}2_{
    ↪ 1} D_{2}^{4} #19 (a^27) & oP108 & $H4_{12}$ & H14NiO11S &
    ↪ Morenosite & H. Ptasiewicz-Bak and I. Olovsson and G. J.
    ↪ McIntyre, Acta Crystallogr. Sect. B Struct. Sci. 53, 325-336 (
    ↪ 1997)
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   0.00000000000000  11.79600000000000   0.00000000000000
   0.00000000000000   0.00000000000000  11.94900000000000
     H    Ni     O     S
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   0.25100000000000   0.04390000000000   0.74980000000000   H   (4a)
   0.24900000000000  -0.04390000000000   1.24980000000000   H   (4a)
  -0.25100000000000   0.54390000000000  -0.24980000000000   H   (4a)
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  -0.68900000000000   0.42310000000000  -0.21750000000000   H   (4a)
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Wülfingite (ϵ-$Zn(OH)_2$, $C31$): A2B2C_oP20_19_2a_2a_a - CIF

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 'H. D. Lutz'
 'W. Kockelmann'
 'H. Jacobs'
_journal_name_full_name
;
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;
_journal_volume 624
_journal_year 1998
_journal_page_first 1130
_journal_page_last 1136
_publ_Section_title
;
 Kristallstrukturen und Wasserstoffbr\"{u}ckenbindungen bei $\beta$-Be(
    ↪ OH)$_{2}$ und $\epsilon$-Zn(OH)$_{2}$
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# Found in The American Mineralogist Crystal Structure Database, 2003

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    ↪ 0.07096,0.64832,0.6242'
_aflow_Strukturbericht '$C31$'
_aflow_Pearson 'oP20'

_symmetry_space_group_name_H-M "P 21 21 21"
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H2 H 4 a 0.28000 0.79000 0.17000 1.00000
O1 O 4 a 0.11550 0.12560 0.07950 1.00000
O2 O 4 a 0.16700 0.31720 0.71980 1.00000
Zn1 Zn 4 a 0.07096 0.64832 0.62420 1.00000
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Wülfingite (ϵ-$Zn(OH)_2$, $C31$): A2B2C_oP20_19_2a_2a_a - POSCAR

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    ↪ 0.07096,0.64832,0.6242 & P2_{1}2_{1}2_{1} D_{2}^{4} #19 (a^5) &
    ↪  oP20 & $C31$ & H2O2Zn & W\"{u}lfingite & R. Stahl et al., Z.
    ↪ Anorg. Allg. Chem. 624, 1130-1136 (1998)
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Ferroelectric $NH_4H_2PO_4$: A6BC4D_oP48_19_6a_a_4a_a - CIF

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 'S. Akahoshi'
 'K. Hukuda'
 'T. Yagi'
_journal_name_full_name
;
 Journal of the Physical Society of Japan
;
_journal_volume 56
_journal_year 1987
_journal_page_first 2223
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;
 Refinement of the Crystal Structure of NH$_{4}$H$_{2}$PO$_{4}$ above
    ↪ and below Antiferroelectric Phase Transition Temperature
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    ↪ x_{7},y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{10},y_
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    ↪ 0.789,0.348,0.122,0.154,0.298,0.892,0.011,0.4837,0.5016,0.082,
    ↪ 0.6443,0.1192,-0.0853,0.3498,0.1145,0.8628,0.5773,0.872,0.1571,
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_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP48'

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_symmetry_Int_Tables_number 19

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_cell_angle_beta   90.00000
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3 -x,y+1/2,-z+1/2
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H2 H 4 a 0.00000 0.38800 0.54700 1.00000
H3 H 4 a -0.06800 0.50400 0.42100 1.00000
H4 H 4 a 0.11300 0.48700 0.45300 1.00000
H5 H 4 a 0.78900 0.34800 0.12200 1.00000
H6 H 4 a 0.15400 0.29800 0.89200 1.00000
N1 N 4 a 0.01100 0.48370 0.50160 1.00000
O1 O 4 a 0.08200 0.64430 0.11920 1.00000
O2 O 4 a -0.08530 0.34980 0.11450 1.00000
O3 O 4 a 0.86280 0.57730 0.87200 1.00000
O4 O 4 a 0.15710 0.41000 0.89740 1.00000
P1 P 4 a 0.00000 0.50190 0.00040 1.00000
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Ferroelectric $NH_4H_2PO_4$: A6BC4D_oP48_19_6a_a_4a_a - POSCAR

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    ↪ 0.487,0.453,0.789,0.348,0.122,0.154,0.298,0.892,0.011,0.4837,
    ↪ 0.5016,0.082,0.6443,0.1192,-0.0853,0.3498,0.1145,0.8628,0.5773,
    ↪ 0.872,0.1571,0.41,0.8974,0.0,0.5019,0.0004 & P2_{1}2_{1}2_{1}
    ↪ D_{2}^{4} #19 (a^12) & oP48 & None & H6NO4P & H6NO4P & T.
    ↪ Fukami et al., J. Phys. Soc. Jpn. 56, 2223-2224 (1987)
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   0.00000000000000    0.00000000000000    7.48800000000000
     H     N     O     P
    24     4    16     4
Direct
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   0.51200000000000   -0.07400000000000   -0.58700000000000   H   (4a)
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β-Arabinose [$(CH_2O)_{20}$]: AB2C_oP80_19_5a_10a_5a - CIF

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 Acta Chemica Scandinavica
;
_journal_volume 15
_journal_year 1961
_journal_page_first 16
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;
 Refinement of the Crystal Structure of $\beta$-Arabinose
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      ↪ {10},z_{10},x_{11},y_{11},z_{11},x_{12},y_{12},z_{12},x_{13},y_
      ↪ {13},z_{13},x_{14},y_{14},z_{14},x_{15},y_{15},z_{15},x_{16},y_
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      ↪ {19},z_{19},x_{20},y_{20},z_{20}'
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      ↪ 0.303,0.811,0.422,-0.043,0.843,0.507,0.024,-0.028,0.564,0.22,-
      ↪ 0.061,0.351,0.6,-0.046,0.38,0.539,0.767,0.892,0.173,0.737,0.567
      ↪ ,0.875,0.852,-0.056,0.5625,-0.0901,0.42,0.6075,0.8019,0.771,
      ↪ 0.191,0.7545,0.737,-0.004,0.8781,-0.074,0.325,-0.0368,0.711'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP80'

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4 -x+1/2,-y,z+1/2

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C3 C 4 a 0.25100 0.81980 0.63100 1.00000
C4 C 4 a 0.06950 0.87120 0.64900 1.00000
C5 C 4 a 0.15000 -0.05860 0.55300 1.00000
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H2 H 4 a 0.37600 0.84700 0.00000 1.00000
H3 H 4 a 0.30300 0.81100 0.42200 1.00000
H4 H 4 a -0.04300 0.84300 0.50700 1.00000
H5 H 4 a 0.02400 -0.02800 0.56400 1.00000
H6 H 4 a 0.22000 -0.06100 0.35100 1.00000
H7 H 4 a 0.60000 -0.04600 0.38000 1.00000
H8 H 4 a 0.53900 0.76700 0.89200 1.00000
H9 H 4 a 0.17300 0.73700 0.56700 1.00000
H10 H 4 a 0.87500 0.85200 -0.05600 1.00000
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O2 O 4 a 0.60750 0.80190 0.77100 1.00000
O3 O 4 a 0.19100 0.75450 0.73700 1.00000
O4 O 4 a -0.00400 0.87810 -0.07400 1.00000
O5 O 4 a 0.32500 -0.03680 0.71100 1.00000
```

β-Arabinose [$(CH_2O)_{20}$]: AB2C_oP80_19_5a_10a_5a - POSCAR

```
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      ↪ x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10,x11,
      ↪ y11,z11,x12,y12,z12,x13,y13,z13,x14,y14,z14,x15,y15,z15,x16,y16
      ↪ ,z16,x17,y17,z17,x18,y18,z18,x19,y19,z19,x20,y20,z20 --params=
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      ↪ 0.422,-0.043,0.843,0.507,0.024,-0.028,0.564,0.22,-0.061,0.351,
      ↪ 0.6,-0.046,0.38,0.539,0.767,0.892,0.173,0.737,0.567,0.875,0.852
      ↪ ,-0.056,0.5625,-0.0901,0.42,0.6075,0.8019,0.771,0.191,0.7545,
      ↪ 0.737,-0.004,0.8781,-0.074,0.325,-0.0368,0.711 & P2_{1}2_{1}2_{
      ↪ 1} D_{2}^{4} #19 (a^20) & oP80 & None & CH2O & CH2O & A.
      ↪ Hordvik, Acta Chem. Scand. 15, 16-30 (1961)
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      20   40   20
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$NaAlCl_4$: AB4C_oP24_19_a_4a_a - CIF

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# Found in The American Mineralogist Crystal Structure Database, 2003

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$NaAlCl_4$: AB4C_oP24_19_a_4a_a - POSCAR

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      ↪ & oP24 & None & AlCl4Na & AlCl4Na & G. Mairesse and P. Barbier
      ↪ and J.-P. Wignacourt, Acta Crystallogr. Sect. B Struct. Sci. 35
      ↪ , 1573-1580 (1979)
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     4    16     4
Direct
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  -0.33503000000000   0.57337000000000  -0.37726000000000   Cl   (4a)
   0.83503000000000   0.42663000000000  -0.87726000000000   Cl   (4a)
   0.52257000000000   0.07458000000000   0.15367000000000   Cl   (4a)
  -0.02257000000000  -0.07458000000000   0.65367000000000   Cl   (4a)
  -0.52257000000000   0.57458000000000   0.34633000000000   Cl   (4a)
   1.02257000000000   0.42542000000000  -0.15367000000000   Cl   (4a)
   0.50874000000000   0.55281000000000  -0.03214000000000   Cl   (4a)
  -0.00874000000000  -0.55281000000000   0.46786000000000   Cl   (4a)
  -0.50874000000000   1.05281000000000   0.53214000000000   Cl   (4a)
   1.00874000000000  -0.05281000000000   0.03214000000000   Cl   (4a)
   0.71343000000000   0.31122000000000   0.37466000000000   Na   (4a)
  -0.21343000000000  -0.31122000000000   0.87466000000000   Na   (4a)
  -0.71343000000000   0.81122000000000   0.12534000000000   Na   (4a)
   1.21343000000000   0.18878000000000  -0.37466000000000   Na   (4a)
```

NaP: AB_oP16_19_2a_2a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'NaP'
_chemical_formula_sum 'Na P'

loop_
_publ_author_name
 'H. G. {von Schnering}'
 'W. H\"{o}nle'
_journal_name_full_name
;
 Zeitschrift fur Anorganische und Allgemeine Chemie
;
_journal_volume 456
_journal_year 1979
_journal_page_first 194
_journal_page_last 206
_publ_Section_title
;
 Zur Chemie und Strukturchemie der Phosphide und Polyphosphide. 20.
       ↪ Darstellung, Struktur und Eigenschaften der
       ↪ Alkalimetallmonophosphide NaP und KP
;

# Found in NaP Crystal Structure, {PAULING FILE in: Inorganic Solid
      ↪ Phases, SpringerMaterials (online database)}, 2016

_aflow_title 'NaP Structure'
_aflow_proto 'AB_oP16_19_2a_2a'
_aflow_params 'a,b/a,c/a,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3},y_{3
      ↪ },z_{3},x_{4},y_{4},z_{4}'
_aflow_params_values '6.038,0.934580987082,1.67969526333,0.4174,-0.0911,
      ↪ 0.0318,0.1338,0.6367,0.3313,0.3086,0.1404,0.2838,0.4287,0.402,
      ↪ 0.1341'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP16'

_symmetry_space_group_name_H-M "P 21 21 21"
_symmetry_Int_Tables_number 19

_cell_length_a     6.03800
_cell_length_b     5.64300
_cell_length_c     10.14200
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z
3 -x,y+1/2,-z+1/2
4 -x+1/2,-y,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Na1 Na 4 a 0.41740 -0.09110 0.03180 1.00000
Na2 Na 4 a 0.13380 0.63670 0.33130 1.00000
P1 P 4 a 0.30860 0.14040 0.28380 1.00000
P2 P 4 a 0.42870 0.40200 0.13410 1.00000
```

NaP: AB_oP16_19_2a_2a - POSCAR

```
AB_oP16_19_2a_2a & a,b/a,c/a,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,z4 --
      ↪ params=6.038,0.934580987082,1.67969526333,0.4174,-0.0911,0.0318
      ↪ ,0.1338,0.6367,0.3313,0.3086,0.1404,0.2838,0.4287,0.402,0.1341
      ↪ & P2_{1}2_{1}2_{1} D_{2}^{4} #19 (a^4) & oP16 & None & NaP &
      ↪ NaP & H. G. {von Schnering} and W. H\"{o}nle, Z. Anorg. Allg.
      ↪ Chem. 456, 194-206 (1979)
   1.00000000000000
   6.03800000000000   0.00000000000000   0.00000000000000
   0.00000000000000   5.64300000000000   0.00000000000000
   0.00000000000000   0.00000000000000  10.14200000000000
    Na    P
     8    8
Direct
   0.41740000000000  -0.09110000000000   0.03180000000000   Na   (4a)
   0.08260000000000   0.09110000000000   0.53180000000000   Na   (4a)
  -0.41740000000000   0.40890000000000   0.46820000000000   Na   (4a)
   0.91740000000000   0.59110000000000  -0.03180000000000   Na   (4a)
   0.13380000000000   0.63670000000000   0.33130000000000   Na   (4a)
   0.36620000000000  -0.63670000000000   0.83130000000000   Na   (4a)
  -0.13380000000000   1.13670000000000   0.16870000000000   Na   (4a)
   0.63380000000000  -0.13670000000000  -0.33130000000000   Na   (4a)
   0.30860000000000   0.14040000000000   0.28380000000000    P   (4a)
   0.19140000000000  -0.14040000000000   0.78380000000000    P   (4a)
  -0.30860000000000   0.64040000000000   0.21620000000000    P   (4a)
   0.80860000000000   0.35960000000000  -0.28380000000000    P   (4a)
   0.42870000000000   0.40200000000000   0.13410000000000    P   (4a)
   0.07130000000000  -0.40200000000000   0.63410000000000    P   (4a)
  -0.42870000000000   0.90200000000000   0.36590000000000    P   (4a)
   0.92870000000000   0.09800000000000  -0.13410000000000    P   (4a)
```

$D0_7$ (CrO_3) (*obsolete*): AB3_oC16_20_a_bc - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'CrO3'
_chemical_formula_sum 'Cr O3'

loop_
```

```
_publ_author_name
 'H. Br\"{a}kken'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 78
_journal_year 1931
_journal_page_first 484
_journal_page_last 488
_publ_Section_title
;
 Die Kristallstrukturen der Trioxyde von Chrom, Molybd\"{a}n und Wolfram
;

# Found in The Crystal Structure of Chromium Trioxide, 1950

_aflow_title '$D0_{7}$ (CrO$_{3}$) ({\em{obsolete}}) Structure'
_aflow_proto 'AB3_oC16_20_a_bc'
_aflow_params 'a,b/a,c/a,x_{1},y_{2},x_{3},y_{3},z_{3}'
_aflow_params_values '8.46,0.563829787234,0.673758865248,0.33333,0.33333
      ↪ ,0.16667,-0.16667,0.25'
_aflow_Strukturbericht '$D0_{7}$'
_aflow_Pearson 'oC16'

_symmetry_space_group_name_H-M "C 2 2 21"
_symmetry_Int_Tables_number 20

_cell_length_a     8.46000
_cell_length_b     4.77000
_cell_length_c     5.70000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z+1/2
4 -x,-y,z+1/2
5 x+1/2,y+1/2,z
6 x+1/2,-y+1/2,-z
7 -x+1/2,y+1/2,-z+1/2
8 -x+1/2,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cr1 Cr 4 a 0.33333 0.00000 0.00000 1.00000
O1 O 4 b 0.00000 0.33333 0.25000 1.00000
O2 O 8 c 0.16667 -0.16667 0.25000 1.00000
```

$D0_7$ (CrO_3) (*obsolete*): AB3_oC16_20_a_bc - POSCAR

```
AB3_oC16_20_a_bc & a,b/a,c/a,x1,y2,x3,y3,z3 --params=8.46,0.563829787234
      ↪ ,0.673758865248,0.33333,0.33333,0.16667,-0.16667,0.25 & C222_{1
      ↪ } D_{2}^{5} #20 (abc) & oC16 & $D0_{7}$ & CrO3 & CrO3 & H. Br
      ↪ \"{a}kken, Zeitschrift f{"u}r Kristallographie - Crystalline
      ↪ Materials 78, 484-488 (1931)
   1.00000000000000
   4.23000000000000  -2.38500000000000   0.00000000000000
   4.23000000000000   2.38500000000000   0.00000000000000
   0.00000000000000   0.00000000000000   5.70000000000000
    Cr    O
     2    6
Direct
   0.33333000000000   0.33333000000000   0.00000000000000  Cr   (4a)
  -0.33333000000000  -0.33333000000000   0.50000000000000  Cr   (4a)
  -0.33333000000000   0.33333000000000   0.25000000000000  O    (4b)
   0.33333000000000  -0.33333000000000   0.75000000000000  O    (4b)
   0.33334000000000   0.00000000000000   0.25000000000000  O    (8c)
  -0.33334000000000   0.00000000000000   0.75000000000000  O    (8c)
   0.00000000000000  -0.33334000000000   0.25000000000000  O    (8c)
   0.00000000000000   0.33334000000000  -0.25000000000000  O    (8c)
```

$AlPO_4$ "low cristobalite type": AB4C_oC24_20_b_2c_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Low cristobalite type'
_chemical_formula_sum 'Al O4 P'

loop_
_publ_author_name
 'R. C. L. Mooney'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 9
_journal_year 1956
_journal_page_first 728
_journal_page_last 734
_publ_Section_title
;
 The crystal structure of aluminium phosphate and gallium phosphate,
      ↪ low-cristobalite type
;

# Found in The $\alpha-\beta$ phase transition in AlPO$_{4}$
      ↪ cristobalite: Symmetry analysis, domain structure and
      ↪ transition dynamics, 1994

_aflow_title 'AlPO$_{4}$ ``low cristobalite type\'\' Structure'
_aflow_proto 'AB4C_oC24_20_b_2c_a'
_aflow_params 'a,b/a,c/a,x_{1},y_{2},x_{3},y_{3},z_{3},x_{4},y_{4},z_{4
      ↪ }'
_aflow_params_values '7.009,1.0,0.999571978884,0.306,0.198,0.179,0.058,
      ↪ 0.172,0.433,0.17,0.941'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oC24'

_symmetry_space_group_name_H-M "C 2 2 21"
_symmetry_Int_Tables_number 20

_cell_length_a     7.00900
_cell_length_b     7.00900
_cell_length_c     7.00600
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z+1/2
4 -x,-y,z+1/2
5 x+1/2,y+1/2,z
6 x+1/2,-y+1/2,-z
7 -x+1/2,y+1/2,-z+1/2
8 -x+1/2,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
P1 P 4 a 0.30600 0.00000 0.00000 1.00000
Al1 Al 4 b 0.00000 0.19800 0.25000 1.00000
O1 O 8 c 0.17900 0.05800 0.17200 1.00000
O2 O 8 c 0.43300 0.17000 0.94100 1.00000
```

$AlPO_4$ "low cristobalite type": AB4C_oC24_20_b_2c_a - POSCAR

```
AB4C_oC24_20_b_2c_a & a,b/a,c/a,x1,y2,x3,y3,z3,x4,y4,z4 --params=7.009,
      ↪ 1.0,0.999571978884,0.306,0.198,0.179,0.058,0.172,0.433,0.17,
      ↪ 0.941 & C222_{1} D_{2}^{5} #20 (abc^2) & oC24 & None & AlO4P &
      ↪ Low cristobalite type & R. C. L. Mooney, Acta Cryst. 9, 728-734
      ↪ (1956)
   1.00000000000000
   3.50450000000000  -3.50450000000000   0.00000000000000
   3.50450000000000   3.50450000000000   0.00000000000000
   0.00000000000000   0.00000000000000   7.00600000000000
    Al    O    P
     2    8    2
Direct
  -0.19800000000000   0.19800000000000   0.25000000000000  Al   (4b)
   0.19800000000000  -0.19800000000000   0.75000000000000  Al   (4b)
   0.12100000000000   0.23700000000000   0.17200000000000  O    (8c)
  -0.12100000000000  -0.23700000000000   0.67200000000000  O    (8c)
  -0.23700000000000  -0.12100000000000   0.32800000000000  O    (8c)
   0.23700000000000   0.12100000000000  -0.17200000000000  O    (8c)
   0.26300000000000   0.60300000000000   0.94100000000000  O    (8c)
  -0.26300000000000  -0.60300000000000   1.44100000000000  O    (8c)
  -0.60300000000000  -0.26300000000000  -0.44100000000000  O    (8c)
   0.60300000000000   0.26300000000000  -0.94100000000000  O    (8c)
   0.30600000000000   0.30600000000000   0.00000000000000  P    (4a)
  -0.30600000000000  -0.30600000000000   0.50000000000000  P    (4a)
```

Tl_2AlF_5 ($K3_3$): AB5C2_oC32_20_b_a2bc_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'AlF5Tl2'
_chemical_formula_sum 'Al F5 Tl2'

loop_
_publ_author_name
 'C. Brosset'
_journal_name_full_name
;
 Zeitschrift fur Anorganische und Allgemeine Chemie
;
_journal_volume 235
_journal_year 1937
_journal_page_first 139
_journal_page_last 147
_publ_Section_title
;
 Herstellung und Kristallbau der Verbindungen TlAlF$_{4}$ und Tl$_{2}
      ↪ $AlF$_{5}$
;
```

```
# Found in A Structural Classification of Fluoaluminates, 1950

_aflow_title 'Tl$_{2}$AlF$_{5}$ ($K3_{3}$) Structure'
_aflow_proto 'AB5C2_oC32_20_b_a2bc_c'
_aflow_params 'a,b/a,c/a,x_{1},y_{2},y_{3},y_{4},x_{5},y_{5},z_{5},x_{6
    ↪ },y_{6},z_{6}'
_aflow_params_values '10.06,0.819085487078,0.741550695825,0.033,0.0,0.23
    ↪ ,0.78,0.19,0.0,0.29,0.29,0.2,0.0'
_aflow_Strukturbericht '$K3_{3}$'
_aflow_Pearson 'oC32'

_symmetry_space_group_name_H-M "C 2 2 21"
_symmetry_Int_Tables_number 20

_cell_length_a      10.06000
_cell_length_b      8.24000
_cell_length_c      7.46000
_cell_angle_alpha   90.00000
_cell_angle_beta    90.00000
_cell_angle_gamma   90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z+1/2
4 -x,-y,z+1/2
5 x+1/2,y+1/2,z
6 x+1/2,-y+1/2,-z
7 -x+1/2,y+1/2,-z+1/2
8 -x+1/2,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
F1 F 4 a 0.03300 0.00000 0.00000 1.00000
Al1 Al 4 b 0.00000 0.00000 0.25000 1.00000
F2 F 4 b 0.00000 0.23000 0.25000 1.00000
F3 F 4 b 0.00000 0.78000 0.25000 1.00000
F4 F 8 c 0.19000 0.00000 0.29000 1.00000
Tl1 Tl 8 c 0.29000 0.20000 0.00000 1.00000
```

Tl_2AlF_5 ($K3_3$): AB5C2_oC32_20_b_a2bc_c - POSCAR

```
AB5C2_oC32_20_b_a2bc_c & a,b/a,c/a,x1,y2,y3,y4,x5,y5,z5,x6,y6,z6 --
    ↪ params=10.06,0.819085487078,0.741550695825,0.033,0.0,0.23,0.78,
    ↪ 0.19,0.0,0.29,0.29,0.2,0.0 & C222_{1}D_{2}^{5}#20 (ab^3c^2) &
    ↪  oC32 & $K3_{3}$ & AlF5Tl2 & AlF5Tl2 & C. Brosset, Z. Anorg.
    ↪ Allg. Chem. 235, 139-147 (1937)
   1.00000000000000
   5.03000000000000  -4.12000000000000   0.00000000000000
   5.03000000000000   4.12000000000000   0.00000000000000
   0.00000000000000   0.00000000000000   7.46000000000000
    Al    F    Tl
     2   10     4
Direct
   0.00000000000000   0.00000000000000   0.25000000000000   Al   (4b)
   0.00000000000000   0.00000000000000   0.75000000000000   Al   (4b)
   0.03300000000000   0.03300000000000   0.00000000000000   F    (4a)
  -0.03300000000000  -0.03300000000000   0.50000000000000   F    (4a)
  -0.23000000000000   0.23000000000000   0.25000000000000   F    (4b)
   0.23000000000000  -0.23000000000000   0.75000000000000   F    (4b)
  -0.78000000000000   0.78000000000000   0.25000000000000   F    (4b)
   0.78000000000000  -0.78000000000000   0.75000000000000   F    (4b)
   0.19000000000000   0.19000000000000   0.29000000000000   F    (8c)
  -0.19000000000000  -0.19000000000000   0.79000000000000   F    (8c)
  -0.19000000000000  -0.19000000000000   0.21000000000000   F    (8c)
   0.19000000000000   0.19000000000000  -0.29000000000000   F    (8c)
   0.09000000000000   0.49000000000000   0.00000000000000   Tl   (8c)
  -0.09000000000000  -0.49000000000000   0.50000000000000   Tl   (8c)
  -0.49000000000000  -0.09000000000000   0.50000000000000   Tl   (8c)
   0.49000000000000   0.09000000000000   0.00000000000000   Tl   (8c)
```

$HoSb_2$: AB2_oC6_21_a_k - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'HoSb2'
_chemical_formula_sum 'Ho Sb2'

loop_
_publ_author_name
 'Q. Johnson'
_journal_name_full_name
;
 Inorganic Chemistry
;
_journal_volume 10
_journal_year 1971
_journal_page_first 2089
_journal_page_last 2090
_publ_Section_title
;
 The Crystal Structure of High-Pressure Synthesized Holmium Diantimonde
;

# Found in The Ho-Sb Alloy System, 1984

_aflow_title 'HoSb$_{2}$ Structure'
_aflow_proto 'AB2_oC6_21_a_k'
_aflow_params 'a,b/a,c/a,z_{2}'
_aflow_params_values '3.343,1.73197726593,2.34519892312,0.34'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oC6'

_symmetry_space_group_name_H-M "C 2 2 2"
_symmetry_Int_Tables_number 21

_cell_length_a      3.34300
_cell_length_b      5.79000
_cell_length_c      7.84000
_cell_angle_alpha   90.00000
_cell_angle_beta    90.00000
_cell_angle_gamma   90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 x+1/2,y+1/2,z
6 x+1/2,-y+1/2,-z
7 -x+1/2,y+1/2,-z
8 -x+1/2,-y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ho1 Ho 2 a 0.00000 0.00000 0.00000 1.00000
Sb1 Sb 4 k 0.25000 0.25000 0.34000 1.00000
```

$HoSb_2$: AB2_oC6_21_a_k - POSCAR

```
AB2_oC6_21_a_k & a,b/a,c/a,z2 --params=3.343,1.73197726593,2.34519892312
    ↪ ,0.34 & C222 D_{2}^{6} #21 (ak) & oC6 & None & HoSb2 & HoSb2 &
    ↪ Q. Johnson, Inorg. Chem. 10, 2089-2090 (1971)
   1.00000000000000
   1.67150000000000  -2.89500000000000   0.00000000000000
   1.67150000000000   2.89500000000000   0.00000000000000
   0.00000000000000   0.00000000000000   7.84000000000000
    Ho    Sb
     1     2
Direct
   0.00000000000000   0.00000000000000   0.00000000000000   Ho   (2a)
   0.00000000000000   0.50000000000000   0.34000000000000   Sb   (4k)
   0.50000000000000   0.00000000000000  -0.34000000000000   Sb   (4k)
```

Predicted Phase IV $Cd_2Re_2O_7$: A2B7C2_oF88_22_k_bdefghij_k - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Cd2O7Re2'
_chemical_formula_sum 'Cd2 O7 Re2'

loop_
_publ_author_name
 'K. J. Kapcia'
 'M. Reedyk'
 'M. Hajialamdari'
 'A. Ptok'
 'P. Piekarz'
 'F. S. Razavi'
 'A. M. Ole\\'{s}'
 'R. K. Kremer'
_journal_year 2019
_publ_Section_title
;
 Low-Temperature Phase of the Cd$_{2}$Re$_{2}$O$_{7}$ Superconductor: {\
    ↪ em Ab initio} Phonon Calculations and Raman Scattering
;

_aflow_title 'Predicted Phase IV Cd$_{2}$Re$_{2}$O$_{7}$ Structure'
_aflow_proto 'A2B7C2_oF88_22_k_bdefghij_k'
_aflow_params 'a,b/a,c/a,x_{3},y_{4},z_{5},z_{6},y_{7},x_{8},x_{9},y_{9
    ↪ },z_{9},x_{10},y_{10},z_{10}'
_aflow_params_values '10.3832,0.999961476231,1.00259072348,0.2,0.2,
    ↪ 0.1689,0.0447,0.4295,0.0706,0.63141,0.63137,0.6114,0.1245,
    ↪ 0.124502,0.126'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oF88'

_symmetry_space_group_name_H-M "F 2 2 2"
_symmetry_Int_Tables_number 22

_cell_length_a      10.38320
_cell_length_b      10.38280
_cell_length_c      10.41010
_cell_angle_alpha   90.00000
_cell_angle_beta    90.00000
_cell_angle_gamma   90.00000
```

```
loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 x,y+1/2,z+1/2
6 x,-y+1/2,-z+1/2
7 -x,y+1/2,-z+1/2
8 -x,-y+1/2,z+1/2
9 x+1/2,y,z+1/2
10 x+1/2,-y,-z+1/2
11 -x+1/2,y,-z+1/2
12 -x+1/2,-y,z+1/2
13 x+1/2,y+1/2,z
14 x+1/2,-y+1/2,-z
15 -x+1/2,y+1/2,-z
16 -x+1/2,-y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 4 b 0.000000 0.000000 0.500000 1.00000
O2 O 4 d 0.250000 0.250000 0.750000 1.00000
O3 O 8 e 0.200000 0.000000 0.000000 1.00000
O4 O 8 f 0.000000 0.200000 0.000000 1.00000
O5 O 8 g 0.000000 0.000000 0.168900 1.00000
O6 O 8 h 0.250000 0.250000 0.044700 1.00000
O7 O 8 i 0.250000 0.429500 0.250000 1.00000
O8 O 8 j 0.070600 0.250000 0.250000 1.00000
Cd1 Cd 16 k 0.631410 0.631370 0.611400 1.00000
Re1 Re 16 k 0.124500 0.124502 0.126000 1.00000
```

Predicted Phase IV $Cd_2Re_2O_7$: A2B7C2_oF88_22_k_bdefghij_k - POSCAR

```
A2B7C2_oF88_22_k_bdefghij_k & a,b/a,c/a,x3,y4,z5,z6,y7,x8,x9,y9,z9,x10,
    ↪ y10,z10 --params=10.3832,0.999961476231,1.00259072348,0.2,0.2,
    ↪ 0.1689,0.0447,0.4295,0.0706,0.63141,0.63137,0.6114,0.1245,
    ↪ 0.124502,0.126 & F222 D_{2}^{7} #22 (bdefghijk^2) & oF88 & None
    ↪ & Cd2O7Re2 & Cd2O7Re2 & K. J. Kapcia et al., (2019)
   1.00000000000000
   0.00000000000000   5.19140000000000   5.20505000000000
   5.19160000000000   0.00000000000000   5.20505000000000
   5.19160000000000   5.19140000000000   0.00000000000000
    Cd    O    Re
     4   14    4
Direct
   0.61136000000000   0.61144000000000   0.65138000000000  Cd  (16k)
   0.61144000000000   0.61136000000000  -1.87418000000000  Cd  (16k)
   0.65138000000000  -1.87418000000000   0.61136000000000  Cd  (16k)
  -1.87418000000000   0.65138000000000   0.61144000000000  Cd  (16k)
   0.50000000000000   0.50000000000000   0.50000000000000  O   (4b)
   0.75000000000000   0.75000000000000   0.75000000000000  O   (4d)
  -0.20000000000000   0.20000000000000   0.20000000000000  O   (8e)
   0.20000000000000  -0.20000000000000  -0.20000000000000  O   (8e)
   0.20000000000000  -0.20000000000000   0.20000000000000  O   (8f)
  -0.20000000000000   0.20000000000000  -0.20000000000000  O   (8f)
   0.16890000000000   0.16890000000000  -0.16890000000000  O   (8g)
  -0.16890000000000  -0.16890000000000   0.16890000000000  O   (8g)
   0.04470000000000   0.04470000000000   0.45530000000000  O   (8h)
   0.45530000000000   0.45530000000000   0.04470000000000  O   (8h)
   0.42950000000000   0.07050000000000   0.42950000000000  O   (8i)
   0.07050000000000   0.42950000000000   0.07050000000000  O   (8i)
   0.42940000000000   0.07060000000000   0.07060000000000  O   (8j)
   0.07060000000000   0.42940000000000   0.42940000000000  O   (8j)
   0.12600200000000   0.12599800000000   0.12300200000000  Re  (16k)
   0.12599800000000   0.12600200000000  -0.37500200000000  Re  (16k)
   0.12300200000000  -0.37500200000000   0.12600200000000  Re  (16k)
  -0.37500200000000   0.12300200000000   0.12599800000000  Re  (16k)
```

Mercury (II) Azide $[Hg(N_3)_2]$: AB6_oP28_29_a_6a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Mercury (ii) azide'
_chemical_formula_sum 'Hg N6'

loop_
_publ_author_name
 'U. M\"{u}ller'
_journal_name_full_name
;
 Zeitschrift fur Anorganische und Allgemeine Chemie
;
_journal_volume 399
_journal_year 1973
_journal_page_first 183
_journal_page_last 192
_publ_Section_title
;
 Die Kristallstruktur von $\alpha$-Quecksilber(II)-Azid
;

# Found in Binary Alloy Phase Diagrams, {Hf-Re to Zn-Zr}, 1990
_aflow_title 'Mercury (II) Azide [Hg(N$_{3}$)$_{2}$] Structure'
_aflow_proto 'AB6_oP28_29_a_6a'
_aflow_params 'a,b/a,c/a,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3},y_{3
    ↪ },z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{6},
    ↪ x_{7},y_{7},z_{7}'
_aflow_params_values '10.632,0.589164785553,0.59471407073,0.03183,
    ↪ 0.23877,0.25,0.082,-0.043,0.106,0.188,0.893,0.121,0.284,0.82,
    ↪ 0.112,0.962,0.529,0.39,0.883,0.609,0.28,0.807,0.7,0.172'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP28'

_symmetry_space_group_name_H-M "P c a 21"
_symmetry_Int_Tables_number 29

_cell_length_a     10.63200
_cell_length_b     6.26400
_cell_length_c     6.32300
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,z+1/2
3 -x+1/2,y,z+1/2
4 x+1/2,-y,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Hg1 Hg 4 a 0.03183 0.23877 0.25000 1.00000
N1 N 4 a 0.08200 -0.04300 0.10600 1.00000
N2 N 4 a 0.18800 0.89300 0.12100 1.00000
N3 N 4 a 0.28400 0.82000 0.11200 1.00000
N4 N 4 a 0.96200 0.52900 0.39000 1.00000
N5 N 4 a 0.88300 0.60900 0.28000 1.00000
N6 N 4 a 0.80700 0.70000 0.17200 1.00000
```

Mercury (II) Azide $[Hg(N_3)_2]$: AB6_oP28_29_a_6a - POSCAR

```
AB6_oP28_29_a_6a & a,b/a,c/a,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,z4,x5,y5,
    ↪ z5,x6,y6,z6,x7,y7,z7 --params=10.632,0.589164785553,
    ↪ 0.59471407073,0.03183,0.23877,0.25,0.082,-0.043,0.106,0.188,
    ↪ 0.893,0.121,0.284,0.82,0.112,0.962,0.529,0.39,0.883,0.609,0.28,
    ↪ 0.807,0.7,0.172 & Pca2_{1} C_{2v}^{5} #29 (a^7) & oP28 & None &
    ↪  HgN6 & Mercury (ii) azide & U. M\"{u}ller, Z. Anorg. Allg.
    ↪ Chem. 399, 183-192 (1973)
   1.00000000000000
  10.63200000000000   0.00000000000000   0.00000000000000
   0.00000000000000   6.26400000000000   0.00000000000000
   0.00000000000000   0.00000000000000   6.32300000000000
    Hg    N
     4   24
Direct
   0.03183000000000   0.23877000000000   0.25000000000000  Hg   (4a)
  -0.03183000000000  -0.23877000000000   0.75000000000000  Hg   (4a)
   0.53183000000000  -0.23877000000000   0.25000000000000  Hg   (4a)
   0.46817000000000   0.23877000000000   0.75000000000000  Hg   (4a)
   0.08200000000000  -0.04300000000000   0.10600000000000  N   (4a)
  -0.08200000000000   0.04300000000000   0.60600000000000  N   (4a)
   0.58200000000000   0.04300000000000   0.10600000000000  N   (4a)
   0.41800000000000  -0.04300000000000   0.60600000000000  N   (4a)
   0.18800000000000   0.89300000000000   0.12100000000000  N   (4a)
  -0.18800000000000  -0.89300000000000   0.62100000000000  N   (4a)
   0.68800000000000  -0.89300000000000   0.12100000000000  N   (4a)
   0.31200000000000   0.89300000000000   0.62100000000000  N   (4a)
   0.28400000000000   0.82000000000000   0.11200000000000  N   (4a)
  -0.28400000000000  -0.82000000000000   0.61200000000000  N   (4a)
   0.78400000000000  -0.82000000000000   0.11200000000000  N   (4a)
   0.21600000000000   0.82000000000000   0.61200000000000  N   (4a)
   0.96200000000000   0.52900000000000   0.39000000000000  N   (4a)
  -0.96200000000000  -0.52900000000000   0.89000000000000  N   (4a)
   1.46200000000000  -0.52900000000000   0.39000000000000  N   (4a)
  -0.46200000000000   0.52900000000000   0.89000000000000  N   (4a)
   0.88300000000000   0.60900000000000   0.28000000000000  N   (4a)
  -0.88300000000000  -0.60900000000000   0.78000000000000  N   (4a)
   1.38300000000000  -0.60900000000000   0.28000000000000  N   (4a)
  -0.38300000000000   0.60900000000000   0.78000000000000  N   (4a)
   0.80700000000000   0.70000000000000   0.17200000000000  N   (4a)
  -0.80700000000000  -0.70000000000000   0.67200000000000  N   (4a)
   1.30700000000000  -0.70000000000000   0.17200000000000  N   (4a)
  -0.30700000000000   0.70000000000000   0.67200000000000  N   (4a)
```

Low-Temperature $(NH_3CH_3)Al(SO_4)_2{\cdot}12H_2O$: ABC30DE20F2_oP220_29_a_a_30a_a_20a_2a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'AlCH30NO20S2'
_chemical_formula_sum 'Al C H30 N O20 S2'

loop_
_publ_author_name
 'R. O. W. Fletcher'
 'H. Steeple'
_journal_name_full_name
```

```
;
 Acta Cristallographica
;
_journal_volume 17
_journal_year 1964
_journal_page_first 290
_journal_page_last 294
_publ_Section_title
;
 The crystal structure of the low-temperature phase of methylammonium
       ↪ alum
;

_aflow_title 'Low-Temperature (NH$_{3}$CH$_{3}$)Al(SO$_{4}$)$_{2}$$\
      ↪ cdot$12H$_{2}$O Structure'
_aflow_proto 'ABC30DE20F2_oP220_29_a_a_30a_a_20a_2a'
_aflow_params 'a,b/a,c/a,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3},y_{3
      ↪ },z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{6},
      ↪ x_{7},y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{10},y_
      ↪ {10},z_{10},x_{11},y_{11},z_{11},x_{12},y_{12},z_{12},x_{13},y_
      ↪ {13},z_{13},x_{14},y_{14},z_{14},x_{15},y_{15},z_{15},x_{16},y_
      ↪ {16},z_{16},x_{17},y_{17},z_{17},x_{18},y_{18},z_{18},x_{19},y_
      ↪ {19},z_{19},x_{20},y_{20},z_{20},x_{21},y_{21},z_{21},x_{22},y_
      ↪ {22},z_{22},x_{23},y_{23},z_{23},x_{24},y_{24},z_{24},x_{25},y_
      ↪ {25},z_{25},x_{26},y_{26},z_{26},x_{27},y_{27},z_{27},x_{28},y_
      ↪ {28},z_{28},x_{29},y_{29},z_{29},x_{30},y_{30},z_{30},x_{31},y_
      ↪ {31},z_{31},x_{32},y_{32},z_{32},x_{33},y_{33},z_{33},x_{34},y_
      ↪ {34},z_{34},x_{35},y_{35},z_{35},x_{36},y_{36},z_{36},x_{37},y_
      ↪ {37},z_{37},x_{38},y_{38},z_{38},x_{39},y_{39},z_{39},x_{40},y_
      ↪ {40},z_{40},x_{41},y_{41},z_{41},x_{42},y_{42},z_{42},x_{43},y_
      ↪ {43},z_{43},x_{44},y_{44},z_{44},x_{45},y_{45},z_{45},x_{46},y_
      ↪ {46},z_{46},x_{47},y_{47},z_{47},x_{48},y_{48},z_{48},x_{49},y_
      ↪ {49},z_{49},x_{50},y_{50},z_{50},x_{51},y_{51},z_{51},x_{52},y_
      ↪ {52},z_{52},x_{53},y_{53},z_{53},x_{54},y_{54},z_{54},x_{55},y_
      ↪ {55},z_{55}'
_aflow_params_values '12.57,0.980906921241,0.984884645982,0.006,0.259,
      ↪ 0.256,0.047,0.77,0.28,0.21,0.237,0.305,0.199,0.277,0.175,0.196,
      ↪ 0.722,0.693,0.193,0.774,0.825,0.076,0.057,0.235,0.048,0.953,
      ↪ 0.781,0.07,0.455,0.22,0.059,0.542,0.768,0.026,0.18,0.45,0.024,
      ↪ 0.698,0.96,0.025,0.192,0.048,0.021,0.68,0.547,0.006,0.012,0.502
      ↪ ,0.136,0.06,0.507,0.247,0.385,0.557,0.243,0.264,0.012,0.225,
      ↪ 0.488,0.752,0.183,0.48,0.885,0.004,0.503,0.502,0.139,0.552,0.49
      ↪ ,0.249,0.757,0.49,0.244,0.893,0.445,0.249,0.01,0.743,0.202,
      ↪ 0.002,0.108,0.12,0.784,0.34,0.033,0.84,0.245,0.083,0.71,0.225,
      ↪ 0.119,0.275,0.665,0.017,0.33,0.755,0.08,0.205,0.775,0.045,0.268
      ↪ ,0.711,0.136,0.04,0.811,0.24,0.022,0.978,0.064,0.096,0.976,
      ↪ 0.199,0.202,0.904,0.169,0.528,0.19,0.225,0.51,0.011,0.061,0.61,
      ↪ 0.046,0.231,0.684,0.097,0.153,0.257,0.236,0.145,0.748,0.759,
      ↪ 0.012,0.105,0.258,0.003,0.405,0.247,0.007,0.242,0.397,0.0,0.256
      ↪ ,0.098,0.228,0.302,0.591,0.154,0.456,0.801,0.235,0.808,0.425,
      ↪ 0.174,0.988,0.182,0.057,0.086,0.521,0.053,0.577,0.472,0.163,
      ↪ 0.088,0.915,0.171,0.581,0.085'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP220'

_symmetry_space_group_name_H-M "P c a 21"
_symmetry_Int_Tables_number 29

_cell_length_a      12.57000
_cell_length_b      12.33000
_cell_length_c      12.38000
_cell_angle_alpha   90.00000
_cell_angle_beta    90.00000
_cell_angle_gamma   90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,z+1/2
3 -x+1/2,y,z+1/2
4 x+1/2,-y,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Al1 Al 4 a 0.00600 0.25900 0.25600 1.00000
C1 C 4 a 0.04700 0.77000 0.28000 1.00000
H1 H 4 a 0.21000 0.23700 0.30500 1.00000
H2 H 4 a 0.19900 0.27700 0.17500 1.00000
H3 H 4 a 0.19600 0.72200 0.69300 1.00000
H4 H 4 a 0.19300 0.77400 0.82500 1.00000
H5 H 4 a 0.07600 0.05700 0.23500 1.00000
H6 H 4 a 0.04800 0.95300 0.78100 1.00000
H7 H 4 a 0.07000 0.45500 0.22000 1.00000
H8 H 4 a 0.05900 0.54200 0.76800 1.00000
H9 H 4 a 0.02600 0.18000 0.45000 1.00000
H10 H 4 a 0.02400 0.69800 0.96000 1.00000
H11 H 4 a 0.02500 0.19200 0.04800 1.00000
H12 H 4 a 0.02100 0.68000 0.54700 1.00000
H13 H 4 a 0.00600 0.01200 0.50200 1.00000
H14 H 4 a 0.13600 0.06000 0.50700 1.00000
H15 H 4 a 0.24700 0.38500 0.55700 1.00000
H16 H 4 a 0.24300 0.26400 0.01200 1.00000
H17 H 4 a 0.22500 0.48800 0.75200 1.00000
H18 H 4 a 0.18300 0.48000 0.88500 1.00000
H19 H 4 a 0.00400 0.50300 0.50200 1.00000
H20 H 4 a 0.13900 0.55200 0.49000 1.00000
H21 H 4 a 0.24900 0.75700 0.49000 1.00000
H22 H 4 a 0.24400 0.89300 0.44500 1.00000
H23 H 4 a 0.24900 0.01000 0.74300 1.00000
H24 H 4 a 0.20200 0.00200 0.10800 1.00000
H25 H 4 a 0.12000 0.78400 0.34000 1.00000
H26 H 4 a 0.03300 0.84000 0.24500 1.00000
H27 H 4 a 0.08300 0.71000 0.22500 1.00000
H28 H 4 a 0.11900 0.27500 0.66500 1.00000
H29 H 4 a 0.01700 0.33000 0.75500 1.00000
H30 H 4 a 0.08000 0.20500 0.77500 1.00000
N1 N 4 a 0.04500 0.26800 0.71100 1.00000
O1 O 4 a 0.13600 0.04000 0.81100 1.00000
O2 O 4 a 0.24000 0.02200 0.97800 1.00000
O3 O 4 a 0.06400 0.09600 0.97600 1.00000
O4 O 4 a 0.19900 0.20200 0.90400 1.00000
O5 O 4 a 0.16900 0.52800 0.19000 1.00000
O6 O 4 a 0.22500 0.51000 0.01100 1.00000
O7 O 4 a 0.06100 0.61000 0.04600 1.00000
O8 O 4 a 0.23100 0.68400 0.09700 1.00000
O9 O 4 a 0.15300 0.25700 0.23600 1.00000
O10 O 4 a 0.14500 0.74800 0.75900 1.00000
O11 O 4 a 0.01200 0.10500 0.25800 1.00000
O12 O 4 a 0.00300 0.40500 0.24700 1.00000
O13 O 4 a 0.00700 0.24200 0.39700 1.00000
O14 O 4 a 0.00000 0.25600 0.09800 1.00000
O15 O 4 a 0.22800 0.30200 0.59100 1.00000
O16 O 4 a 0.15400 0.45600 0.80100 1.00000
O17 O 4 a 0.23500 0.80800 0.42500 1.00000
O18 O 4 a 0.17400 0.98800 0.18200 1.00000
O19 O 4 a 0.05700 0.08600 0.52100 1.00000
O20 O 4 a 0.05300 0.57700 0.47200 1.00000
S1 S 4 a 0.16300 0.08800 0.91500 1.00000
S2 S 4 a 0.17100 0.58100 0.08500 1.00000
```

Low-Temperature $(NH_3CH_3)Al(SO_4)_2 \cdot 12H_2O$: ABC30DE20F2_oP220_29_a_a_30a_a_20a_2a - POSCAR

```
ABC30DE20F2_oP220_29_a_a_30a_a_20a_2a & a,b/a,c/a,x1,y1,z1,x2,y2,z2,x3,
      ↪ y3,z3,x4,y4,z4,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10
      ↪ ,y10,z10,x11,y11,z11,x12,y12,z12,x13,y13,z13,x14,y14,z14,x15,
      ↪ y15,z15,x16,y16,z16,x17,y17,z17,x18,y18,z18,x19,y19,z19,x20,y20
      ↪ ,z20,x21,y21,z21,x22,y22,z22,x23,y23,z23,x24,y24,z24,x25,y25,
      ↪ z25,x26,y26,z26,x27,y27,z27,x28,y28,z28,x29,y29,z29,x30,y30,z30
      ↪ ,x31,y31,z31,x32,y32,z32,x33,y33,z33,x34,y34,z34,x35,y35,z35,
      ↪ x36,y36,z36,x37,y37,z37,x38,y38,z38,x39,y39,z39,x40,y40,z40,x41
      ↪ ,y41,z41,x42,y42,z42,x43,y43,z43,x44,y44,z44,x45,y45,z45,x46,
      ↪ y46,z46,x47,y47,z47,x48,y48,z48,x49,y49,z49,x50,y50,z50,x51,y51
      ↪ ,z51,x52,y52,z52,x53,y53,z53,x54,y54,z54,x55,y55,z55 --params=
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      ↪ 0.025,0.192,0.048,0.021,0.68,0.547,0.006,0.012,0.502,0.136,0.06
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      ↪ ,0.784,0.34,0.033,0.84,0.245,0.083,0.71,0.225,0.119,0.275,0.665
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      ↪ 0.581,0.085 & Pca2_{1} C_{2v}^{5} #29 (a^55) & oP220 & None &
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      ↪ Acta Cryst. 17, 290-294 (1964)
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Orthorhombic Co_4Al_{13}: A13B4_oP102_31_17a11b_8a2b - CIF

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      ↪ 0.2291,0.422,0.4252,0.2133,0.5174,0.7338,0.2319,0.2763,0.2849,
      ↪ 0.2257,-0.0901,0.7346,0.2181,0.5969,0.0984'
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_aflow_Pearson 'oP102'

_symmetry_space_group_name_H-M "P m n 21"
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2 -x+1/2,-y,z+1/2
3 -x,y,z
4 x+1/2,-y,z+1/2

loop_
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_atom_site_symmetry_multiplicity
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_atom_site_occupancy
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Al3  Al  2  a  0.00000  -0.09470  0.62260  1.00000
Al4  Al  2  a  0.00000  -0.09150  0.44310  1.00000
Al5  Al  2  a  0.00000  0.13940  0.67280  1.00000
Al6  Al  2  a  0.00000  0.81130  0.15640  1.00000
Al7  Al  2  a  0.00000  0.80920  0.83600  1.00000
Al8  Al  2  a  0.00000  0.25200  0.41400  1.00000
Al9  Al  2  a  0.00000  0.40700  0.85200  1.00000
Al10  Al  2  a  0.00000  0.40100  0.14910  1.00000
Al11  Al  2  a  0.00000  0.68370  0.67490  1.00000
Al12  Al  2  a  0.00000  0.59000  0.39600  1.00000
Al13  Al  2  a  0.00000  0.52710  0.00540  1.00000
Al14  Al  2  a  0.00000  0.59860  0.21490  1.00000
Al15  Al  2  a  0.00000  0.71110  -0.03320  1.00000
Al16  Al  2  a  0.00000  0.10470  0.23800  1.00000
Al17  Al  2  a  0.00000  0.41790  0.58810  1.00000
Co1  Co  2  a  0.00000  0.89780  0.00000  1.00000
Co2  Co  2  a  0.00000  0.09010  0.51140  1.00000
Co3  Co  2  a  0.00000  0.19770  0.82500  1.00000
Co4  Co  2  a  0.00000  0.77200  0.31670  1.00000
Co5  Co  2  a  0.00000  0.59860  0.82480  1.00000
Co6  Co  2  a  0.00000  0.41190  0.31370  1.00000
Co7  Co  2  a  0.00000  0.73090  0.51800  1.00000
Co8  Co  2  a  0.00000  0.28770  0.00760  1.00000
Al18  Al  4  b  0.21350  0.21380  0.09960  1.00000
Al19  Al  4  b  0.21380  -0.09800  0.28310  1.00000
Al20  Al  4  b  0.24010  0.08290  0.40880  1.00000
Al21  Al  4  b  0.25340  0.25500  -0.08280  1.00000
Al22  Al  4  b  0.22510  -0.01900  0.09350  1.00000
Al23  Al  4  b  0.22140  0.22290  0.55000  1.00000
Al24  Al  4  b  0.22590  0.58850  0.54220  1.00000
Al25  Al  4  b  0.22440  0.29460  0.73520  1.00000
Al26  Al  4  b  0.22910  0.42200  0.42520  1.00000
Al27  Al  4  b  0.21330  0.51740  0.73380  1.00000
Al28  Al  4  b  0.23190  0.27630  0.28490  1.00000
Co9  Co  4  b  0.22570  -0.09010  0.73460  1.00000
Co10  Co  4  b  0.21810  0.59690  0.09840  1.00000
```

Orthorhombic Co_4Al_{13}: A13B4_oP102_31_17a11b_8a2b - POSCAR

```
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      ↪ z6,y7,z7,y8,z8,y9,z9,y10,z10,y11,z11,y12,z12,y13,z13,y14,z14,
      ↪ y15,z15,y16,z16,y17,z17,y18,z18,y19,z19,y20,z20,y21,z21,y22,z22
      ↪ ,y23,z23,y24,z24,y25,z25,x26,y26,z26,x27,y27,z27,x28,y28,z28,
      ↪ x29,y29,z29,x30,y30,z30,x31,y31,z31,x32,y32,z32,x33,y33,z33,x34
      ↪ ,y34,z34,x35,y35,z35,x36,y36,z36,x37,y37,z37,x38,y38,z38 --
      ↪ params=8.158,1.51287080167,1.77151262564,-0.0055,0.8151,0.0914
      ↪ ,-0.0431,-0.0947,0.6226,-0.0915,0.4431,0.1394,0.6728,0.8113,
      ↪ 0.1564,0.8092,0.836,0.252,0.414,0.407,0.852,0.401,0.1491,0.6837
      ↪ ,0.6749,0.59,0.396,0.5271,0.0054,0.5986,0.2149,0.7111,-0.0332,
      ↪ 0.1047,0.238,0.4179,0.5881,0.8978,0.0,0.0901,0.5114,0.1977,
      ↪ 0.825,0.772,0.3167,0.5986,0.8248,0.4119,0.3137,0.7309,0.518,
      ↪ 0.2877,0.0076,0.2135,0.2138,0.0996,0.2138,-0.098,0.2831,0.2401,
      ↪ 0.0829,0.4088,0.2534,0.255,-0.0828,0.2251,-0.019,0.0935,0.2214,
      ↪ 0.2229,0.55,0.2259,0.5885,0.5422,0.2244,0.2946,0.7352,0.2291,
      ↪ 0.422,0.4252,0.2133,0.5174,0.7338,0.2319,0.2763,0.2849,0.2257,-
      ↪ 0.0901,0.7346,0.2181,0.5969,0.0984 & Pmn2_{1} C_{2v}^{7} #31 (a
      ↪ ^25b^13) & oP102 & None & Al13Co4 & Al13Co4 & J. Grin et al.,
      ↪ J. Alloys Compd. 206, 243-247 (1994)
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  8.15800000000000     0.00000000000000     0.00000000000000
  0.00000000000000    12.34200000000000     0.00000000000000
  0.00000000000000     0.00000000000000    14.45200000000000
   Al   Co
   78   24
Direct
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  0.23190000000000   0.27630000000000   0.28490000000000   Al   (4b)
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  0.00000000000000   0.09010000000000   0.51140000000000   Co   (2a)
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  0.00000000000000   0.28770000000000   0.00760000000000  Co   (2a)
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 −0.22570000000000  −0.09010000000000   0.73460000000000  Co   (4b)
  0.21810000000000   0.59690000000000   0.09840000000000  Co   (4b)
  0.28190000000000  −0.59690000000000   0.59840000000000  Co   (4b)
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 −0.21810000000000   0.59690000000000   0.09840000000000  Co   (4b)
```

$Mg(ClO_4)_2{\cdot}6H_2O$ ($H4_{11}$): A2B6CD8_oP34_31_2a_2a2b_a_4a2b - CIF

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 'C. D. West'
_journal_name_full_name
;
 Zeitschrift f\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 91
_journal_year 1935
_journal_page_first 480
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 Crystal Structures of Hydrated Compounds II. Structure Type Mg(ClO$_{4}
      ↪ $)$_{2}$$\cdot$6H$_{2}$O
;

# Found in Stability of phases in the Mg(ClO$_{4}$)$_{2}$$\cdot$n$H$_{2
      ↪ }$O system and implications for perchlorate occurrences on Mars
      ↪ , 2011

_aflow_title 'Mg(ClO$_{4}$)$_{2}$$\cdot$6H$_{2}$O ($H4_{11}$) Structure'
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Cl2 Cl 2 a 0.00000 0.58333 0.00000 1.00000
H2O1 H2O 2 a 0.00000 0.12500 0.00000 1.00000
H2O2 H2O 2 a 0.00000 0.37500 0.50000 1.00000
Mg1 Mg 2 a 0.00000 0.25000 0.75000 1.00000
O1 O 2 a 0.00000 -0.08333 0.77778 1.00000
O2 O 2 a 0.00000 0.58333 0.27778 1.00000
O3 O 2 a 0.00000 0.81111 0.40833 1.00000
O4 O 2 a 0.00000 0.68889 -0.09167 1.00000
H2O3 H2O 4 b 0.81250 0.18750 0.50000 1.00000
H2O4 H2O 4 b 0.81250 0.31250 0.00000 1.00000
O5 O 4 b 0.84167 -0.03056 0.40833 1.00000
O6 O 4 b 0.84167 0.53056 -0.09167 1.00000
```

$Mg(ClO_4)_2{\cdot}6H_2O$ ($H4_{11}$): A2B6CD8_oP34_31_2a_2a2b_a_4a2b - POSCAR

```
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      ↪ ,y6,z6,y7,z7,y8,z8,y9,z9,x10,y10,z10,x11,y11,z11,x12,y12,z12,
      ↪ x13,y13,z13 --params=7.76,1.73453608247,0.677835051546,-0.08333
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      ↪ ,0.58333,0.27778,0.81111,0.40833,0.68889,-0.09167,0.8125,0.1875
      ↪ ,0.5,0.8125,0.3125,0.0,0.84167,-0.03056,0.40833,0.84167,0.53056
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      ↪  & Cl2(H2O)6MgO8 & Cl2(H2O)6MgO8 & C. D. West, Zeitschrift f{"u
      ↪ }r Kristallographie - Crystalline Materials 91, 480-493 (1935)
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   0.00000000000000    13.46000000000000     0.00000000000000
   0.00000000000000     0.00000000000000     5.26000000000000
     Cl    H2O     Mg      O
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Direct
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   0.50000000000000  -0.58333000000000   0.77778000000000   O    (2a)
   0.00000000000000   0.81111000000000   0.40833000000000   O    (2a)
   0.50000000000000  -0.81111000000000   0.90833000000000   O    (2a)
   0.00000000000000   0.68889000000000  -0.09167000000000   O    (2a)
   0.50000000000000  -0.68889000000000   0.40833000000000   O    (2a)
   0.84167000000000  -0.03056000000000   0.40833000000000   O    (4b)
  -0.34167000000000   0.03056000000000   0.90833000000000   O    (4b)
   1.34167000000000   0.03056000000000   0.90833000000000   O    (4b)
  -0.84167000000000  -0.03056000000000   0.40833000000000   O    (4b)
   0.84167000000000   0.53056000000000  -0.09167000000000   O    (4b)
  -0.34167000000000  -0.53056000000000   0.40833000000000   O    (4b)
   1.34167000000000  -0.53056000000000   0.40833000000000   O    (4b)
  -0.84167000000000   0.53056000000000  -0.09167000000000   O    (4b)
```

B_4SrO_7: A4B7C_oP24_31_2b_a3b_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'B4O7Sr'
_chemical_formula_sum 'B4 O7 Sr'

loop_
_publ_author_name
 'J. {Krogh-Moe}'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 18
_journal_year 1964
_journal_page_first 2055
_journal_page_last 2060
_publ_Section_title
;
 The Crystal Structure of Strontium Diborate, SrO$\cdot$2B$_{2}$O$_{3}$
;

_aflow_title 'B$_{4}$SrO$_{7}$ Structure'
_aflow_proto 'A4B7C_oP24_31_2b_a3b_a'
_aflow_params 'a,b/a,c/a,y_{1},z_{1},y_{2},z_{2},x_{3},y_{3},z_{3},x_{4
      ↪ },y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{6},x_{7},y_{7},
      ↪ z_{7}'
_aflow_params_values '10.711,0.4133134161l4,0.395387918962,0.728,0.454,
      ↪ 0.289,0.0,0.379,0.174,-0.024,0.246,0.671,-0.037,0.359,0.857,
      ↪ 0.064,0.221,0.631,0.335,0.365,0.226,0.335'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP24'

_symmetry_space_group_name_H-M "P m n 21"
_symmetry_Int_Tables_number 31

_cell_length_a     10.71100
_cell_length_b     4.42700
_cell_length_c     4.23500
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
```

```
2 -x+1/2,-y,z+1/2
3 -x,y,z
4 x+1/2,-y,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 2 a 0.00000 0.72800 0.45400 1.00000
Sr1 Sr 2 a 0.00000 0.28900 0.00000 1.00000
B1 B 4 b 0.37900 0.17400 -0.02400 1.00000
B2 B 4 b 0.24600 0.67100 -0.03700 1.00000
O2 O 4 b 0.35900 0.85700 0.06400 1.00000
O3 O 4 b 0.22100 0.63100 0.33500 1.00000
O4 O 4 b 0.36500 0.22600 0.33500 1.00000
```

B_4SrO_7: A4B7C_oP24_31_2b_a3b_a - POSCAR

```
A4B7C_oP24_31_2b_a3b_a & a,b/a,c/a,y1,z1,y2,z2,x3,y3,z3,x4,y4,z4,x5,y5,
    ↪ z5,x6,y6,z6,x7,y7,z7 --params=10.711,0.413313416114,
    ↪ 0.395387918962,0.728,0.454,0.289,0.0,0.379,0.174,-0.024,0.246,
    ↪ 0.671,-0.037,0.359,0.857,0.064,0.221,0.631,0.335,0.365,0.226,
    ↪ 0.335 & Pmn2_{1} C_{2v}^{7} #31 (a^2b^5) & oP24 & None & B4O7Sr
    ↪ & B4O7Sr & J. {Krogh-Moe}, Acta Chem. Scand. 18, 2055-2060 (
    ↪ 1964)
   1.00000000000000
  10.71100000000000    0.00000000000000    0.00000000000000
   0.00000000000000    4.42700000000000    0.00000000000000
   0.00000000000000    0.00000000000000    4.23500000000000
     B     O     Sr
     8    14     2
Direct
   0.37900000000000    0.17400000000000   -0.02400000000000    B    (4b)
   0.12100000000000   -0.17400000000000    0.47600000000000    B    (4b)
   0.87900000000000   -0.17400000000000    0.47600000000000    B    (4b)
  -0.37900000000000    0.17400000000000   -0.02400000000000    B    (4b)
   0.24600000000000    0.67100000000000   -0.03700000000000    B    (4b)
   0.25400000000000   -0.67100000000000    0.46300000000000    B    (4b)
   0.74600000000000   -0.67100000000000    0.46300000000000    B    (4b)
  -0.24600000000000    0.67100000000000   -0.03700000000000    B    (4b)
   0.00000000000000    0.72800000000000    0.45400000000000    O    (2a)
   0.50000000000000   -0.72800000000000    0.95400000000000    O    (2a)
   0.35900000000000    0.85700000000000    0.06400000000000    O    (4b)
   0.14100000000000   -0.85700000000000    0.56400000000000    O    (4b)
   0.85900000000000   -0.85700000000000    0.56400000000000    O    (4b)
  -0.35900000000000    0.85700000000000    0.06400000000000    O    (4b)
   0.22100000000000    0.63100000000000    0.33500000000000    O    (4b)
   0.27900000000000   -0.63100000000000    0.83500000000000    O    (4b)
   0.72100000000000   -0.63100000000000    0.83500000000000    O    (4b)
  -0.22100000000000    0.63100000000000    0.33500000000000    O    (4b)
   0.36500000000000    0.22600000000000    0.33500000000000    O    (4b)
   0.13500000000000   -0.22600000000000    0.83500000000000    O    (4b)
   0.86500000000000   -0.22600000000000    0.83500000000000    O    (4b)
  -0.36500000000000    0.22600000000000    0.33500000000000    O    (4b)
   0.00000000000000    0.28900000000000    0.00000000000000   Sr    (2a)
   0.50000000000000   -0.28900000000000    0.50000000000000   Sr    (2a)
```

$D8_7$ (Shcherbinaite, V_2O_5) (*obsolete*): A5B2_oP14_31_a2b_b - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Shcherbinaite'
_chemical_formula_sum 'O5 V2'

loop_
_publ_author_name
 'J. A. A. Ketelaar'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 95
_journal_year 1936
_journal_page_first 9
_journal_page_last 27
_publ_Section_title
;
 Die Kristallstruktur des Vanadinpentoxyds
;

# Found in A Refinement of the Structure of V$_{2}$O$_{5}$, 1986

_aflow_title '$D8_{7}$ (Shcherbinaite, V$_{2}$O$_{5}$) ({\em{obsolete}})
    ↪ Structure'
_aflow_proto 'A5B2_oP14_31_a2b_b'
_aflow_params 'a,b/a,c/a,y_{1},z_{1},x_{2},y_{2},z_{2},x_{3},y_{3},z_{3
    ↪ },x_{4},y_{4},z_{4}'
_aflow_params_values '11.48,0.379790940767,0.309233449477,0.08,0.89,
    ↪ 0.148,0.45,-0.08,0.2,0.03,0.46,0.148,0.097,0.0'
_aflow_Strukturbericht '$D8_{7}$'
_aflow_Pearson 'oP14'

_symmetry_space_group_name_H-M "P m n 21"
_symmetry_Int_Tables_number 31

_cell_length_a    11.48000
_cell_length_b    4.36000
_cell_length_c    3.55000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x+1/2,-y,z+1/2
3 -x,y,z
4 x+1/2,-y,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 2 a 0.00000 0.08000 0.89000 1.00000
O2 O 4 b 0.14800 0.45000 -0.08000 1.00000
O3 O 4 b 0.20000 0.03000 0.46000 1.00000
V1 V 4 b 0.14800 0.09700 0.00000 1.00000
```

$D8_7$ (Shcherbinaite, V_2O_5) (*obsolete*): A5B2_oP14_31_a2b_b - POSCAR

```
A5B2_oP14_31_a2b_b & a,b/a,c/a,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,z4 --params
    ↪ =11.48,0.379790940767,0.309233449477,0.08,0.89,0.148,0.45,-0.08
    ↪ ,0.2,0.03,0.46,0.148,0.097,0.0 & Pmn2_{1} C_{2v}^{7} #31 (ab^3)
    ↪ & oP14 & $D8_{7}$ & O5V2 & Shcherbinaite & J. A. A. Ketelaar,
    ↪ Zeitschrift f{"u}r Kristallographie - Crystalline Materials 95,
    ↪ 9-27 (1936)
   1.00000000000000
  11.48000000000000    0.00000000000000    0.00000000000000
   0.00000000000000    4.36000000000000    0.00000000000000
   0.00000000000000    0.00000000000000    3.55000000000000
     O     V
    10     4
Direct
   0.00000000000000    0.08000000000000    0.89000000000000    O    (2a)
   0.50000000000000   -0.08000000000000    1.39000000000000    O    (2a)
   0.14800000000000    0.45000000000000   -0.08000000000000    O    (4b)
   0.35200000000000   -0.45000000000000    0.42000000000000    O    (4b)
   0.64800000000000   -0.45000000000000    0.42000000000000    O    (4b)
  -0.14800000000000    0.45000000000000   -0.08000000000000    O    (4b)
   0.20000000000000    0.03000000000000    0.46000000000000    O    (4b)
   0.30000000000000   -0.03000000000000    0.96000000000000    O    (4b)
   0.70000000000000   -0.03000000000000    0.96000000000000    O    (4b)
  -0.20000000000000    0.03000000000000    0.46000000000000    O    (4b)
   0.14800000000000    0.09700000000000    0.00000000000000    V    (4b)
   0.35200000000000   -0.09700000000000    0.50000000000000    V    (4b)
   0.64800000000000   -0.09700000000000    0.50000000000000    V    (4b)
  -0.14800000000000    0.09700000000000    0.00000000000000    V    (4b)
```

$Mo_{17}O_{47}$: A17B47_oP128_32_a8c_a23c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Mo17O47'
_chemical_formula_sum 'Mo17 O47'

loop_
_publ_author_name
 'L. Kihlborg'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 17
_journal_year 1963
_journal_page_first 1485
_journal_page_last 1487
_publ_Section_title
;
 Least Squares Refinement of the Structure of Mo$_{17}$O$_{47}$
;

_aflow_title 'Mo$_{17}$O$_{47}$ Structure'
_aflow_proto 'A17B47_oP128_32_a8c_a23c'
_aflow_params 'a,b/a,c/a,z_{1},z_{2},x_{3},y_{3},z_{3},x_{4},y_{4},z_{4
    ↪ },x_{5},y_{5},z_{5},x_{6},y_{6},z_{6},x_{7},y_{7},z_{7},x_{8},
    ↪ y_{8},z_{8},x_{9},y_{9},z_{9},x_{10},y_{10},z_{10},x_{11},y_{11
    ↪ },z_{11},x_{12},y_{12},z_{12},x_{13},y_{13},z_{13},x_{14},y_{14
    ↪ },z_{14},x_{15},y_{15},z_{15},x_{16},y_{16},z_{16},x_{17},y_{17
    ↪ },z_{17},x_{18},y_{18},z_{18},x_{19},y_{19},z_{19},x_{20},y_{20
    ↪ },z_{20},x_{21},y_{21},z_{21},x_{22},y_{22},z_{22},x_{23},y_{23
    ↪ },z_{23},x_{24},y_{24},z_{24},x_{25},y_{25},z_{25},x_{26},y_{26
    ↪ },z_{26},x_{27},y_{27},z_{27},x_{28},y_{28},z_{28},x_{29},y_{29
    ↪ },z_{29},x_{30},y_{30},z_{30},x_{31},y_{31},z_{31},x_{32},y_{32
    ↪ },z_{32},x_{33},y_{33},z_{33}'
_aflow_params_values '21.61,0.908375751967,0.182832022212,0.579,0.033,
    ↪ 0.02398,0.26126,0.586,0.12899,0.11845,0.4177,0.13648,0.39897,
    ↪ 0.4329,0.2426,0.25684,0.579,0.28825,0.06556,0.5848,0.38285,
    ↪ 0.19333,0.4334,0.38519,0.36637,0.4205,0.46561,0.05514,0.5551,
    ↪ 0.0225,0.263,0.027,0.131,0.1206,-0.052,0.1349,0.4038,-0.011,
    ↪ 0.2465,0.2532,0.007,0.2882,0.0716,0.003,0.381,0.1938,-0.008,
    ↪ 0.3919,0.3568,-0.047,0.4649,0.0579,-0.001,0.0767,0.0429,0.505,
    ↪ 0.0612,0.179,0.514,0.093,0.3224,0.521,0.0585,0.4581,0.483,
    ↪ 0.2048,0.0683,0.486,0.1788,0.1999,0.503,0.2064,0.3388,0.503,
    ↪ 0.1892,0.4792,0.538,0.2969,0.1672,0.49,0.3297,0.2778,0.481,
```

```
      ↪ 0.3266,0.4217,0.465,0.3764,0.0925,0.494,0.4697,0.1567,0.477,
      ↪ 0.4396,0.2746,0.483,0.4595,0.4132,0.481'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP128'

_symmetry_space_group_name_H-M "P b a 2"
_symmetry_Int_Tables_number 32

_cell_length_a 21.61000
_cell_length_b 19.63000
_cell_length_c 3.95100
_cell_angle_alpha 90.00000
_cell_angle_beta 90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,z
3 -x+1/2,y+1/2,z
4 x+1/2,-y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Mo1 Mo 2 a 0.00000 0.00000 0.57900 1.00000
O1 O 2 a 0.00000 0.00000 0.03300 1.00000
Mo2 Mo 4 c 0.02398 0.26126 0.58600 1.00000
Mo3 Mo 4 c 0.12899 0.11845 0.41770 1.00000
Mo4 Mo 4 c 0.13648 0.39897 0.43290 1.00000
Mo5 Mo 4 c 0.24260 0.25684 0.57900 1.00000
Mo6 Mo 4 c 0.28825 0.06556 0.58480 1.00000
Mo7 Mo 4 c 0.38285 0.19333 0.43340 1.00000
Mo8 Mo 4 c 0.38519 0.36637 0.42050 1.00000
Mo9 Mo 4 c 0.46561 0.05514 0.55510 1.00000
O2 O 4 c 0.02250 0.26300 0.02700 1.00000
O3 O 4 c 0.13100 0.12060 -0.05200 1.00000
O4 O 4 c 0.13490 0.40380 -0.01100 1.00000
O5 O 4 c 0.24650 0.25320 0.00700 1.00000
O6 O 4 c 0.28820 0.07160 0.00300 1.00000
O7 O 4 c 0.38100 0.19380 -0.00800 1.00000
O8 O 4 c 0.39190 0.35680 -0.04700 1.00000
O9 O 4 c 0.46490 0.05790 -0.00100 1.00000
O10 O 4 c 0.07670 0.04290 0.50500 1.00000
O11 O 4 c 0.06120 0.17900 0.51400 1.00000
O12 O 4 c 0.09300 0.32240 0.52100 1.00000
O13 O 4 c 0.05850 0.45810 0.48300 1.00000
O14 O 4 c 0.20480 0.06830 0.48600 1.00000
O15 O 4 c 0.17880 0.19990 0.50300 1.00000
O16 O 4 c 0.20640 0.33880 0.50300 1.00000
O17 O 4 c 0.18920 0.47920 0.53800 1.00000
O18 O 4 c 0.29690 0.16720 0.49000 1.00000
O19 O 4 c 0.32970 0.27780 0.48100 1.00000
O20 O 4 c 0.32660 0.42170 0.46500 1.00000
O21 O 4 c 0.37640 0.09250 0.49400 1.00000
O22 O 4 c 0.46970 0.15670 0.47700 1.00000
O23 O 4 c 0.43960 0.27460 0.48300 1.00000
O24 O 4 c 0.45950 0.41320 0.48100 1.00000
```

$Mo_{17}O_{47}$: A17B47_oP128_32_a8c_a23c - POSCAR

```
A17B47_oP128_32_a8c_a23c & a,b/a,c/a,z1,z2,x3,y3,z3,x4,y4,z4,x5,y5,z5,x6
      ↪ ,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10,x11,y11,z11,x12,
      ↪ y12,z12,x13,y13,z13,x14,y14,z14,x15,y15,z15,x16,y16,z16,x17,y17
      ↪ ,z17,x18,y18,z18,x19,y19,z19,x20,y20,z20,x21,y21,z21,x22,y22,
      ↪ z22,x23,y23,z23,x24,y24,z24,x25,y25,z25,x26,y26,z26,x27,y27,z27
      ↪ ,x28,y28,z28,x29,y29,z29,x30,y30,z30,x31,y31,z31,x32,y32,z32,
      ↪ x33,y33,z33 --params=21.61,0.908375751967,0.182832022212,0.579,
      ↪ 0.033,0.02398,0.26126,0.586,0.12899,0.11845,0.4177,0.13648,
      ↪ 0.39897,0.4329,0.2426,0.25684,0.579,0.28825,0.06556,0.5848,
      ↪ 0.38285,0.19333,0.4334,0.38519,0.36637,0.4205,0.46561,0.05514,
      ↪ 0.5551,0.0225,0.263,0.027,0.131,0.1206,-0.052,0.1349,0.4038,-
      ↪ 0.011,0.2465,0.2532,0.007,0.2882,0.0716,0.003,0.381,0.1938,-
      ↪ 0.008,0.3919,0.3568,-0.047,0.4649,0.0579,-0.001,0.0767,0.0429,
      ↪ 0.505,0.0612,0.179,0.514,0.093,0.3224,0.521,0.0585,0.4581,0.483
      ↪ ,0.2048,0.0683,0.486,0.1788,0.1999,0.503,0.2064,0.3388,0.503,
      ↪ 0.1892,0.4792,0.538,0.2969,0.1672,0.49,0.3297,0.2778,0.481,
      ↪ 0.3266,0.4217,0.465,0.3764,0.0925,0.494,0.4697,0.1567,0.477,
      ↪ 0.4396,0.2746,0.483,0.4595,0.4132,0.481 & Pba2 C_{2v}^{8} #32 (
      ↪ a^2c^31) & oP128 & None & Mo17O47 & Mo17O47 & L. Kihlborg, Acta
      ↪ Chem. Scand. 17, 1485-1487 (1963)
   1.00000000000000
  21.61000000000000    0.00000000000000    0.00000000000000
   0.00000000000000   19.63000000000000    0.00000000000000
   0.00000000000000    0.00000000000000    3.95100000000000
     Mo      O
     34      94
Direct
   0.00000000000000    0.00000000000000    0.57900000000000   Mo   (2a)
   0.50000000000000    0.50000000000000    0.57900000000000   Mo   (2a)
   0.02398000000000    0.26126000000000    0.58600000000000   Mo   (4c)
  -0.02398000000000   -0.26126000000000    0.58600000000000   Mo   (4c)
   0.52398000000000    0.23874000000000    0.58600000000000   Mo   (4c)
   0.47602000000000    0.76126000000000    0.58600000000000   Mo   (4c)
   0.12899000000000    0.11845000000000    0.41770000000000   Mo   (4c)
  -0.12899000000000   -0.11845000000000    0.41770000000000   Mo   (4c)
   0.62899000000000    0.38155000000000    0.41770000000000   Mo   (4c)
   0.37101000000000    0.61845000000000    0.41770000000000   Mo   (4c)
   0.13648000000000    0.39897000000000    0.43290000000000   Mo   (4c)
  -0.13648000000000   -0.39897000000000    0.43290000000000   Mo   (4c)
   0.63648000000000    0.10103000000000    0.43290000000000   Mo   (4c)
   0.36352000000000    0.89897000000000    0.43290000000000   Mo   (4c)
   0.24260000000000    0.25684000000000    0.57900000000000   Mo   (4c)
  -0.24260000000000   -0.25684000000000    0.57900000000000   Mo   (4c)
   0.74260000000000    0.24316000000000    0.57900000000000   Mo   (4c)
   0.25740000000000    0.75684000000000    0.57900000000000   Mo   (4c)
   0.28825000000000    0.06556000000000    0.58480000000000   Mo   (4c)
  -0.28825000000000   -0.06556000000000    0.58480000000000   Mo   (4c)
   0.78825000000000    0.43444000000000    0.58480000000000   Mo   (4c)
   0.21175000000000    0.56556000000000    0.58480000000000   Mo   (4c)
   0.38285000000000    0.19333000000000    0.43340000000000   Mo   (4c)
  -0.38285000000000   -0.19333000000000    0.43340000000000   Mo   (4c)
   0.88285000000000    0.30667000000000    0.43340000000000   Mo   (4c)
   0.11715000000000    0.69333000000000    0.43340000000000   Mo   (4c)
   0.38519000000000    0.36637000000000    0.42050000000000   Mo   (4c)
  -0.38519000000000   -0.36637000000000    0.42050000000000   Mo   (4c)
   0.88519000000000    0.13363000000000    0.42050000000000   Mo   (4c)
   0.11481000000000    0.86637000000000    0.42050000000000   Mo   (4c)
   0.46561000000000    0.05514000000000    0.55510000000000   Mo   (4c)
  -0.46561000000000   -0.05514000000000    0.55510000000000   Mo   (4c)
   0.96561000000000    0.44486000000000    0.55510000000000   Mo   (4c)
   0.03439000000000    0.55514000000000    0.55510000000000   Mo   (4c)
   0.00000000000000    0.00000000000000    0.03300000000000   O    (2a)
   0.50000000000000    0.50000000000000    0.03300000000000   O    (2a)
   0.02250000000000    0.26300000000000    0.02700000000000   O    (4c)
  -0.02250000000000   -0.26300000000000    0.02700000000000   O    (4c)
   0.52250000000000    0.23700000000000    0.02700000000000   O    (4c)
   0.47750000000000    0.76300000000000    0.02700000000000   O    (4c)
   0.13100000000000    0.12060000000000   -0.05200000000000   O    (4c)
  -0.13100000000000   -0.12060000000000   -0.05200000000000   O    (4c)
   0.63100000000000    0.37940000000000   -0.05200000000000   O    (4c)
   0.36900000000000    0.62060000000000   -0.05200000000000   O    (4c)
   0.13490000000000    0.40380000000000   -0.01100000000000   O    (4c)
  -0.13490000000000   -0.40380000000000   -0.01100000000000   O    (4c)
   0.63490000000000    0.09620000000000   -0.01100000000000   O    (4c)
   0.36510000000000    0.90380000000000   -0.01100000000000   O    (4c)
   0.24650000000000    0.25320000000000    0.00700000000000   O    (4c)
  -0.24650000000000   -0.25320000000000    0.00700000000000   O    (4c)
   0.74650000000000    0.24680000000000    0.00700000000000   O    (4c)
   0.25350000000000    0.75320000000000    0.00700000000000   O    (4c)
   0.28820000000000    0.07160000000000    0.00300000000000   O    (4c)
  -0.28820000000000   -0.07160000000000    0.00300000000000   O    (4c)
   0.78820000000000    0.42840000000000    0.00300000000000   O    (4c)
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  -0.38100000000000   -0.19380000000000   -0.00800000000000   O    (4c)
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   0.11900000000000    0.69380000000000   -0.00800000000000   O    (4c)
   0.39190000000000    0.35680000000000   -0.04700000000000   O    (4c)
  -0.39190000000000   -0.35680000000000   -0.04700000000000   O    (4c)
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   0.10810000000000    0.85680000000000   -0.04700000000000   O    (4c)
   0.46490000000000    0.05790000000000   -0.00100000000000   O    (4c)
  -0.46490000000000   -0.05790000000000   -0.00100000000000   O    (4c)
   0.96490000000000    0.44210000000000   -0.00100000000000   O    (4c)
   0.03510000000000    0.55790000000000   -0.00100000000000   O    (4c)
   0.07670000000000    0.04290000000000    0.50500000000000   O    (4c)
  -0.07670000000000   -0.04290000000000    0.50500000000000   O    (4c)
   0.57670000000000    0.45710000000000    0.50500000000000   O    (4c)
   0.42330000000000    0.54290000000000    0.50500000000000   O    (4c)
   0.06120000000000    0.17900000000000    0.51400000000000   O    (4c)
  -0.06120000000000   -0.17900000000000    0.51400000000000   O    (4c)
   0.56120000000000    0.32100000000000    0.51400000000000   O    (4c)
   0.43880000000000    0.67900000000000    0.51400000000000   O    (4c)
   0.09300000000000    0.32240000000000    0.52100000000000   O    (4c)
  -0.09300000000000   -0.32240000000000    0.52100000000000   O    (4c)
   0.59300000000000    0.17760000000000    0.52100000000000   O    (4c)
   0.40700000000000    0.82240000000000    0.52100000000000   O    (4c)
   0.05850000000000    0.45810000000000    0.48300000000000   O    (4c)
  -0.05850000000000   -0.45810000000000    0.48300000000000   O    (4c)
   0.55850000000000    0.04190000000000    0.48300000000000   O    (4c)
   0.44150000000000    0.95810000000000    0.48300000000000   O    (4c)
   0.20480000000000    0.06830000000000    0.48600000000000   O    (4c)
  -0.20480000000000   -0.06830000000000    0.48600000000000   O    (4c)
   0.70480000000000    0.43170000000000    0.48600000000000   O    (4c)
   0.29520000000000    0.56830000000000    0.48600000000000   O    (4c)
   0.17880000000000    0.19990000000000    0.50300000000000   O    (4c)
  -0.17880000000000   -0.19990000000000    0.50300000000000   O    (4c)
   0.67880000000000    0.30010000000000    0.50300000000000   O    (4c)
   0.32120000000000    0.69990000000000    0.50300000000000   O    (4c)
   0.20640000000000    0.33880000000000    0.50300000000000   O    (4c)
  -0.20640000000000   -0.33880000000000    0.50300000000000   O    (4c)
   0.70640000000000    0.16120000000000    0.50300000000000   O    (4c)
   0.29360000000000    0.83880000000000    0.50300000000000   O    (4c)
   0.18920000000000    0.47920000000000    0.53800000000000   O    (4c)
  -0.18920000000000   -0.47920000000000    0.53800000000000   O    (4c)
   0.68920000000000    0.02080000000000    0.53800000000000   O    (4c)
   0.31080000000000    0.97920000000000    0.53800000000000   O    (4c)
   0.29690000000000    0.16720000000000    0.49000000000000   O    (4c)
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   0.79690000000000    0.33280000000000    0.49000000000000   O    (4c)
   0.20310000000000    0.66720000000000    0.49000000000000   O    (4c)
   0.32970000000000    0.27780000000000    0.48100000000000   O    (4c)
  -0.32970000000000   -0.27780000000000    0.48100000000000   O    (4c)
   0.82970000000000    0.22220000000000    0.48100000000000   O    (4c)
   0.17030000000000    0.77780000000000    0.48100000000000   O    (4c)
   0.32660000000000    0.42170000000000    0.46500000000000   O    (4c)
  -0.32660000000000   -0.42170000000000    0.46500000000000   O    (4c)
   0.82660000000000    0.07830000000000    0.46500000000000   O    (4c)
   0.17340000000000    0.92170000000000    0.46500000000000   O    (4c)
   0.37640000000000    0.09250000000000    0.49400000000000   O    (4c)
  -0.37640000000000   -0.09250000000000    0.49400000000000   O    (4c)
   0.87640000000000    0.40750000000000    0.49400000000000   O    (4c)
```

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   0.12360000000000   0.59250000000000   0.49400000000000   O   (4c)
   0.46970000000000   0.15670000000000   0.47700000000000   O   (4c)
  -0.46970000000000  -0.15670000000000   0.47700000000000   O   (4c)
   0.96970000000000   0.34330000000000   0.47700000000000   O   (4c)
   0.03030000000000   0.65670000000000   0.47700000000000   O   (4c)
   0.43960000000000   0.27460000000000   0.48300000000000   O   (4c)
  -0.43960000000000  -0.27460000000000   0.48300000000000   O   (4c)
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   0.06040000000000   0.77460000000000   0.48300000000000   O   (4c)
   0.45950000000000   0.41320000000000   0.48100000000000   O   (4c)
  -0.45950000000000  -0.41320000000000   0.48100000000000   O   (4c)
   0.95950000000000   0.08680000000000   0.48100000000000   O   (4c)
   0.04050000000000   0.91320000000000   0.48100000000000   O   (4c)
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Possible δ-$Gd_2Si_2O_7$: A2B7C2_oP44_33_2a_7a_2a - CIF

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_chemical_formula_sum 'Gd2 O7 Si2'

loop_
_publ_author_name
 'Y. I. Smolin'
 'Y. F. Shepelev'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 26
_journal_year 1970
_journal_page_first 484
_journal_page_last 492
_publ_Section_title
;
 The Crystal Structures of the Rare Earth Pyrosilicates
;

# Found in The crystal structure of $\delta$-yttrium pyrosilicate, $\
    ↪ delta$-Y$_{2}$Si$_{2}$O$_{7}$, 1990

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    ↪ x_{7},y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{10},y_
    ↪ {10},z_{10},x_{11},y_{11},z_{11}'
_aflow_params_values '13.87,0.365753424658,0.600576784427,0.12551,0.3373
    ↪ ,-0.00169,0.12564,0.33739,0.51409,0.2715,0.4769,0.0876,0.2658,
    ↪ 0.4857,0.413,0.3457,0.0706,0.2465,0.4211,0.5557,0.2448,0.5472,
    ↪ 0.7858,0.0866,0.5456,0.7882,0.4206,0.5988,0.3526,0.256,0.3205,
    ↪ 0.3744,0.2505,0.539,0.6253,0.2498'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP44'

_symmetry_space_group_name_H-M "P n a 21"
_symmetry_Int_Tables_number 33

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_cell_length_b    5.07300
_cell_length_c    8.33000
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

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_space_group_symop_operation_xyz
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2 -x,-y,z+1/2
3 -x+1/2,y+1/2,z+1/2
4 x+1/2,-y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
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Gd2 Gd 4 a 0.12564 0.33739 0.51409 1.00000
O1 O 4 a 0.27150 0.47690 0.08760 1.00000
O2 O 4 a 0.26580 0.48570 0.41300 1.00000
O3 O 4 a 0.34570 0.07060 0.24650 1.00000
O4 O 4 a 0.42110 0.55570 0.24480 1.00000
O5 O 4 a 0.54720 0.78580 0.08660 1.00000
O6 O 4 a 0.54560 0.78820 0.42060 1.00000
O7 O 4 a 0.59880 0.35260 0.25600 1.00000
Si1 Si 4 a 0.32050 0.37440 0.25050 1.00000
Si2 Si 4 a 0.53900 0.62530 0.24980 1.00000
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Possible δ-$Gd_2Si_2O_7$: A2B7C2_oP44_33_2a_7a_2a - POSCAR

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A2B7C2_oP44_33_2a_7a_2a & a,b/a,c/a,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,z4,
    ↪ x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10,x11,
    ↪ y11,z11 --params=13.87,0.365753424658,0.600576784427,0.12551,
    ↪ 0.3373,-0.00169,0.12564,0.33739,0.51409,0.2715,0.4769,0.0876,
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    ↪ 0.5472,0.7858,0.0866,0.5456,0.7882,0.4206,0.5988,0.3526,0.256,
    ↪ 0.3205,0.3744,0.2505,0.539,0.6253,0.2498 & Pna2_{1} C_{2v}^{9}
    ↪ #33 (a^11) & oP44 & None & Gd2O7Si2 & Gd2O7Si2 & Y. I. Smolin
    ↪ and Y. F. Shepelev, Acta Crystallogr. Sect. B Struct. Sci. 26,
    ↪ 484-492 (1970)
  1.00000000000000
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  0.00000000000000   5.07300000000000   0.00000000000000
  0.00000000000000   0.00000000000000   8.33000000000000
   Gd    O    Si
    8    28    8
Direct
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  -0.12551000000000  -0.33730000000000   0.49831000000000   Gd   (4a)
   0.62551000000000   0.16270000000000  -0.00169000000000   Gd   (4a)
   0.37449000000000   0.83730000000000   0.49831000000000   Gd   (4a)
   0.12564000000000   0.33739000000000   0.51409000000000   Gd   (4a)
  -0.12564000000000  -0.33739000000000   1.01409000000000   Gd   (4a)
   0.62564000000000   0.16261000000000   0.51409000000000   Gd   (4a)
   0.37436000000000   0.83739000000000   1.01409000000000   Gd   (4a)
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  -0.27150000000000  -0.47690000000000   0.58760000000000   O    (4a)
   0.77150000000000   0.02310000000000   0.08760000000000   O    (4a)
   0.22850000000000   0.97690000000000   0.58760000000000   O    (4a)
   0.26580000000000   0.48570000000000   0.41300000000000   O    (4a)
  -0.26580000000000  -0.48570000000000   0.91300000000000   O    (4a)
   0.76580000000000   0.01430000000000   0.41300000000000   O    (4a)
   0.23420000000000   0.98570000000000   0.91300000000000   O    (4a)
   0.34570000000000   0.07060000000000   0.24650000000000   O    (4a)
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   0.15430000000000   0.57060000000000   0.74650000000000   O    (4a)
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   0.53900000000000   0.62530000000000   0.24980000000000   Si   (4a)
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CaB_2O_4 (III): A2BC4_oP84_33_6a_3a_12a - CIF

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_audit_creation_method FINDSYM

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loop_
_publ_author_name
 'M. Marezio'
 'J. P. Remeika'
 'P. D. Dernier'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 25
_journal_year 1969
_journal_page_first 955
_journal_page_last 964
_publ_Section_title
;
 The crystal structure of the high-pressure phase CaB$_{2}$O$_{4}$(III)
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    ↪ x_{7},y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{10},y_
    ↪ {10},z_{10},x_{11},y_{11},z_{11},x_{12},y_{12},z_{12},x_{13},y_
    ↪ {13},z_{13},x_{14},y_{14},z_{14},x_{15},y_{15},z_{15},x_{16},y_
    ↪ {16},z_{16},x_{17},y_{17},z_{17},x_{18},y_{18},z_{18},x_{19},y_
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    ↪ 0.9442,0.4658,0.0183,0.48,0.4748,0.3758,0.2477,0.07351,0.05043,
    ↪ 0.0,0.03257,0.65995,0.23386,0.25309,0.11288,0.27955,0.2657,
    ↪ 0.1944,0.0459,0.2171,0.4704,0.1734,0.0717,0.0297,0.2009,0.4547,
    ↪ 0.1715,0.2567,0.2009,0.375,0.4486,0.1456,0.0322,0.5116,0.0061,
    ↪ 0.3009,0.5305,0.1734,0.2946,0.6635,0.1069,0.022,0.7981,0.287,
    ↪ 0.2046,0.8336,0.4563,0.137,0.9458,0.1504,0.4768,0.9668'
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_aflow_Pearson 'oP84'

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_symmetry_Int_Tables_number 33
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_cell_angle_gamma  90.00000

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_space_group_symop_operation_xyz
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3 -x+1/2,y+1/2,z+1/2
4 x+1/2,-y+1/2,z

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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
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B2 B 4 a 0.25100 0.41930 0.05150 1.00000
B3 B 4 a 0.21310 0.11910 0.74030 1.00000
B4 B 4 a 0.32220 0.10060 0.94420 1.00000
B5 B 4 a 0.46580 0.01830 0.48000 1.00000
B6 B 4 a 0.47480 0.37580 0.24770 1.00000
Ca1 Ca 4 a 0.07351 0.05043 0.00000 1.00000
Ca2 Ca 4 a 0.03257 0.65995 0.23386 1.00000
Ca3 Ca 4 a 0.25309 0.11288 0.27955 1.00000
O1 O 4 a 0.26570 0.19440 0.04590 1.00000
O2 O 4 a 0.21710 0.47040 0.17340 1.00000
O3 O 4 a 0.07170 0.02970 0.20090 1.00000
O4 O 4 a 0.45470 0.17150 0.25670 1.00000
O5 O 4 a 0.20090 0.37500 0.44860 1.00000
O6 O 4 a 0.14560 0.03220 0.51160 1.00000
O7 O 4 a 0.00610 0.30090 0.53050 1.00000
O8 O 4 a 0.17340 0.29460 0.66350 1.00000
O9 O 4 a 0.10690 0.02200 0.79810 1.00000
O10 O 4 a 0.28700 0.20460 0.83360 1.00000
O11 O 4 a 0.45630 0.13700 0.94580 1.00000
O12 O 4 a 0.15040 0.47680 0.96680 1.00000
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CaB_2O_4 (III): A2BC4_oP84_33_6a_3a_12a - POSCAR

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    ↪ x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10,x11,
    ↪ y11,z11,x12,y12,z12,x13,y13,z13,x14,y14,z14,x15,y15,z15,x16,y16
    ↪ ,z16,x17,y17,z17,x18,y18,z18,x19,y19,z19,x20,y20,z20,x21,y21,
    ↪ z21 --params=11.38,0.554217926186,0.993321616872,0.1401,0.2529,
    ↪ 0.539,0.251,0.4193,0.0515,0.2131,0.1191,0.7403,0.3222,0.1006,
    ↪ 0.9442,0.4658,0.0183,0.48,0.4748,0.3758,0.2477,0.07351,0.05043,
    ↪ 0.0,0.03257,0.65995,0.23386,0.25309,0.11288,0.27955,0.2657,
    ↪ 0.1944,0.0459,0.2171,0.4704,0.1734,0.0717,0.0297,0.2009,0.4547,
    ↪ 0.1715,0.2567,0.2009,0.375,0.4486,0.1456,0.0322,0.5116,0.0061,
    ↪ 0.3009,0.5305,0.1734,0.2946,0.6635,0.1069,0.022,0.7981,0.287,
    ↪ 0.2046,0.8336,0.4563,0.137,0.9458,0.1504,0.4768,0.9668 & Pna2_{
    ↪ 1} C_{2v}^{9} #33 (a^21) & oP84 & None & B2CaO4 & B2CaO4 & M.
    ↪ Marezio and J. P. Remeika and P. D. Dernier, Acta Crystallogr.
    ↪ Sect. B Struct. Sci. 25, 955-964 (1969)
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    24     12     48
Direct
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Cervantite (α-Sb_2O_4): A2B_oP24_33_4a_2a - CIF

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_audit_creation_method FINDSYM

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_chemical_formula_sum 'O2 Sb'

loop_
_publ_author_name
 'G. Thornton'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 33
_journal_year 1977
_journal_page_first 1271
_journal_page_last 1273
_publ_Section_title
;
 A Neutron Diffraction Study of $\alpha$-Sb$_{2}$O$_{3}$
;

# Found in The American Mineralogist Crystal Structure Database, 2003

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_aflow_proto 'A2B_oP24_33_4a_2a'
_aflow_params 'a,b/a,c/a,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3},y_{3
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_aflow_params_values '5.456,0.882331378299,2.16037390029,0.34,0.177,
    ↪ 0.0962,0.159,0.71,0.195,0.086,0.208,0.312,0.33,0.84,0.41,-0.032
    ↪ ,0.036,0.009,0.366,0.016,0.253'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP24'

_symmetry_space_group_name_H-M "P n a 21"
_symmetry_Int_Tables_number 33

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_cell_length_c      11.78700
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_cell_angle_gamma  90.00000

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3 -x+1/2,y+1/2,z+1/2
4 x+1/2,-y+1/2,z
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_atom_site_Wyckoff_label
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_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
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O2 O 4 a 0.15900 0.71000 0.19500 1.00000
O3 O 4 a 0.08600 0.20800 0.31200 1.00000
O4 O 4 a 0.33000 0.84000 0.41000 1.00000
Sb1 Sb 4 a −0.03200 0.03600 0.00900 1.00000
Sb2 Sb 4 a 0.36600 0.01600 0.25300 1.00000
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Cervantite (α-Sb_2O_4): A2B_oP24_33_4a_2a - POSCAR

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    ↪ 0.177,0.0962,0.159,0.71,0.195,0.086,0.208,0.312,0.33,0.84,0.41
    ↪ ,−0.032,0.036,0.009,0.366,0.016,0.253 & Pna2_{1} C_{2v}^{9} #33
    ↪  (a^6) & oP24 & None & O2Sb & Cervantite & G. Thornton, Acta
    ↪ Crystallogr. Sect. B Struct. Sci. 33, 1271−1273 (1977)
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  5.45600000000000   0.00000000000000   0.00000000000000
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    O     Sb
   16      8
Direct
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CsB_4O_6F: A4BCD6_oP48_33_4a_a_a_6a - CIF

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 'Y. Wang'
 'B. Zhang'
 'F. Zhang'
 'Z. Yang'
 'S. Pan'
_journal_name_full_name
;
 Angewandte Chemie
;
_journal_volume 129
_journal_year 2017
_journal_page_first 14307
_journal_page_last 14311
_publ_Section_title
;
 CsB${4}$O${6}$F: A Congruent-Melting Deep-Ultraviolet Nonlinear
    ↪ Optical Material by Combining Superior Functional Units
;

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_aflow_proto 'A4BCD6_oP48_33_4a_a_a_6a'
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    ↪ x_{7},y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{10},y_
    ↪ {10},z_{10},x_{11},y_{11},z_{11},x_{12},y_{12},z_{12}'
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    ↪ 0.348,0.49584,0.22117,0.852,0.0566,0.4371,0.3269,0.306,0.4658,
    ↪ 0.8049,0.2888,0.6655,0.701,0.3,0.5194,0.4588,0.2626,0.3109,
    ↪ 0.4526,0.3038,0.4165,0.1445,0.2784,0.6167,0.0497'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP48'

_symmetry_space_group_name_H-M "P n a 21"
_symmetry_Int_Tables_number 33

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3 −x+1/2,y+1/2,z+1/2
4 x+1/2,−y+1/2,z

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_atom_site_fract_x
_atom_site_fract_y
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B3 B 4 a 0.23040 0.20030 0.39420 1.00000
B4 B 4 a 0.23270 0.41890 0.34800 1.00000
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F1 F 4 a 0.05660 0.43710 0.32690 1.00000
O1 O 4 a 0.30600 0.46580 0.80490 1.00000
O2 O 4 a 0.28880 0.66550 0.70100 1.00000
O3 O 4 a 0.30000 0.51940 0.45880 1.00000
O4 O 4 a 0.26260 0.31090 0.45260 1.00000
O5 O 4 a 0.30380 0.41650 0.14450 1.00000
O6 O 4 a 0.27840 0.61670 0.04970 1.00000
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CsB_4O_6F: A4BCD6_oP48_33_4a_a_a_6a - POSCAR

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    ↪ 0.840953546775,0.2971,0.5446,0.6555,0.2977,0.5006,0.0044,0.2304
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    ↪ B4CsFO6 & B4CsFO6 & X. Wang et al., Angew. Chem. 129,
    ↪ 14307−14311 (2017)
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 −0.30000000000000  −0.51940000000000   0.95880000000000   O   (4a)
  0.80000000000000  −0.01940000000000   0.45880000000000   O   (4a)
  0.20000000000000   1.01940000000000   0.95880000000000   O   (4a)
  0.26260000000000   0.31090000000000   0.45260000000000   O   (4a)
 −0.26260000000000  −0.31090000000000   0.95260000000000   O   (4a)
  0.76260000000000   0.18910000000000   0.45260000000000   O   (4a)
  0.23740000000000   0.81090000000000   0.95260000000000   O   (4a)
  0.30380000000000   0.41650000000000   0.14450000000000   O   (4a)
 −0.30380000000000  −0.41650000000000   0.64450000000000   O   (4a)
  0.80380000000000   0.08350000000000   0.14450000000000   O   (4a)
  0.19620000000000   0.91650000000000   0.64450000000000   O   (4a)
  0.27840000000000   0.61670000000000   0.04970000000000   O   (4a)
 −0.27840000000000  −0.61670000000000   0.54970000000000   O   (4a)
  0.77840000000000  −0.11670000000000   0.04970000000000   O   (4a)
  0.22160000000000   1.11670000000000   0.54970000000000   O   (4a)
```

$LiGaO_2$: ABC2_oP16_33_a_a_2a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'GaLiO2'
_chemical_formula_sum 'Ga Li O2'

loop_
_publ_author_name
 'M. Marezio'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 18
_journal_year 1965
_journal_page_first 481
_journal_page_last 484
_publ_Section_title
;
 The Crystal Structure of LiGaO$_{2}$
;

_aflow_title 'LiGaO$_{2}$ Structure'
_aflow_proto 'ABC2_oP16_33_a_a_2a'
_aflow_params 'a,b/a,c/a,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3},y_{3
    ↪ },z_{3},x_{4},y_{4},z_{4}'
_aflow_params_values '5.402,1.17956312477,0.926878933728,0.0821,0.1263,
    ↪ 0.0,0.4207,0.1267,0.4936,0.4066,0.1388,0.8972,0.0697,0.1121,
    ↪ 0.3708'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP16'

_symmetry_space_group_name_H-M "P n a 21"
_symmetry_Int_Tables_number 33

_cell_length_a    5.40200
_cell_length_b    6.37200
_cell_length_c    5.00700
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,z+1/2
3 -x+1/2,y+1/2,z+1/2
4 x+1/2,-y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ga1 Ga 4 a 0.08210 0.12630 0.00000 1.00000
Li1 Li 4 a 0.42070 0.12670 0.49360 1.00000
O1 O 4 a 0.40660 0.13880 0.89720 1.00000
O2 O 4 a 0.06970 0.11210 0.37080 1.00000
```

$LiGaO_2$: ABC2_oP16_33_a_a_2a - POSCAR

```
ABC2_oP16_33_a_a_2a & a,b/a,c/a,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,z4 --
    ↪ params=5.402,1.17956312477,0.926878933728,0.0821,0.1263,0.0,
    ↪ 0.4207,0.1267,0.4936,0.4066,0.1388,0.8972,0.0697,0.1121,0.3708
    ↪ & Pna2_{1} C_{2v}^{9} #33 (a^4) & oP16 & None & GaLiO2 & GaLiO2
    ↪ & M. Marezio, Acta Cryst. 18, 481-484 (1965)
   1.00000000000000
   5.40200000000000   0.00000000000000   0.00000000000000
   0.00000000000000   6.37200000000000   0.00000000000000
   0.00000000000000   0.00000000000000   5.00700000000000
    Ga   Li   O
     4    4   8
Direct
   0.08210000000000   0.12630000000000   0.00000000000000  Ga  (4a)
  -0.08210000000000  -0.12630000000000   0.50000000000000  Ga  (4a)
   0.58210000000000   0.37370000000000   0.00000000000000  Ga  (4a)
   0.41790000000000   0.62630000000000   0.50000000000000  Ga  (4a)
   0.42070000000000   0.12670000000000   0.49360000000000  Li  (4a)
  -0.42070000000000  -0.12670000000000   0.99360000000000  Li  (4a)
   0.92070000000000   0.37330000000000   0.49360000000000  Li  (4a)
   0.07930000000000   0.62670000000000   0.99360000000000  Li  (4a)
   0.40660000000000   0.13880000000000   0.89720000000000  O  (4a)
  -0.40660000000000  -0.13880000000000   1.39720000000000  O  (4a)
   0.90660000000000   0.36120000000000   0.89720000000000  O  (4a)
   0.09340000000000   0.63880000000000   1.39720000000000  O  (4a)
   0.06970000000000   0.11210000000000   0.37080000000000  O  (4a)
  -0.06970000000000  -0.11210000000000   0.87080000000000  O  (4a)
   0.56970000000000   0.38790000000000   0.37080000000000  O  (4a)
   0.43030000000000   0.61210000000000   0.87080000000000  O  (4a)
```

γ-$LiIO_3$: ABC3_oP20_33_a_a_3a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'ILiO3'
_chemical_formula_sum 'I Li O3'

loop_
_publ_author_name
 'R. Liminga'
 'C. Svensson'
 'J. Albertsson'
 'S. C. Abrahams'
_journal_name_full_name
;
 Journal of Chemical Physics
;
_journal_volume 77
_journal_year 1982
_journal_page_first 4222
_journal_page_last 4226
_publ_Section_title
;
 Gamma-lithium iodate structure at 515~K and the $\alpha$-LiIO$_{3}$ to
    ↪ $\gamma$-LiIO$_{3}$, $\gamma$-LiIO$_{3}$ to $\beta$-LiIO$_{3}$
    ↪  phase transitions
;

_aflow_title '$\gamma$-LiIO$_{3}$ Structure'
_aflow_proto 'ABC3_oP20_33_a_a_3a'
_aflow_params 'a,b/a,c/a,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3},y_{3
    ↪ },z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{5}'
_aflow_params_values '9.422,0.622054765443,0.562619401401,0.3179,0.0751,
    ↪ 0.0,0.0,0.0,0.0706,0.1298,0.0556,0.8087,0.4117,0.8806,0.8382,
    ↪ 0.3663,0.3439,0.8147'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP20'

_symmetry_space_group_name_H-M "P n a 21"
_symmetry_Int_Tables_number 33

_cell_length_a    9.42200
_cell_length_b    5.86100
_cell_length_c    5.30100
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,z+1/2
3 -x+1/2,y+1/2,z+1/2
4 x+1/2,-y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
I1 I 4 a 0.31790 0.07510 0.00000 1.00000
Li1 Li 4 a 0.00000 0.00000 0.07060 1.00000
O1 O 4 a 0.12980 0.05560 0.80870 1.00000
O2 O 4 a 0.41170 0.88060 0.83820 1.00000
O3 O 4 a 0.36630 0.34390 0.81470 1.00000
```

γ-$LiIO_3$: ABC3_oP20_33_a_a_3a - POSCAR

```
ABC3_oP20_33_a_a_3a & a,b/a,c/a,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,z4,x5,
    ↪ y5,z5 --params=9.422,0.622054765443,0.562619401401,0.3179,
    ↪ 0.0751,0.0,0.0,0.0,0.0706,0.1298,0.0556,0.8087,0.4117,0.8806,
    ↪ 0.8382,0.3663,0.3439,0.8147 & Pna2_{1} C_{2v}^{9} #33 (a^5) &
    ↪ oP20 & None & ILiO3 & ILiO3 & R. Liminga et al., J. Chem. Phys.
    ↪  77, 4222-4226 (1982)
   1.00000000000000
   9.42200000000000   0.00000000000000   0.00000000000000
   0.00000000000000   5.86100000000000   0.00000000000000
   0.00000000000000   0.00000000000000   5.30100000000000
    I   Li   O
    4    4  12
Direct
   0.31790000000000   0.07510000000000   0.00000000000000  I  (4a)
  -0.31790000000000  -0.07510000000000   0.50000000000000  I  (4a)
   0.81790000000000   0.42490000000000   0.00000000000000  I  (4a)
   0.18210000000000   0.57510000000000   0.50000000000000  I  (4a)
   0.00000000000000   0.00000000000000   0.07060000000000  Li  (4a)
   0.00000000000000   0.00000000000000   0.57060000000000  Li  (4a)
   0.50000000000000   0.50000000000000   0.07060000000000  Li  (4a)
   0.50000000000000   0.50000000000000   0.57060000000000  Li  (4a)
   0.12980000000000   0.05560000000000   0.80870000000000  O  (4a)
  -0.12980000000000  -0.05560000000000   1.30870000000000  O  (4a)
   0.62980000000000   0.44440000000000   0.80870000000000  O  (4a)
   0.37020000000000   0.55560000000000   1.30870000000000  O  (4a)
   0.41170000000000   0.88060000000000   0.83820000000000  O  (4a)
  -0.41170000000000  -0.88060000000000   1.33820000000000  O  (4a)
   0.91170000000000  -0.38060000000000   0.83820000000000  O  (4a)
   0.08830000000000   1.38060000000000   1.33820000000000  O  (4a)
   0.36630000000000   0.34390000000000   0.81470000000000  O  (4a)
  -0.36630000000000  -0.34390000000000   1.31470000000000  O  (4a)
   0.86630000000000   0.15610000000000   0.81470000000000  O  (4a)
   0.13370000000000   0.84390000000000   1.31470000000000  O  (4a)
```

$MnF_{2-x}(OH)_x$: A2BC2_oP10_34_c_a_c - CIF

```
# CIF file
data_findsym-output
```

```
_audit_creation_method FINDSYM

_chemical_name_mineral 'F_{2-x}HxMnOx'
_chemical_formula_sum 'H2 Mn O2'

loop_
_publ_author_name
 'H. B. Yahia'
 'M. Shikano'
 'H. Kobayashi'
 'M. Avdeev'
 'S. Liu'
 'C. D. Ling'
_journal_name_full_name
;
 Physical Chemistry Chemical Physics
;
_journal_volume 15
_journal_year 2013
_journal_page_first 13061
_journal_page_last 13069
_publ_Section_title
;
 Synthesis and characterization of the crystal structure and magnetic
      ↪ properties of the hydroxyfluoride MnF$_{2-x}$(OH)$_{x}$ ($x \
      ↪ approx 0.8$)
;

_aflow_title 'MnF$_{2-x}$(OH)$_{x}$ Structure'
_aflow_proto 'A2BC2_oP10_34_c_a_c'
_aflow_params 'a,b/a,c/a,z_{1},x_{2},y_{2},z_{2},x_{3},y_{3},z_{3}'
_aflow_params_values '4.71143,1.11290415012,0.689459463475,0.5,0.425,
      ↪ 0.0287,0.543,0.2511,0.1533,0.504'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP10'

_symmetry_space_group_name_H-M "P n n 2"
_symmetry_Int_Tables_number 34

_cell_length_a    4.71143
_cell_length_b    5.24337
_cell_length_c    3.24834
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,z
3 -x+1/2,y+1/2,z+1/2
4 x+1/2,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Mn1 Mn 2 a 0.00000 0.00000 0.50000 1.00000
H1 H 4 c 0.42500 0.02870 0.54300 1.00000
O1 O 4 c 0.25110 0.15330 0.50400 1.00000
```

$MnF_{2-x}(OH)_x$: A2BC2_oP10_34_c_a_c - POSCAR

```
A2BC2_oP10_34_c_a_c & a,b/a,c/a,z1,x2,y2,z2,x3,y3,z3 --params=4.71143,
      ↪ 1.11290415012,0.689459463475,0.5,0.425,0.0287,0.543,0.2511,
      ↪ 0.1533,0.504 & Pnn2 C_{2v}^{10} #34 (ac^2) & oP10 & None &
      ↪ xHxMnOx & F_{2 & H. B. Yahia et al., Phys. Chem. Chem. Phys. 15
      ↪ , 13061-13069 (2013)
   1.00000000000000
   4.71143000000000   0.00000000000000   0.00000000000000
   0.00000000000000   5.24337000000000   0.00000000000000
   0.00000000000000   0.00000000000000   3.24834000000000
     H     Mn     O
     4     2     4
Direct
   0.42500000000000   0.02870000000000   0.54300000000000   H   (4c)
  -0.42500000000000  -0.02870000000000   0.54300000000000   H   (4c)
   0.92500000000000   0.47130000000000   1.04300000000000   H   (4c)
   0.07500000000000   0.52870000000000   1.04300000000000   H   (4c)
   0.00000000000000   0.00000000000000   0.50000000000000   Mn  (2a)
   0.50000000000000   0.50000000000000   1.00000000000000   Mn  (2a)
   0.25110000000000   0.15330000000000   0.50400000000000   O   (4c)
  -0.25110000000000  -0.15330000000000   0.50400000000000   O   (4c)
   0.75110000000000   0.34670000000000   1.00400000000000   O   (4c)
   0.24890000000000   0.65330000000000   1.00400000000000   O   (4c)
```

Si_2N_2O: A2BC2_oC20_36_b_a_b - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'N2OSi2'
_chemical_formula_sum 'N2 O Si2'

loop_
_publ_author_name
 'I. Idrestedt'
 'C. Brosset'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 18
_journal_year 1964
_journal_page_first 1879
_journal_page_last 1886
_publ_Section_title
;
 Structure of Si$_{2}$N$_{2}$O
;

_aflow_title 'Si$_{2}$N$_{2}$O Structure'
_aflow_proto 'A2BC2_oC20_36_b_a_b'
_aflow_params 'a,b/a,c/a,y_{1},z_{1},x_{2},y_{2},z_{2},x_{3},y_{3},z_{3
      ↪ }'
_aflow_params_values '8.843,0.618907610539,0.546760149271,0.214,0.23,
      ↪ 0.218,0.121,0.642,0.1763,0.1509,0.2898'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oC20'

_symmetry_space_group_name_H-M "C m c 21"
_symmetry_Int_Tables_number 36

_cell_length_a    8.84300
_cell_length_b    5.47300
_cell_length_c    4.83500
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,z+1/2
3 -x,y,z
4 x,-y,z+1/2
5 x+1/2,y+1/2,z
6 -x+1/2,-y+1/2,z+1/2
7 -x+1/2,y+1/2,z
8 x+1/2,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 4 a 0.00000 0.21400 0.23000 1.00000
N1 N 8 b 0.21800 0.12100 0.64200 1.00000
Si1 Si 8 b 0.17630 0.15090 0.28980 1.00000
```

Si_2N_2O: A2BC2_oC20_36_b_a_b - POSCAR

```
A2BC2_oC20_36_b_a_b & a,b/a,c/a,y1,z1,x2,y2,z2,x3,y3,z3 --params=8.843,
      ↪ 0.618907610539,0.546760149271,0.214,0.23,0.218,0.121,0.642,
      ↪ 0.1763,0.1509,0.2898 & Cmc2_{1} C_{2v}^{12} #36 (ab^2) & oC20 &
      ↪  None & N2OSi2 & N2OSi2 & I. Idrestedt and C. Brosset, Acta
      ↪ Chem. Scand. 18, 1879-1886 (1964)
   1.00000000000000
   4.42150000000000  -2.73650000000000   0.00000000000000
   4.42150000000000   2.73650000000000   0.00000000000000
   0.00000000000000   0.00000000000000   4.83500000000000
     N     O     Si
     4     2     4
Direct
   0.09700000000000   0.33900000000000   0.64200000000000   N   (8b)
  -0.09700000000000  -0.33900000000000   1.14200000000000   N   (8b)
   0.33900000000000   0.09700000000000   1.14200000000000   N   (8b)
  -0.33900000000000  -0.09700000000000   0.64200000000000   N   (8b)
  -0.21400000000000   0.21400000000000   0.23000000000000   O   (4a)
   0.21400000000000  -0.21400000000000   0.73000000000000   O   (4a)
   0.02540000000000   0.32720000000000   0.28980000000000   Si  (8b)
  -0.02540000000000  -0.32720000000000   0.78980000000000   Si  (8b)
   0.32720000000000   0.02540000000000   0.78980000000000   Si  (8b)
  -0.32720000000000  -0.02540000000000   0.28980000000000   Si  (8b)
```

Bi_2GeO_5: A2BC5_oC32_36_b_a_a2b - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Bi2GeO5'
_chemical_formula_sum 'Bi2 Ge O5'

loop_
_publ_author_name
 'B. Aurivillius'
 'C.-I. Lindblom'
 'P. St\\'{e}nson'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 18
_journal_year 1964
_journal_page_first 1555
_journal_page_last 1557
_publ_Section_title
```

```
;
 The Crystal Structure of Bi$_{2}$GeO$_{5}$
;

_aflow_title 'Bi$_{2}$GeO$_{5}$ Structure'
_aflow_proto 'A2BC5_oC32_36_b_a_a2b'
_aflow_params 'a,b/a,c/a,y_{1},z_{1},y_{2},z_{2},x_{3},y_{3},z_{3},x_{4
      ↪ },y_{4},z_{4},x_{5},y_{5},z_{5}'
_aflow_params_values '15.69,0.350031867431,0.343084767368,0.68572,
      ↪ 0.20003,0.5803,0.4908,0.16758,0.21738,0.25,0.0949,0.1368,0.6482
      ↪ ,0.2541,0.4777,0.4621'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oC32'

_symmetry_space_group_name_H-M "C m c 21"
_symmetry_Int_Tables_number 36

_cell_length_a     15.69000
_cell_length_b     5.49200
_cell_length_c     5.38300
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,z+1/2
3 -x,y,z
4 x,-y,z+1/2
5 x+1/2,y+1/2,z
6 -x+1/2,-y+1/2,z+1/2
7 -x+1/2,y+1/2,z
8 x+1/2,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ge1 Ge 4 a 0.00000 0.68572 0.20003 1.00000
O1 O 4 a 0.00000 0.58030 0.49080 1.00000
Bi1 Bi 8 b 0.16758 0.21738 0.25000 1.00000
O2 O 8 b 0.09490 0.13680 0.64820 1.00000
O3 O 8 b 0.25410 0.47770 0.46210 1.00000
```

Bi_2GeO_5: A2BC5_oC32_36_b_a_a2b - POSCAR

```
A2BC5_oC32_36_b_a_a2b & a,b/a,c/a,y1,z1,y2,z2,x3,y3,z3,x4,y4,z4,x5,y5,z5
      ↪ --params=15.69,0.350031867431,0.343084767368,0.68572,0.20003,
      ↪ 0.5803,0.4908,0.16758,0.21738,0.25,0.0949,0.1368,0.6482,0.2541,
      ↪ 0.4777,0.4621 & Cmc2_{1} C_{2v}^{12} #36 (a^2b^3) & oC32 & None
      ↪ & Bi2GeO5 & Bi2GeO5 & B. Aurivillius and C.-I. Lindblom and P.
      ↪ St\'{e}nson, Acta Chem. Scand. 18, 1555-1557 (1964)
   1.00000000000000
   7.84500000000000  -2.74600000000000   0.00000000000000
   7.84500000000000   2.74600000000000   0.00000000000000
   0.00000000000000   0.00000000000000   5.38300000000000
     Bi    Ge    O
      4     2    10
Direct
  -0.04980000000000   0.38496000000000   0.25000000000000  Bi   (8b)
   0.04980000000000  -0.38496000000000   0.75000000000000  Bi   (8b)
   0.38496000000000  -0.04980000000000   0.75000000000000  Bi   (8b)
  -0.38496000000000   0.04980000000000   0.25000000000000  Bi   (8b)
  -0.68572000000000   0.68572000000000   0.20003000000000  Ge   (4a)
   0.68572000000000  -0.68572000000000   0.70003000000000  Ge   (4a)
  -0.58030000000000   0.58030000000000   0.49080000000000  O    (4a)
   0.58030000000000  -0.58030000000000   0.99080000000000  O    (4a)
  -0.04190000000000   0.23170000000000   0.64820000000000  O    (8b)
   0.04190000000000  -0.23170000000000   1.14820000000000  O    (8b)
   0.23170000000000  -0.04190000000000   1.14820000000000  O    (8b)
  -0.23170000000000   0.04190000000000   0.64820000000000  O    (8b)
  -0.22360000000000   0.73180000000000   0.46210000000000  O    (8b)
   0.22360000000000  -0.73180000000000   0.96210000000000  O    (8b)
   0.73180000000000  -0.22360000000000   0.96210000000000  O    (8b)
  -0.73180000000000   0.22360000000000   0.46210000000000  O    (8b)
```

Ni_3Si_2: A3B2_oC80_36_4a4b_2a3b - CIF

```
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_audit_creation_method FINDSYM

_chemical_name_mineral 'Ni3Si2'
_chemical_formula_sum 'Ni3 Si2'

loop_
_publ_author_name
 'G. Pilstr\"{o}m'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 15
_journal_year 1961
_journal_page_first 893
_journal_page_last 902
_publ_Section_title
;
```

```
 The Crystal Structure of Ni$_{3}$Si$_{2}$ with some Notes on Ni$_{5}
      ↪ $Si$_{2}$
;

_aflow_title 'Ni$_{3}$Si$_{2}$ Structure'
_aflow_proto 'A3B2_oC80_36_4a4b_2a3b'
_aflow_params 'a,b/a,c/a,y_{1},z_{1},y_{2},z_{2},y_{3},z_{3},y_{4},z_{4
      ↪ },y_{5},z_{5},y_{6},z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},z_{8},
      ↪ x_{9},y_{9},z_{9},x_{10},y_{10},z_{10},x_{11},y_{11},z_{11},x_{
      ↪ 12},y_{12},z_{12},x_{13},y_{13},z_{13}'
_aflow_params_values '12.22901,0.883554760361,0.566194646991,0.0,0.0,
      ↪ 0.2345,0.024,0.233,0.4,0.3814,0.714,0.157,0.712,0.409,0.218,
      ↪ 0.1732,0.1177,0.518,0.1723,0.1189,-0.1,0.1972,0.2467,0.217,
      ↪ 0.1824,0.4975,0.225,0.12,0.059,0.214,0.152,0.344,0.506,0.151,
      ↪ 0.343,-0.08'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oC80'

_symmetry_space_group_name_H-M "C m c 21"
_symmetry_Int_Tables_number 36

_cell_length_a     12.22901
_cell_length_b     10.80500
_cell_length_c     6.92400
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

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_space_group_symop_operation_xyz
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2 -x,-y,z+1/2
3 -x,y,z
4 x,-y,z+1/2
5 x+1/2,y+1/2,z
6 -x+1/2,-y+1/2,z+1/2
7 -x+1/2,y+1/2,z
8 x+1/2,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ni1 Ni 4 a 0.00000 0.00000 0.00000 1.00000
Ni2 Ni 4 a 0.00000 0.23450 0.02400 1.00000
Ni3 Ni 4 a 0.00000 0.23300 0.40000 1.00000
Ni4 Ni 4 a 0.00000 0.38140 0.71400 1.00000
Si1 Si 4 a 0.00000 0.15700 0.71200 1.00000
Si2 Si 4 a 0.00000 0.40900 0.21800 1.00000
Ni5 Ni 8 b 0.17320 0.11770 0.51800 1.00000
Ni6 Ni 8 b 0.17230 0.11890 -0.10000 1.00000
Ni7 Ni 8 b 0.19720 0.24670 0.21700 1.00000
Ni8 Ni 8 b 0.18240 0.49750 0.22500 1.00000
Si3 Si 8 b 0.12000 0.05900 0.21400 1.00000
Si4 Si 8 b 0.15200 0.34400 0.50600 1.00000
Si5 Si 8 b 0.15100 0.34300 -0.08000 1.00000
```

Ni_3Si_2: A3B2_oC80_36_4a4b_2a3b - POSCAR

```
A3B2_oC80_36_4a4b_2a3b & a,b/a,c/a,y1,z1,y2,z2,y3,z3,y4,z4,y5,z5,y6,z6,
      ↪ x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10,x11,y11,z11,x12,y12,z12,
      ↪ x13,y13,z13 --params=12.22901,0.883554760361,0.566194646991,0.0
      ↪ ,0.0,0.2345,0.024,0.233,0.4,0.3814,0.714,0.157,0.712,0.409,
      ↪ 0.218,0.1732,0.1177,0.518,0.1723,0.1189,-0.1,0.1972,0.2467,
      ↪ 0.217,0.1824,0.4975,0.225,0.12,0.059,0.214,0.152,0.344,0.506,
      ↪ 0.151,0.343,-0.08 & Cmc2_{1} C_{2v}^{12} #36 (a^6b^7) & oC80 &
      ↪ None & Ni3Si2 & Ni3Si2 & G. Pilstr\"{o}m, Acta Chem. Scand. 15,
      ↪ 893-902 (1961)
   1.00000000000000
   6.11450500000000  -5.40250000000000   0.00000000000000
   6.11450500000000   5.40250000000000   0.00000000000000
   0.00000000000000   0.00000000000000   6.92400000000000
     Ni    Si
     24    16
Direct
   0.00000000000000   0.00000000000000   0.00000000000000  Ni   (4a)
   0.00000000000000   0.00000000000000   0.50000000000000  Ni   (4a)
  -0.23450000000000   0.23450000000000   0.02400000000000  Ni   (4a)
   0.23450000000000  -0.23450000000000   0.52400000000000  Ni   (4a)
  -0.23300000000000   0.23300000000000   0.40000000000000  Ni   (4a)
   0.23300000000000  -0.23300000000000   0.90000000000000  Ni   (4a)
  -0.38140000000000   0.38140000000000   0.71400000000000  Ni   (4a)
   0.38140000000000  -0.38140000000000   1.21400000000000  Ni   (4a)
   0.05550000000000   0.29090000000000   0.51800000000000  Ni   (8b)
  -0.05550000000000  -0.29090000000000   1.01800000000000  Ni   (8b)
   0.29090000000000   0.05550000000000   1.01800000000000  Ni   (8b)
  -0.29090000000000  -0.05550000000000   0.51800000000000  Ni   (8b)
   0.05340000000000   0.29120000000000  -0.10000000000000  Ni   (8b)
  -0.05340000000000  -0.29120000000000   0.40000000000000  Ni   (8b)
   0.29120000000000   0.05340000000000   0.40000000000000  Ni   (8b)
  -0.29120000000000  -0.05340000000000  -0.10000000000000  Ni   (8b)
  -0.04950000000000   0.44390000000000   0.21700000000000  Ni   (8b)
   0.04950000000000  -0.44390000000000   0.71700000000000  Ni   (8b)
   0.44390000000000  -0.04950000000000   0.71700000000000  Ni   (8b)
  -0.44390000000000   0.04950000000000   0.21700000000000  Ni   (8b)
  -0.31510000000000   0.67990000000000   0.22500000000000  Ni   (8b)
   0.31510000000000  -0.67990000000000   0.72500000000000  Ni   (8b)
   0.67990000000000  -0.31510000000000   0.72500000000000  Ni   (8b)
  -0.67990000000000   0.31510000000000   0.22500000000000  Ni   (8b)
```

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 −0.15700000000000   0.15700000000000   0.71200000000000   Si   (4a)
  0.15700000000000  −0.15700000000000   1.21200000000000   Si   (4a)
 −0.40900000000000   0.40900000000000   0.21800000000000   Si   (4a)
  0.40900000000000  −0.40900000000000   0.71800000000000   Si   (4a)
  0.06100000000000   0.17900000000000   0.21400000000000   Si   (8b)
 −0.06100000000000  −0.17900000000000   0.71400000000000   Si   (8b)
  0.17900000000000   0.06100000000000   0.71400000000000   Si   (8b)
 −0.17900000000000  −0.06100000000000   0.21400000000000   Si   (8b)
 −0.19200000000000   0.49600000000000   0.50600000000000   Si   (8b)
  0.19200000000000  −0.49600000000000   1.00600000000000   Si   (8b)
  0.49600000000000  −0.19200000000000   1.00600000000000   Si   (8b)
 −0.49600000000000   0.19200000000000   0.50600000000000   Si   (8b)
 −0.19200000000000   0.49400000000000  −0.08000000000000   Si   (8b)
  0.19200000000000  −0.49400000000000   0.42000000000000   Si   (8b)
  0.49400000000000  −0.19200000000000   0.42000000000000   Si   (8b)
 −0.49400000000000   0.19200000000000  −0.08000000000000   Si   (8b)
```

Bertrandite ($Be_4Si_2O_7(OH)_2$, $S4_6$): A4B7C2D2_oC60_36_2b_a3b_2a_b - CIF

```
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_chemical_name_mineral 'Bertrandite'
_chemical_formula_sum 'Be4 O7 (OH)2 Si2'

loop_
_publ_author_name
 'R. M. Hazen'
 'A. Y. Au'
_journal_name_full_name
;
 Physics and Chemistry of Minerals
;
_journal_volume 13
_journal_year 1986
_journal_page_first 69
_journal_page_last 78
_publ_Section_title
;
 High-pressure crystal chemistry of phenakite (Be$_{2}$SiO$_{4}$) and
      ↪ bertrandite (Be$_{4}$Si$_{2}$O$_{7}$(OH)$_{2}$)
;

# Found in The power of databases: the RRUFF project, 2015

_aflow_title 'Bertrandite (Be$_{4}$Si$_{2}$O$_{7}$(OH)$_{2}$, $S4_{6}$)
      ↪ Structure'
_aflow_proto 'A4B7C2D2_oC60_36_2b_a3b_2a_b'
_aflow_params 'a,b/a,c/a,y_{1},z_{1},y_{2},z_{2},y_{3},z_{3},x_{4},y_{4
      ↪ },z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{6},x_{7},y_{7},z_{7},
      ↪ x_{8},y_{8},z_{8},x_{9},y_{9},z_{9}'
_aflow_params_values '8.7135,1.75222356114,0.524278418546,0.5852,0.5951,
      ↪ 0.7553,0.0886,0.0872,0.0993,0.1735,0.0527,0.1628,0.3264,0.2203,
      ↪ 0.1517,0.2897,0.1244,0.0,0.2101,0.0431,0.5074,0.2934,0.2093,
      ↪ 0.5024,0.3254,0.1144,0.6523'
_aflow_Strukturbericht '$S4_{6}$'
_aflow_Pearson 'oC60'

_symmetry_space_group_name_H-M "C m c 21"
_symmetry_Int_Tables_number 36

_cell_length_a     8.71350
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_cell_length_c     4.56830
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_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

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_space_group_symop_id
_space_group_symop_operation_xyz
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3 -x,y,z
4 x,-y,z+1/2
5 x+1/2,y+1/2,z
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7 -x+1/2,y+1/2,z
8 x+1/2,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 4 a 0.00000 0.58520 0.59510 1.00000
OH1 OH 4 a 0.00000 0.75530 0.08860 1.00000
OH2 OH 4 a 0.00000 0.08720 0.09930 1.00000
Be1 Be 8 b 0.17350 0.05270 0.16280 1.00000
Be2 Be 8 b 0.32640 0.22030 0.15170 1.00000
O2 O 8 b 0.28970 0.12440 0.00000 1.00000
O3 O 8 b 0.21010 0.04310 0.50740 1.00000
O4 O 8 b 0.29340 0.20930 0.50240 1.00000
Si1 Si 8 b 0.32540 0.11440 0.65230 1.00000
```

Bertrandite ($Be_4Si_2O_7(OH)_2$, $S4_6$): A4B7C2D2_oC60_36_2b_a3b_2a_b - POSCAR

```
A4B7C2D2_oC60_36_2b_a3b_2a_b & a,b/a,c/a,y1,z1,y2,z2,y3,z3,x4,y4,z4,x5,
      ↪ y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9 --params=8.7135,
      ↪ 1.75222356114,0.524278418546,0.5852,0.5951,0.7553,0.0886,0.0872
      ↪ ,0.0993,0.1735,0.0527,0.1628,0.3264,0.2203,0.1517,0.2897,0.1244
      ↪ ,0.0,0.2101,0.0431,0.5074,0.2934,0.2093,0.5024,0.3254,0.1144,
      ↪ 0.6523 & Cmc2_{1} C_{2v}^{12} #36 (a^3b^6) & oC60 & $S4_{6}$ &
      ↪ Be4O7(OH)2Si2 & Bertrandite & R. M. Hazen and A. Y. Au, Phys.
      ↪ Chem. Miner. 13, 69-78 (1986)
   1.00000000000000
   4.35675000000000  -7.63400000000000   0.00000000000000
   4.35675000000000   7.63400000000000   0.00000000000000
   0.00000000000000   0.00000000000000   4.56830000000000
    Be    O    OH    Si
     8   14     4     4
Direct
   0.12080000000000   0.22620000000000   0.16280000000000   Be   (8b)
  -0.12080000000000  -0.22620000000000   0.66280000000000   Be   (8b)
   0.22620000000000   0.12080000000000   0.66280000000000   Be   (8b)
  -0.22620000000000  -0.12080000000000   0.16280000000000   Be   (8b)
   0.10610000000000   0.54670000000000   0.15170000000000   Be   (8b)
  -0.10610000000000  -0.54670000000000   0.65170000000000   Be   (8b)
   0.54670000000000   0.10610000000000   0.65170000000000   Be   (8b)
  -0.54670000000000  -0.10610000000000   0.15170000000000   Be   (8b)
  -0.58520000000000   0.58520000000000   0.59510000000000   O    (4a)
   0.58520000000000  -0.58520000000000   1.09510000000000   O    (4a)
   0.16530000000000   0.41410000000000   0.00000000000000   O    (8b)
  -0.16530000000000  -0.41410000000000   0.50000000000000   O    (8b)
   0.41410000000000   0.16530000000000   0.50000000000000   O    (8b)
  -0.41410000000000  -0.16530000000000   0.00000000000000   O    (8b)
   0.16700000000000   0.25320000000000   0.50740000000000   O    (8b)
  -0.16700000000000  -0.25320000000000   1.00740000000000   O    (8b)
   0.25320000000000   0.16700000000000   1.00740000000000   O    (8b)
  -0.25320000000000  -0.16700000000000   0.50740000000000   O    (8b)
   0.08410000000000   0.50270000000000   0.50240000000000   O    (8b)
  -0.08410000000000  -0.50270000000000   1.00240000000000   O    (8b)
   0.50270000000000   0.08410000000000   1.00240000000000   O    (8b)
  -0.50270000000000  -0.08410000000000   0.50240000000000   O    (8b)
  -0.75530000000000   0.75530000000000   0.08860000000000   OH   (4a)
   0.75530000000000  -0.75530000000000   0.58860000000000   OH   (4a)
  -0.08720000000000   0.08720000000000   0.09930000000000   OH   (4a)
   0.08720000000000  -0.08720000000000   0.59930000000000   OH   (4a)
   0.21100000000000   0.43980000000000   0.65230000000000   Si   (8b)
  -0.21100000000000  -0.43980000000000   1.15230000000000   Si   (8b)
   0.43980000000000   0.21100000000000   1.15230000000000   Si   (8b)
  -0.43980000000000  -0.21100000000000   0.65230000000000   Si   (8b)
```

MoP_2: AB2_oC12_36_a_2a - CIF

```
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_audit_creation_method FINDSYM

_chemical_name_mineral 'MoP2'
_chemical_formula_sum 'Mo P2'

loop_
_publ_author_name
 'S. Rundqvist'
 'T. Lundstr\"{o}m'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 17
_journal_year 1963
_journal_page_first 37
_journal_page_last 46
_publ_Section_title
;
 X-Ray Studies of Molybdenum and Tungsten Phosphides
;

_aflow_title 'MoP$_{2}$ Structure'
_aflow_proto 'AB2_oC12_36_a_2a'
_aflow_params 'a,b/a,c/a,y_{1},z_{1},y_{2},z_{2},y_{3},z_{3}'
_aflow_params_values '3.145,3.55612082671,1.84702384738,0.0934,0.0,0.294
      ↪ ,0.803,0.426,0.121'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oC12'

_symmetry_space_group_name_H-M "C m c 21"
_symmetry_Int_Tables_number 36

_cell_length_a     3.14500
_cell_length_b     11.18400
_cell_length_c     5.80889
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

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_space_group_symop_id
_space_group_symop_operation_xyz
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2 -x,-y,z+1/2
3 -x,y,z
4 x,-y,z+1/2
5 x+1/2,y+1/2,z
6 -x+1/2,-y+1/2,z+1/2
7 -x+1/2,y+1/2,z
8 x+1/2,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
```

```
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
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P1 P 4 a 0.00000 0.29400 0.80300 1.00000
P2 P 4 a 0.00000 0.42600 0.12100 1.00000
```

MoP_2: AB2_oC12_36_a_2a - POSCAR

```
AB2_oC12_36_a_2a & a,b/a,c/a,y1,z1,y2,z2,y3,z3 --params=3.145,
      ↪ 3.55612082671,1.84702384738,0.0934,0.0,0.294,0.803,0.426,0.121
      ↪ & Cmc2_{1} C_{2v}^{12} #36 (a^3) & oC12 & None & MoP2 & MoP2 &
      ↪ S. Rundqvist and T. Lundstr\"{o}m, Acta Chem. Scand. 17, 37-46
      ↪ (1963)
   1.00000000000000
   1.57250000000000  -5.59200000000000   0.00000000000000
   1.57250000000000   5.59200000000000   0.00000000000000
   0.00000000000000   0.00000000000000   5.80889000000000
     Mo     P
      2     4
Direct
  -0.09340000000000   0.09340000000000   0.00000000000000  Mo   (4a)
   0.09340000000000  -0.09340000000000   0.50000000000000  Mo   (4a)
  -0.29400000000000   0.29400000000000   0.80300000000000  P    (4a)
   0.29400000000000  -0.29400000000000   1.30300000000000  P    (4a)
  -0.42600000000000   0.42600000000000   0.12100000000000  P    (4a)
   0.42600000000000  -0.42600000000000   0.62100000000000  P    (4a)
```

α-Potassium Nitrate (KNO_3) II: ABC3_oC80_36_2ab_2ab_2a5b - CIF

```
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_chemical_name_mineral 'KNO3'
_chemical_formula_sum 'K N O3'

loop_
_publ_author_name
 'G. Adiwidjaja'
 'D. Pohl'
_journal_name_full_name
;
 Acta Crystallographica Section C: Structural Chemistry
;
_journal_volume 59
_journal_year 2003
_journal_page_first i139
_journal_page_last i140
_publ_Section_title
;
 Superstructure of $\alpha$-phase potassium nitrate
;

_aflow_title '$\alpha$-Potassium Nitrate (KNO$_{3}$) II Structure'
_aflow_proto 'ABC3_oC80_36_2ab_2ab_2a5b'
_aflow_params 'a,b/a,c/a,y_{1},z_{1},y_{2},z_{2},y_{3},z_{3},y_{4},z_{4
      ↪ },y_{5},z_{5},y_{6},z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},z_{8},
      ↪ x_{9},y_{9},z_{9},x_{10},y_{10},z_{10},x_{11},y_{11},z_{11},x_{
      ↪ 12},y_{12},z_{12},x_{13},y_{13},z_{13}'
_aflow_params_values '10.825,1.69524249423,0.594457274827,0.41703,0.2604
      ↪ ,-0.08321,0.2614,0.2483,0.4215,0.7469,0.4208,0.1807,0.414,
      ↪ 0.6788,0.4138,0.2498,0.1667,0.25104,0.2497,0.4979,0.0911,0.0999
      ↪ ,0.283,0.4206,0.3987,0.2809,0.4248,0.6506,0.0313,0.0984,0.8479,
      ↪ 0.0321,0.0896,0.2478,0.4296,0.0937'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oC80'

_symmetry_space_group_name_H-M "C m c 21"
_symmetry_Int_Tables_number 36

_cell_length_a     10.82500
_cell_length_b     18.35100
_cell_length_c     6.43500
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,z+1/2
3 -x,y,z
4 x,-y,z+1/2
5 x+1/2,y+1/2,z
6 -x+1/2,-y+1/2,z+1/2
7 -x+1/2,y+1/2,z
8 x+1/2,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
K1 K 4 a 0.00000 0.41703 0.26040 1.00000
K2 K 4 a 0.00000 -0.08321 0.26140 1.00000
N1 N 4 a 0.00000 0.24830 0.42150 1.00000
N2 N 4 a 0.00000 0.74690 0.42080 1.00000
O1 O 4 a 0.00000 0.18070 0.41400 1.00000
O2 O 4 a 0.00000 0.67880 0.41380 1.00000
K3 K 8 b 0.24980 0.16670 0.25104 1.00000
N3 N 8 b 0.24970 0.49790 0.09110 1.00000
O3 O 8 b 0.09990 0.28300 0.42060 1.00000
O4 O 8 b 0.39870 0.28090 0.42480 1.00000
O5 O 8 b 0.65060 0.03130 0.09840 1.00000
O6 O 8 b 0.84790 0.03210 0.08960 1.00000
O7 O 8 b 0.24780 0.42960 0.09370 1.00000
```

α-Potassium Nitrate (KNO_3) II: ABC3_oC80_36_2ab_2ab_2a5b - POSCAR

```
ABC3_oC80_36_2ab_2ab_2a5b & a,b/a,c/a,y1,z1,y2,z2,y3,z3,y4,z4,y5,z5,y6,
      ↪ z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10,x11,y11,z11,x12,y12,
      ↪ z12,x13,y13,z13 --params=10.825,1.69524249423,0.594457274827,
      ↪ 0.41703,0.2604,-0.08321,0.2614,0.2483,0.4215,0.7469,0.4208,
      ↪ 0.1807,0.414,0.6788,0.4138,0.2498,0.1667,0.25104,0.2497,0.4979,
      ↪ 0.0911,0.0999,0.283,0.4206,0.3987,0.2809,0.4248,0.6506,0.0313,
      ↪ 0.0984,0.8479,0.0321,0.0896,0.2478,0.4296,0.0937 & Cmc2_{1} C_{
      ↪ 2v}^{12} #36 (a^6b^7) & oC80 & None & KNO3 & KNO3 & G.
      ↪ Adiwidjaja and D. Pohl, Acta Crystallogr. C 59, i139-i140 (2003
      ↪ )
   1.00000000000000
   5.41250000000000  -9.17550000000000   0.00000000000000
   5.41250000000000   9.17550000000000   0.00000000000000
   0.00000000000000   0.00000000000000   6.43500000000000
     K     N     O
     8     8    24
Direct
  -0.41703000000000   0.41703000000000   0.26040000000000  K    (4a)
   0.41703000000000  -0.41703000000000   0.76040000000000  K    (4a)
   0.08321000000000  -0.08321000000000   0.26140000000000  K    (4a)
  -0.08321000000000   0.08321000000000   0.76140000000000  K    (4a)
   0.08310000000000   0.41650000000000   0.25104000000000  K    (8b)
  -0.08310000000000  -0.41650000000000   0.75104000000000  K    (8b)
   0.41650000000000   0.08310000000000   0.75104000000000  K    (8b)
  -0.41650000000000  -0.08310000000000   0.25104000000000  K    (8b)
  -0.24830000000000   0.24830000000000   0.42150000000000  N    (4a)
   0.24830000000000  -0.24830000000000   0.92150000000000  N    (4a)
  -0.74690000000000   0.74690000000000   0.42080000000000  N    (4a)
   0.74690000000000  -0.74690000000000   0.92080000000000  N    (4a)
  -0.24820000000000   0.74760000000000   0.09110000000000  N    (8b)
   0.24820000000000  -0.74760000000000   0.59110000000000  N    (8b)
   0.74760000000000  -0.24820000000000   0.59110000000000  N    (8b)
  -0.74760000000000   0.24820000000000   0.09110000000000  N    (8b)
  -0.18070000000000   0.18070000000000   0.41400000000000  O    (4a)
   0.18070000000000  -0.18070000000000   0.91400000000000  O    (4a)
  -0.67880000000000   0.67880000000000   0.41380000000000  O    (4a)
   0.67880000000000  -0.67880000000000   0.91380000000000  O    (4a)
  -0.18310000000000   0.38290000000000   0.42060000000000  O    (8b)
   0.18310000000000  -0.38290000000000   0.92060000000000  O    (8b)
   0.38290000000000  -0.18310000000000   0.92060000000000  O    (8b)
  -0.38290000000000   0.18310000000000   0.42060000000000  O    (8b)
   0.11780000000000   0.67960000000000   0.42480000000000  O    (8b)
  -0.11780000000000  -0.67960000000000   0.92480000000000  O    (8b)
   0.67960000000000   0.11780000000000   0.92480000000000  O    (8b)
  -0.67960000000000  -0.11780000000000   0.42480000000000  O    (8b)
   0.61930000000000   0.68190000000000   0.09840000000000  O    (8b)
  -0.61930000000000  -0.68190000000000   0.59840000000000  O    (8b)
   0.68190000000000   0.61930000000000   0.59840000000000  O    (8b)
  -0.68190000000000  -0.61930000000000   0.09840000000000  O    (8b)
   0.81580000000000   0.88000000000000   0.08960000000000  O    (8b)
  -0.81580000000000  -0.88000000000000   0.58960000000000  O    (8b)
   0.88000000000000   0.81580000000000   0.58960000000000  O    (8b)
  -0.88000000000000  -0.81580000000000   0.08960000000000  O    (8b)
  -0.18180000000000   0.67740000000000   0.09370000000000  O    (8b)
   0.18180000000000  -0.67740000000000   0.59370000000000  O    (8b)
   0.67740000000000  -0.18180000000000   0.59370000000000  O    (8b)
  -0.67740000000000   0.18180000000000   0.09370000000000  O    (8b)
```

Ta_3Ti_{13} (BCC SQS-16): A3B13_oC32_38_ac_a2bcdef - CIF

```
# AFLOW.org Repositories
# TaTi/A3B13_oC32_38_ac_a2bcdef-001.AB params=6.5421326204,1.41421356237
      ↪ ,1.41421356237,0.5,0.0,0.5,0.0,0.75,0.25,0.75,0.75,0.25,0.25,
      ↪ 0.25,0.25,0.75,0.75,0.0 SG=38 [ANRL doi: 10.1016/
      ↪ j.commatsci.2017.01.017 (part 1), doi: 10.1016/
      ↪ j.commatsci.2018.10.043 (part 2)]
data_TaTi
_pd_phase_name A3B13_oC32_38_ac_a2bcdef-001.AB

_chemical_name_mineral 'Ta3Ti13'
_chemical_formula_sum 'Ta3 Ti13'

loop_
_publ_author_name
 'T. Chakraborty'
 'J. Rogal'
 'R. Drautz'
_journal_name_full_name
;
 Physical Review B
;
_journal_volume 94
_journal_year 2016
_journal_page_first 224104
_journal_page_last 224104
_publ_Section_title
;
 Unraveling the composition dependence of the martensitic transformation
      ↪ temperature: A first-principles study of Ti-Ta alloys
;

_aflow_title 'Ta$_{3}$Ti$_{13}$ (BCC SQS-16) Structure'
_aflow_proto 'A3B13_oC32_38_ac_a2bcdef'
_aflow_params 'a,b/a,c/a,z_{1},z_{2},z_{3},z_{4},x_{5},z_{5},x_{6},z_{6
      ↪ },y_{7},z_{7},y_{8},z_{8},x_{9},y_{9},z_{9}'
```

```
_aflow_params_values '6.5421326204,1.41421356237,1.41421356237,0.5,0.0,
    ↪ 0.5,0.0,0.75,0.75,0.75,0.25,0.75,0.75,0.75,0.75,0.75,0.25,0.0'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oC32'

_cell_length_a 6.5421326204
_cell_length_b 9.2519726786
_cell_length_c 9.2519726786
_cell_angle_alpha 90.0000000000
_cell_angle_beta 90.0000000000
_cell_angle_gamma 90.0000000000
_symmetry_space_group_name_H-M 'Amm2'
_symmetry_Int_Tables_Number 38
loop_
_symmetry_equiv_pos_site_id
_symmetry_equiv_pos_as_xyz
 1 x,y,z
 2 -x,-y,z
 3 x,-y,z
 4 -x,y,z
 5 x,y+1/2,z+1/2
 6 -x,-y+1/2,z+1/2
 7 x,-y+1/2,z+1/2
 8 -x,y+1/2,z+1/2
loop_
 _atom_site_label
 _atom_site_occupancy
 _atom_site_fract_x
 _atom_site_fract_y
 _atom_site_fract_z
 _atom_site_thermal_displace_type
 _atom_site_B_iso_or_equiv
 _atom_site_type_symbol
 _atom_site_symmetry_multiplicity
 _atom_site_Wyckoff_label
Ta1 1.0000000000 0.0000000000 -0.0000000000 0.5000000000 Biso 1.0 Ta 2 a
Ti1 1.0000000000 0.0000000000 0.0000000000 0.0000000000 Biso 1.0 Ti 2 a
Ti2 1.0000000000 0.5000000000 -0.0000000000 0.5000000000 Biso 1.0 Ti 2 b
Ti3 1.0000000000 0.5000000000 0.0000000000 0.0000000000 Biso 1.0 Ti 2 b
Ta2 1.0000000000 0.7500000000 -0.0000000000 0.7500000000 Biso 1.0 Ta 4 c
Ti4 1.0000000000 0.7500000000 -0.0000000000 0.2500000000 Biso 1.0 Ti 4 c
Ti5 1.0000000000 0.0000000000 0.7500000000 0.7500000000 Biso 1.0 Ti 4 d
Ti6 1.0000000000 0.5000000000 0.7500000000 0.7500000000 Biso 1.0 Ti 4 e
Ti7 1.0000000000 0.7500000000 0.2500000000 0.0000000000 Biso 1.0 Ti 8 f
```

Ta_3Ti_{13} (BCC SQS-16): A3B13_oC32_38_ac_a2bcdef - POSCAR

```
A3B13_oC32_38_ac_a2bcdef & a,b/a,c/a,z1,z2,z3,z4,x5,z5,x6,z6,y7,z7,y8,z8
    ↪ ,x9,y9,z9 --params=6.5421326204,1.41421356237,1.41421356237,0.5
    ↪ ,0.0,0.5,0.0,0.75,0.75,0.75,0.25,0.75,0.75,0.75,0.75,0.75,0.25,
    ↪ 0.0 & Amm2 C_{2v}^{14} #38 (a^2b^2c^2def) & oC32 & None &
    ↪ Ta3Ti13 & Ta3Ti13 & T. Chakraborty and J. Rogal and R. Drautz,
    ↪ Phys. Rev. B 94, 224104(2016)
   1.00000000000000
   6.54213262040000   0.00000000000000   0.00000000000000
   0.00000000000000   4.62598633930000  -4.62598633930000
   0.00000000000000   4.62598633930000   4.62598633930000
     Ta    Ti
      3    13
Direct
   0.00000000000000  -0.50000000000000   0.50000000000000   Ta   (2a)
   0.75000000000000  -0.75000000000000   0.75000000000000   Ta   (4c)
  -0.75000000000000  -0.75000000000000   0.75000000000000   Ta   (4c)
   0.00000000000000   0.00000000000000   0.00000000000000   Ti   (2a)
   0.50000000000000  -0.50000000000000   0.50000000000000   Ti   (2b)
   0.50000000000000   0.00000000000000   0.00000000000000   Ti   (2b)
   0.75000000000000  -0.25000000000000   0.25000000000000   Ti   (4c)
  -0.75000000000000  -0.25000000000000   0.25000000000000   Ti   (4c)
   0.00000000000000   0.00000000000000   1.50000000000000   Ti   (4d)
   0.00000000000000  -1.50000000000000   0.00000000000000   Ti   (4d)
   0.50000000000000   0.00000000000000   1.50000000000000   Ti   (4e)
   0.50000000000000  -1.50000000000000   0.00000000000000   Ti   (4e)
   0.75000000000000   0.25000000000000   0.25000000000000   Ti   (8f)
  -0.75000000000000  -0.25000000000000  -0.25000000000000   Ti   (8f)
   0.75000000000000  -0.25000000000000  -0.25000000000000   Ti   (8f)
  -0.75000000000000   0.25000000000000   0.25000000000000   Ti   (8f)
```

Ta_3Ti_5 (BCC SQS-16): A3B5_oC32_38_abce_abcdf - CIF

```
# AFLOW.org Repositories
# TaTi/A3B5_oC32_38_abce_abcdf-001.AB params=6.5421326204,1.41421356237,
    ↪ 1.41421356237,0.5,-0.0,-0.0,0.5,0.25,0.75,0.25,0.25,0.75,0.25,
    ↪ 0.75,0.25,0.25,0.75,-0.0 SG=38 [ANRL doi: 10.1016/
    ↪ j.commatsci.2017.01.017 (part 1), doi: 10.1016/
    ↪ j.commatsci.2018.10.043 (part 2)]
data_TaTi
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_chemical_name_mineral 'Ta3Ti5'
_chemical_formula_sum 'Ta3 Ti5'

loop_
_publ_author_name
 'T. Chakraborty'
 'J. Rogal'
 'R. Drautz'
_journal_name_full_name
;
 Physical Review B
;
_journal_volume 94
_journal_year 2016
_journal_page_first 224104
_journal_page_last 224104
_publ_Section_title
;
 Unraveling the composition dependence of the martensitic transformation
    ↪ temperature: A first-principles study of Ti-Ta alloys
;

_aflow_title 'Ta${_}{3}$Ti${_}{5}$ (BCC SQS-16) Structure'
_aflow_proto 'A3B5_oC32_38_abce_abcdf'
_aflow_params 'a,b/a,c/a,z_{1},z_{2},z_{3},z_{4},x_{5},z_{5},x_{6},z_{6
    ↪ },y_{7},z_{7},y_{8},z_{8},x_{9},y_{9},z_{9}'
_aflow_params_values '6.5421326204,1.41421356237,1.41421356237,0.5,0.0,
    ↪ 0.0,0.5,0.25,0.25,0.25,0.75,0.75,0.25,0.75,0.25,0.25,0.25,0.0'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oC32'

_cell_length_a 6.5421326204
_cell_length_b 9.2519726786
_cell_length_c 9.2519726786
_cell_angle_alpha 90.0000000000
_cell_angle_beta 90.0000000000
_cell_angle_gamma 90.0000000000
_symmetry_space_group_name_H-M 'Amm2'
_symmetry_Int_Tables_Number 38
loop_
_symmetry_equiv_pos_site_id
_symmetry_equiv_pos_as_xyz
 1 x,y,z
 2 -x,-y,z
 3 x,-y,z
 4 -x,y,z
 5 x,y+1/2,z+1/2
 6 -x,-y+1/2,z+1/2
 7 x,-y+1/2,z+1/2
 8 -x,y+1/2,z+1/2
loop_
 _atom_site_label
 _atom_site_occupancy
 _atom_site_fract_x
 _atom_site_fract_y
 _atom_site_fract_z
 _atom_site_thermal_displace_type
 _atom_site_B_iso_or_equiv
 _atom_site_type_symbol
 _atom_site_symmetry_multiplicity
 _atom_site_Wyckoff_label
Ta1 1.0000000000 0.0000000000 -0.0000000000 0.5000000000 Biso 1.0 Ta 2 a
Ti1 1.0000000000 0.0000000000 0.0000000000 0.0000000000 Biso 1.0 Ti 2 a
Ta2 1.0000000000 0.5000000000 0.0000000000 0.0000000000 Biso 1.0 Ta 2 b
Ti2 1.0000000000 0.5000000000 -0.0000000000 0.5000000000 Biso 1.0 Ti 2 b
Ta3 1.0000000000 0.2500000000 -0.0000000000 0.2500000000 Biso 1.0 Ta 4 c
Ti3 1.0000000000 0.2500000000 -0.0000000000 0.7500000000 Biso 1.0 Ti 4 c
Ti4 1.0000000000 0.0000000000 0.7500000000 0.2500000000 Biso 1.0 Ti 4 d
Ta4 1.0000000000 0.5000000000 0.7500000000 0.2500000000 Biso 1.0 Ta 4 e
Ti5 1.0000000000 0.2500000000 0.2500000000 0.0000000000 Biso 1.0 Ti 8 f
```

Ta_3Ti_5 (BCC SQS-16): A3B5_oC32_38_abce_abcdf - POSCAR

```
A3B5_oC32_38_abce_abcdf & a,b/a,c/a,z1,z2,z3,z4,x5,z5,x6,z6,y7,z7,y8,z8,
    ↪ x9,y9,z9 --params=6.5421326204,1.41421356237,1.41421356237,0.5,
    ↪ 0.0,0.0,0.5,0.25,0.25,0.25,0.75,0.75,0.25,0.75,0.25,0.25,0.25,
    ↪ 0.0 & Amm2 C_{2v}^{14} #38 (a^2b^2c^2def) & oC32 & None &
    ↪ Ta3Ti5 & Ta3Ti5 & T. Chakraborty and J. Rogal and R. Drautz,
    ↪ Phys. Rev. B 94, 224104(2016)
   1.00000000000000
   6.54213262040000   0.00000000000000   0.00000000000000
   0.00000000000000   4.62598633930000  -4.62598633930000
   0.00000000000000   4.62598633930000   4.62598633930000
     Ta    Ti
      6    10
Direct
   0.00000000000000  -0.50000000000000   0.50000000000000   Ta   (2a)
   0.50000000000000   0.00000000000000   0.00000000000000   Ta   (2b)
   0.25000000000000  -0.25000000000000   0.25000000000000   Ta   (4c)
  -0.25000000000000  -0.25000000000000   0.25000000000000   Ta   (4c)
   0.50000000000000   0.50000000000000   1.00000000000000   Ta   (4e)
   0.50000000000000  -1.00000000000000  -0.50000000000000   Ta   (4e)
   0.00000000000000   0.00000000000000   0.00000000000000   Ti   (2a)
   0.50000000000000  -0.50000000000000   0.50000000000000   Ti   (2b)
   0.25000000000000  -0.75000000000000   0.75000000000000   Ti   (4c)
  -0.25000000000000  -0.75000000000000   0.75000000000000   Ti   (4c)
   0.00000000000000   0.50000000000000   1.00000000000000   Ti   (4d)
   0.00000000000000  -1.00000000000000  -0.50000000000000   Ti   (4d)
   0.25000000000000   0.25000000000000   0.25000000000000   Ti   (8f)
  -0.25000000000000  -0.25000000000000  -0.25000000000000   Ti   (8f)
   0.25000000000000  -0.25000000000000  -0.25000000000000   Ti   (8f)
  -0.25000000000000   0.25000000000000   0.25000000000000   Ti   (8f)
```

$NaNb_6O_{15}F$: ABC6D15_oC46_38_b_b_2a2d_2ab4d2e - CIF

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_chemical_formula_sum 'F Na Nb6 O15'

loop_
_publ_author_name
 'S. Andersson'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 19
_journal_year 1965
_journal_page_first 2285
```

```
_journal_page_last 2290
_publ_Section_title
;
 The Crystal Structure of NaNb$_{6}$O$_{15}$F and NaNb$_{6}$O$_{15}$OH
;

_aflow_title 'NaNb$_{6}$O$_{15}$F Structure'
_aflow_proto 'ABC6D15_oC46_38_b_b_2a2d_2ab4d2e'
_aflow_params 'a,b/a,c/a,z_{1},z_{2},z_{3},z_{4},z_{5},z_{6},z_{7},y_{8
      ↪ },z_{8},y_{9},z_{9},y_{10},z_{10},y_{11},z_{11},y_{12},z_{12},
      ↪ y_{13},z_{13},y_{14},z_{14},y_{15},z_{15}'
_aflow_params_values '3.949,2.58090655862,3.7277791846,0.0,0.2142,0.856,
      ↪ 0.519,0.0,0.404,0.226,0.3186,0.062,0.3153,0.3154,0.119,0.105,
      ↪ 0.133,0.311,0.305,0.447,0.351,0.178,0.324,0.05,0.323,0.307'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oC46'

_symmetry_space_group_name_H-M "A m m 2"
_symmetry_Int_Tables_number 38

_cell_length_a      3.94900
_cell_length_b      10.19200
_cell_length_c      14.72100
_cell_angle_alpha   90.00000
_cell_angle_beta    90.00000
_cell_angle_gamma   90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,z
3 x,-y,z
4 -x,y,z
5 x,y+1/2,z+1/2
6 -x,-y+1/2,z+1/2
7 x,-y+1/2,z+1/2
8 -x,y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
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Nb2 Nb 2 a 0.00000 0.00000 0.21420 1.00000
O1 O 2 a 0.00000 0.00000 0.85600 1.00000
O2 O 2 a 0.00000 0.00000 0.51900 1.00000
F1 F 2 b 0.50000 0.00000 0.00000 1.00000
Na1 Na 2 b 0.50000 0.00000 0.40400 1.00000
O3 O 2 b 0.50000 0.00000 0.22600 1.00000
Nb3 Nb 4 d 0.00000 0.31860 0.06200 1.00000
Nb4 Nb 4 d 0.00000 0.31530 0.31540 1.00000
O4 O 4 d 0.00000 0.11900 0.10500 1.00000
O5 O 4 d 0.00000 0.13300 0.31100 1.00000
O6 O 4 d 0.00000 0.30500 0.44700 1.00000
O7 O 4 d 0.00000 0.35100 0.17800 1.00000
O8 O 4 e 0.50000 0.32400 0.05000 1.00000
O9 O 4 e 0.50000 0.32300 0.30700 1.00000
```

$NaNb_6O_{15}F$: ABC6D15_oC46_38_b_b_2a2d_2ab4d2e - POSCAR

```
ABC6D15_oC46_38_b_b_2a2d_2ab4d2e & a,b/a,c/a,z1,z2,z3,z4,z5,z6,z7,y8,z8,
      ↪ y9,z9,y10,z10,y11,z11,y12,z12,y13,z13,y14,z14,y15,z15 --params=
      ↪ 3.949,2.58090655862,3.7277791846,0.0,0.2142,0.856,0.519,0.0,
      ↪ 0.404,0.226,0.3186,0.062,0.3153,0.3154,0.119,0.105,0.133,0.311,
      ↪ 0.305,0.447,0.351,0.178,0.324,0.05,0.323,0.307 & Amm2 C_{2v}^{
      ↪ 14} #38 (a^4b^3d^6e^2) & oC46 & None & FNaNb6O15 & FNaNb6O15 &
      ↪ S. Andersson, Acta Chem. Scand. 19, 2285-2290 (1965)
   1.00000000000000
   3.94900000000000    0.00000000000000    0.00000000000000
   0.00000000000000    5.09600000000000   -7.36050000000000
   0.00000000000000    5.09600000000000    7.36050000000000
     F     Na     Nb     O
     1      1      6     15
Direct
   0.50000000000000    0.00000000000000    0.00000000000000    F    (2b)
   0.50000000000000   -0.40400000000000    0.40400000000000    Na   (2b)
   0.00000000000000    0.00000000000000    0.00000000000000    Nb   (2a)
   0.00000000000000   -0.21420000000000    0.21420000000000    Nb   (2a)
   0.00000000000000    0.25660000000000    0.38060000000000    Nb   (4d)
   0.00000000000000   -0.38060000000000   -0.25660000000000    Nb   (4d)
   0.00000000000000   -0.00010000000000    0.63070000000000    Nb   (4d)
   0.00000000000000   -0.63070000000000    0.00010000000000    Nb   (4d)
   0.00000000000000   -0.85600000000000    0.85600000000000    O    (2a)
   0.00000000000000   -0.51900000000000    0.51900000000000    O    (2a)
   0.50000000000000   -0.22600000000000    0.22600000000000    O    (2b)
   0.00000000000000    0.01400000000000    0.22400000000000    O    (4d)
   0.00000000000000   -0.22400000000000   -0.01400000000000    O    (4d)
   0.00000000000000   -0.17800000000000    0.44400000000000    O    (4d)
   0.00000000000000   -0.44400000000000    0.17800000000000    O    (4d)
   0.00000000000000   -0.14200000000000    0.75200000000000    O    (4d)
   0.00000000000000   -0.75200000000000    0.14200000000000    O    (4d)
   0.00000000000000    0.17300000000000    0.52900000000000    O    (4d)
   0.00000000000000   -0.52900000000000   -0.17300000000000    O    (4d)
   0.50000000000000    0.27400000000000    0.37400000000000    O    (4e)
   0.50000000000000   -0.37400000000000   -0.27400000000000    O    (4e)
   0.50000000000000    0.01600000000000    0.63000000000000    O    (4e)
   0.50000000000000   -0.63000000000000   -0.01600000000000    O    (4e)
```

$Rb_2Mo_2O_7$: A2B7C2_oC88_40_abc_2b6c_a3b - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Mo2O7Rb2'
_chemical_formula_sum 'Mo2 O7 Rb2'

loop_
_publ_author_name
 'Z. A. Solodovnikova'
 'S. F. Solodovnikov'
_journal_name_full_name
;
 Acta Crystallographica Section C: Structural Chemistry
;
_journal_volume 62
_journal_year 2006
_journal_page_first i53
_journal_page_last i56
_publ_Section_title
;
 Rubidium dimolybdate, Rb$_{2}$Mo$_{2}$O$_{7}$, and caesium dimolybdate,
      ↪ Cs$_{2}$Mo$_{2}$O$_{7}$
;

_aflow_title 'Rb$_{2}$Mo$_{2}$O$_{7}$ Structure'
_aflow_proto 'A2B7C2_oC88_40_abc_2b6c_a3b'
_aflow_params 'a,b/a,c/a,z_{1},z_{2},y_{3},z_{3},y_{4},z_{4},y_{5},z_{5
      ↪ },y_{6},z_{6},y_{7},z_{7},y_{8},z_{8},x_{9},y_{9},z_{9},x_{10},
      ↪ y_{10},z_{10},x_{11},y_{11},z_{11},x_{12},y_{12},z_{12},x_{13},
      ↪ y_{13},z_{13},x_{14},y_{14},z_{14},x_{15},y_{15},z_{15}'
_aflow_params_values '11.8887,1.07920125834,0.861860422081,0.56117,0.0,
      ↪ 0.47105,0.28712,0.3413,0.244,0.4846,0.4559,0.48386,0.7361,
      ↪ 0.21144,0.4933,0.2554,-0.0226,0.01169,0.25276,0.22891,0.1103,
      ↪ 0.5171,-0.0426,0.0319,0.3571,0.1036,0.1236,0.5306,0.228,0.8759,
      ↪ 0.205,0.2201,0.1013,0.1507,0.1898,0.0325,0.2957,0.3879'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oC88'

_symmetry_space_group_name_H-M "A m a 2"
_symmetry_Int_Tables_number 40

_cell_length_a      11.88870
_cell_length_b      12.83030
_cell_length_c      10.24640
_cell_angle_alpha   90.00000
_cell_angle_beta    90.00000
_cell_angle_gamma   90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,z
3 x+1/2,-y,z
4 -x+1/2,y,z
5 x,y+1/2,z+1/2
6 -x,-y+1/2,z+1/2
7 x+1/2,-y+1/2,z+1/2
8 -x+1/2,y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Mo1 Mo 4 a 0.00000 0.00000 0.56117 1.00000
Rb1 Rb 4 a 0.00000 0.00000 0.00000 1.00000
Mo2 Mo 4 b 0.25000 0.47105 0.28712 1.00000
O1 O 4 b 0.25000 0.34130 0.24400 1.00000
O2 O 4 b 0.25000 0.48460 0.45590 1.00000
Rb2 Rb 4 b 0.25000 0.48386 0.73610 1.00000
Rb3 Rb 4 b 0.25000 0.21144 0.49330 1.00000
Rb4 Rb 4 b 0.25000 0.25540 -0.02260 1.00000
Mo3 Mo 8 c 0.01169 0.25276 0.22891 1.00000
O3 O 8 c 0.11030 0.51710 -0.04260 1.00000
O4 O 8 c 0.03190 0.35710 0.10360 1.00000
O5 O 8 c 0.12360 0.53060 0.22800 1.00000
O6 O 8 c 0.87590 0.20500 0.22010 1.00000
O7 O 8 c 0.10130 0.15070 0.18980 1.00000
O8 O 8 c 0.03250 0.29570 0.38790 1.00000
```

$Rb_2Mo_2O_7$: A2B7C2_oC88_40_abc_2b6c_a3b - POSCAR

```
A2B7C2_oC88_40_abc_2b6c_a3b & a,b/a,c/a,z1,z2,y3,z3,y4,z4,y5,z5,y6,z6,y7
      ↪ ,z7,y8,z8,x9,y9,z9,x10,y10,z10,x11,y11,z11,x12,y12,z12,x13,y13,
      ↪ z13,x14,y14,z14,x15,y15,z15 --params=11.8887,1.07920125834,
      ↪ 0.861860422081,0.56117,0.0,0.47105,0.28712,0.3413,0.244,0.4846,
      ↪ 0.4559,0.48386,0.7361,0.21144,0.4933,0.2554,-0.0226,0.01169,
      ↪ 0.25276,0.22891,0.1103,0.5171,-0.0426,0.0319,0.3571,0.1036,
      ↪ 0.1236,0.5306,0.228,0.8759,0.205,0.2201,0.1013,0.1507,0.1898,
      ↪ 0.0325,0.2957,0.3879 & Ama2 C_{2v}^{16} #40 (a^2b^6c^7) & oC88
      ↪ & None & Mo2O7Rb2 & Mo2O7Rb2 & Z. A. Solodovnikova and S. F.
      ↪ Solodovnikov, Acta Crystallogr. C 62, i53-i56 (2006)
   1.00000000000000
  11.88870000000000    0.00000000000000    0.00000000000000
   0.00000000000000    6.41515000000000   -5.12320000000000
   0.00000000000000    6.41515000000000    5.12320000000000
     Mo     O     Rb
      8    28     8
Direct
```

```
    0.00000000000000  −0.56117000000000   0.56117000000000   Mo   (4a)
    0.50000000000000  −0.56117000000000   0.56117000000000   Mo   (4a)
    0.25000000000000   0.18393000000000   0.75817000000000   Mo   (4b)
    0.75000000000000  −0.75817000000000  −0.18393000000000   Mo   (4b)
    0.01169000000000   0.02385000000000   0.48167000000000   Mo   (8c)
   −0.01169000000000  −0.48167000000000  −0.02385000000000   Mo   (8c)
    0.51169000000000  −0.48167000000000  −0.02385000000000   Mo   (8c)
    0.48831000000000   0.02385000000000   0.48167000000000   Mo   (8c)
    0.25000000000000   0.09730000000000   0.58530000000000   O    (4b)
    0.75000000000000  −0.58530000000000  −0.09730000000000   O    (4b)
    0.25000000000000   0.02870000000000   0.94050000000000   O    (4b)
    0.75000000000000  −0.94050000000000  −0.02870000000000   O    (4b)
    0.11030000000000   0.55970000000000   0.47450000000000   O    (8c)
   −0.11030000000000  −0.47450000000000  −0.55970000000000   O    (8c)
    0.61030000000000  −0.47450000000000  −0.55970000000000   O    (8c)
    0.38970000000000   0.55970000000000   0.47450000000000   O    (8c)
    0.03190000000000   0.25350000000000   0.46070000000000   O    (8c)
   −0.03190000000000  −0.46070000000000  −0.25350000000000   O    (8c)
    0.53190000000000  −0.46070000000000  −0.25350000000000   O    (8c)
    0.46810000000000   0.25350000000000   0.46070000000000   O    (8c)
    0.12360000000000   0.30260000000000   0.75860000000000   O    (8c)
   −0.12360000000000  −0.75860000000000  −0.30260000000000   O    (8c)
    0.62360000000000  −0.75860000000000  −0.30260000000000   O    (8c)
    0.37640000000000   0.30260000000000   0.75860000000000   O    (8c)
    0.87590000000000  −0.01510000000000   0.42510000000000   O    (8c)
   −0.87590000000000  −0.42510000000000   0.01510000000000   O    (8c)
    1.37590000000000  −0.42510000000000   0.01510000000000   O    (8c)
   −0.37590000000000  −0.01510000000000   0.42510000000000   O    (8c)
    0.10130000000000  −0.03910000000000   0.34050000000000   O    (8c)
   −0.10130000000000  −0.34050000000000   0.03910000000000   O    (8c)
    0.60130000000000  −0.34050000000000   0.03910000000000   O    (8c)
    0.39870000000000  −0.03910000000000   0.34050000000000   O    (8c)
    0.03250000000000  −0.09220000000000   0.68360000000000   O    (8c)
   −0.03250000000000  −0.68360000000000   0.09220000000000   O    (8c)
    0.53250000000000  −0.68360000000000   0.09220000000000   O    (8c)
    0.46750000000000  −0.09220000000000   0.68360000000000   O    (8c)
    0.00000000000000   0.00000000000000   0.00000000000000   Rb   (4a)
    0.50000000000000   0.00000000000000   0.00000000000000   Rb   (4a)
    0.25000000000000  −0.25224000000000   1.21996000000000   Rb   (4b)
    0.75000000000000  −1.21996000000000   0.25224000000000   Rb   (4b)
    0.25000000000000  −0.28186000000000   0.70474000000000   Rb   (4b)
    0.75000000000000  −0.70474000000000   0.28186000000000   Rb   (4b)
    0.25000000000000   0.27800000000000   0.23280000000000   Rb   (4b)
    0.75000000000000  −0.23280000000000  −0.27800000000000   Rb   (4b)
```

Orthorhombic CrO_3: AB3_oC16_40_b_a2b - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'CrO3'
_chemical_formula_sum 'Cr O3'

loop_
_publ_author_name
 'A. Bystr\"{o}m'
 'K.-A. Wilhelmi'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 4
_journal_year 1950
_journal_page_first 1131
_journal_page_last 1141
_publ_Section_title
;
 The Crystal Structure of Chromium Trioxide
;

_aflow_title 'Orthorhombic CrO$_{3}$ Structure'
_aflow_proto 'AB3_oC16_40_b_a2b'
_aflow_params 'a,b/a,c/a,z_{1},y_{2},z_{2},y_{3},z_{3},y_{4},z_{4}'
_aflow_params_values '5.743,1.54199895525,0.833884729236,0.444,0.097,
      ↪ 0.306,0.778,0.0,0.222,0.0'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oC16'

_symmetry_space_group_name_H-M "A m a 2"
_symmetry_Int_Tables_number 40

_cell_length_a    5.74300
_cell_length_b    8.85570
_cell_length_c    4.78900
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,z
3 x+1/2,-y,z
4 -x+1/2,y,z
5 x,y+1/2,z+1/2
6 -x,-y+1/2,z+1/2
7 x+1/2,-y+1/2,z+1/2
8 -x+1/2,y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 4 a 0.00000 0.00000 0.44400 1.00000
Cr1 Cr 4 b 0.25000 0.09700 0.30600 1.00000
O2 O 4 b 0.25000 0.77800 0.00000 1.00000
O3 O 4 b 0.25000 0.22200 0.00000 1.00000
```

Orthorhombic CrO_3: AB3_oC16_40_b_a2b - POSCAR

```
AB3_oC16_40_b_a2b & a,b/a,c/a,z1,y2,z2,y3,z3,y4,z4 --params=5.743,
      ↪ 1.54199895525,0.833884729236,0.444,0.097,0.306,0.778,0.0,0.222,
      ↪ 0.0 & Ama2 C_{2v}^{16} #40 (ab^3) & oC16 & None & CrO3 & CrO3 &
      ↪  A. Bystr\"{o}m and K.-A. Wilhelmi, Acta Chem. Scand. 4,
      ↪ 1131-1141 (1950)
   1.00000000000000
     5.74300000000000    0.00000000000000    0.00000000000000
     0.00000000000000    4.42785000000000   -2.39450000000000
     0.00000000000000    4.42785000000000    2.39450000000000
      Cr      O
       2      6
Direct
    0.25000000000000  -0.20900000000000   0.40300000000000   Cr   (4b)
    0.75000000000000  -0.40300000000000   0.20900000000000   Cr   (4b)
    0.00000000000000  -0.44400000000000   0.44400000000000   O    (4a)
    0.50000000000000  -0.44400000000000   0.44400000000000   O    (4a)
    0.25000000000000   0.77800000000000   0.77800000000000   O    (4b)
    0.75000000000000  -0.77800000000000  -0.77800000000000   O    (4b)
    0.25000000000000   0.22200000000000   0.22200000000000   O    (4b)
    0.75000000000000  -0.22200000000000  -0.22200000000000   O    (4b)
```

Santite ($KB_5O_8{\cdot}4H_2O$, $K3_5$): A5B8CD12_oC104_41_a2b_4b_a_6b - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Santite'
_chemical_formula_sum 'B5 H8 K O12'

loop_
_publ_author_name
 'W. H. Zachariasen'
 'H. A. Plettinger'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 16
_journal_year 1963
_journal_page_first 376
_journal_page_last 379
_publ_Section_title
;
 Refinement of the structure of potassium pentaborate tetrahydrate
;

_aflow_title 'Santite (KB$_{5}$O$_{8}$$\cdot$4H$_{2}$O, $K3_{5}$)
      ↪ Structure'
_aflow_proto 'A5B8CD12_oC104_41_a2b_4b_a_6b'
_aflow_params 'a,b/a,c/a,z_{1},z_{2},x_{3},y_{3},z_{3},x_{4},y_{4},z_{4
      ↪ },x_{5},y_{5},z_{5},x_{6},y_{6},z_{6},x_{7},y_{7},z_{7},x_{8},
      ↪ y_{8},z_{8},x_{9},y_{9},z_{9},x_{10},y_{10},z_{10},x_{11},y_{11
      ↪ },z_{11},x_{12},y_{12},z_{12},x_{13},y_{13},z_{13},x_{14},y_{14
      ↪ },z_{14}'
_aflow_params_values '11.062,1.01021605496,0.817302476948,0.4041,0.0,
      ↪ 0.189,0.0943,0.3126,0.2042,0.9432,0.4952,0.978,0.247,0.804,
      ↪ 0.967,0.111,0.649,0.161,0.159,0.741,0.251,0.316,0.193,0.0672,
      ↪ 0.0843,0.3082,0.0832,0.9313,0.4991,0.2591,0.0254,0.4033,0.2448,
      ↪ 0.1753,0.2235,0.2816,0.8801,0.5802,0.0151,0.1707,0.7588'
_aflow_Strukturbericht '$K3_{5}$'
_aflow_Pearson 'oC104'

_symmetry_space_group_name_H-M "A b a 2"
_symmetry_Int_Tables_number 41

_cell_length_a    11.06200
_cell_length_b    11.17501
_cell_length_c    9.04100
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,z
3 x+1/2,-y,z+1/2
4 -x+1/2,y,z+1/2
5 x,y+1/2,z+1/2
6 -x,-y+1/2,z+1/2
7 x+1/2,-y+1/2,z
8 -x+1/2,y+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
```

```
_atom_site_occupancy
B1 B 4 a 0.00000 0.00000 0.40410 1.00000
K1 K 4 a 0.00000 0.00000 0.00000 1.00000
B2 B 8 b 0.18900 0.09430 0.31260 1.00000
B3 B 8 b 0.20420 0.94320 0.49520 1.00000
H1 H 8 b 0.97800 0.24700 0.80400 1.00000
H2 H 8 b 0.96700 0.11100 0.64900 1.00000
H3 H 8 b 0.16100 0.15900 0.74100 1.00000
H4 H 8 b 0.25100 0.31600 0.19300 1.00000
O1 O 8 b 0.06720 0.08430 0.30820 1.00000
O2 O 8 b 0.08320 0.93130 0.49910 1.00000
O3 O 8 b 0.25910 0.02540 0.40330 1.00000
O4 O 8 b 0.24480 0.17530 0.22350 1.00000
O5 O 8 b 0.28160 0.88010 0.58020 1.00000
O6 O 8 b 0.01510 0.17070 0.75880 1.00000
```

Santite ($KB_5O_8{\cdot}4H_2O$, $K3_5$): A5B8CD12_oC104_41_a2b_4b_a_6b - POSCAR

```
A5B8CD12_oC104_41_a2b_4b_a_6b & a,b/a,c/a,z1,z2,x3,y3,z3,x4,y4,z4,x5,y5,
    ↪ z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10,x11,y11,z11,
    ↪ x12,y12,z12,x13,y13,z13,x14,y14,z14 --params=11.062,
    ↪ 1.01021605496,0.817302476948,0.4041,0.0,0.189,0.0943,0.3126,
    ↪ 0.2042,0.9432,0.4952,0.978,0.247,0.804,0.967,0.111,0.649,0.161,
    ↪ 0.159,0.741,0.251,0.316,0.193,0.0672,0.0843,0.3082,0.0832,
    ↪ 0.9313,0.4991,0.2591,0.0254,0.4033,0.2448,0.1753,0.2235,0.2816,
    ↪ 0.8801,0.5802,0.0151,0.1707,0.7588 & Aba2 C_{2v}^{17} #41 (a^2b
    ↪ ^12) & oC104 & $K3_{5}$ & B5H8KO12 & Santite & W. H.
    ↪ Zachariasen and H. A. Plettinger, Acta Cryst. 16, 376-379 (1963
    ↪ )
   1.00000000000000
  11.06200000000000    0.00000000000000    0.00000000000000
   0.00000000000000    5.58750500000000   -4.52050000000000
   0.00000000000000    5.58750500000000    4.52050000000000
     B     H     K     O
    10    16     2    24
Direct
   0.00000000000000   -0.40410000000000    0.40410000000000   B   (4a)
   0.50000000000000    0.09590000000000    0.90410000000000   B   (4a)
   0.18900000000000   -0.21830000000000    0.40690000000000   B   (8b)
  -0.18900000000000   -0.40690000000000    0.21830000000000   B   (8b)
   0.68900000000000    0.09310000000000    0.71830000000000   B   (8b)
   0.31100000000000    0.28170000000000    0.90690000000000   B   (8b)
   0.20420000000000    0.44800000000000    1.43840000000000   B   (8b)
  -0.20420000000000   -1.43840000000000   -0.44800000000000   B   (8b)
   0.70420000000000   -0.93840000000000    0.05200000000000   B   (8b)
   0.29580000000000    0.94800000000000    1.93840000000000   B   (8b)
   0.97800000000000   -0.55700000000000    1.05100000000000   H   (8b)
  -0.97800000000000   -1.05100000000000    0.55700000000000   H   (8b)
   1.47800000000000   -0.55100000000000    1.05700000000000   H   (8b)
  -0.47800000000000   -0.05700000000000    1.55100000000000   H   (8b)
   0.96700000000000   -0.53800000000000    0.76000000000000   H   (8b)
  -0.96700000000000   -0.76000000000000    0.53800000000000   H   (8b)
   1.46700000000000   -0.26000000000000    1.03800000000000   H   (8b)
  -0.46700000000000   -0.03800000000000    1.26000000000000   H   (8b)
   0.16100000000000   -0.58200000000000    0.90000000000000   H   (8b)
  -0.16100000000000   -0.90000000000000    0.58200000000000   H   (8b)
   0.66100000000000   -0.40000000000000    1.08200000000000   H   (8b)
   0.33900000000000   -0.08200000000000    1.40000000000000   H   (8b)
   0.25100000000000    0.12300000000000    0.50900000000000   H   (8b)
  -0.25100000000000   -0.50900000000000   -0.12300000000000   H   (8b)
   0.75100000000000   -0.00900000000000    0.37700000000000   H   (8b)
   0.24900000000000    0.62300000000000    1.00900000000000   H   (8b)
   0.00000000000000    0.00000000000000    0.00000000000000   K   (4a)
   0.50000000000000    0.50000000000000    0.50000000000000   K   (4a)
   0.06720000000000   -0.22390000000000    0.39250000000000   O   (8b)
  -0.06720000000000   -0.39250000000000    0.22390000000000   O   (8b)
   0.56720000000000    0.10750000000000    0.72390000000000   O   (8b)
   0.43280000000000    0.27610000000000    0.89250000000000   O   (8b)
   0.08320000000000    0.43220000000000    1.43040000000000   O   (8b)
  -0.08320000000000   -1.43040000000000   -0.43220000000000   O   (8b)
   0.58320000000000   -0.93040000000000    0.06780000000000   O   (8b)
   0.41680000000000    0.93220000000000    1.93040000000000   O   (8b)
   0.25910000000000   -0.37790000000000    0.42870000000000   O   (8b)
  -0.25910000000000   -0.42870000000000    0.37790000000000   O   (8b)
   0.75910000000000    0.07130000000000    0.87790000000000   O   (8b)
   0.24090000000000    0.12210000000000    0.92870000000000   O   (8b)
   0.24480000000000   -0.04820000000000    0.39880000000000   O   (8b)
  -0.24480000000000   -0.39880000000000    0.04820000000000   O   (8b)
   0.74480000000000    0.10120000000000    0.54820000000000   O   (8b)
   0.25520000000000    0.45180000000000    0.89880000000000   O   (8b)
   0.28160000000000    0.29990000000000    1.46030000000000   O   (8b)
  -0.28160000000000   -1.46030000000000   -0.29990000000000   O   (8b)
   0.78160000000000   -0.96030000000000    0.20010000000000   O   (8b)
   0.21840000000000    0.79990000000000    1.96030000000000   O   (8b)
   0.01510000000000   -0.58810000000000    0.92950000000000   O   (8b)
  -0.01510000000000   -0.92950000000000    0.58810000000000   O   (8b)
   0.51510000000000   -0.42950000000000    1.08810000000000   O   (8b)
   0.48490000000000   -0.08810000000000    1.42950000000000   O   (8b)
```

Ag_2O_3: A2B3_oF40_43_b_ab - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Ag2O3'
_chemical_formula_sum 'Ag2 O3'

loop_
_publ_author_name
 'B. Standke'
 'M. Jansen'
_journal_name_full_name
;
 Zeitschrift fur Anorganische und Allgemeine Chemie
;
_journal_volume 535
_journal_year 1986
_journal_page_first 39
_journal_page_last 46
_publ_Section_title
;
 Darstellung und Kristallstruktur von Ag$_{2}$O$_{3}$
;

_aflow_title 'Ag$_{2}$O$_{3}$ Structure'
_aflow_proto 'A2B3_oF40_43_b_ab'
_aflow_params 'a,b/a,c/a,z_{1},x_{2},y_{2},z_{2},x_{3},y_{3},z_{3}'
_aflow_params_values '12.869,0.815137151294,0.284699665864,0.8402,-
    ↪ 0.0467,0.1325,0.5,0.083,0.2333,0.6153'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oF40'

_symmetry_space_group_name_H-M "F d d 2"
_symmetry_Int_Tables_number 43

_cell_length_a      12.86900
_cell_length_b      10.49000
_cell_length_c      3.66380
_cell_angle_alpha   90.00000
_cell_angle_beta    90.00000
_cell_angle_gamma   90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,z
3 -x+1/4,y+1/4,z+1/4
4 x+1/4,-y+1/4,z+1/4
5 x,y+1/2,z+1/2
6 -x,-y+1/2,z+1/2
7 -x+1/4,y+3/4,z+3/4
8 x+1/4,-y+3/4,z+3/4
9 x+1/2,y,z+1/2
10 -x+1/2,-y,z+1/2
11 -x+3/4,y+1/4,z+3/4
12 x+3/4,-y+1/4,z+3/4
13 x+1/2,y+1/2,z
14 -x+1/2,-y+1/2,z
15 -x+3/4,y+3/4,z+1/4
16 x+3/4,-y+3/4,z+1/4

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 8 a 0.00000 0.00000 0.84020 1.00000
Ag1 Ag 16 b -0.04670 0.13250 0.50000 1.00000
O2 O 16 b 0.08300 0.23330 0.61530 1.00000
```

Ag_2O_3: A2B3_oF40_43_b_ab - POSCAR

```
A2B3_oF40_43_b_ab & a,b/a,c/a,z1,x2,y2,z2,x3,y3,z3 --params=12.869,
    ↪ 0.815137151294,0.284699665864,0.8402,-0.0467,0.1325,0.5,0.083,
    ↪ 0.2333,0.6153 & Fdd2 C_{2v}^{19} #43 (ab^2) & oF40 & None &
    ↪ Ag2O3 & Ag2O3 & B. Standke and M. Jansen, Z. Anorg. Allg. Chem.
    ↪ 535, 39-46 (1986)
   1.00000000000000
   0.00000000000000    5.24500000000000    1.83190000000000
   6.43450000000000    0.00000000000000    1.83190000000000
   6.43450000000000    5.24500000000000    0.00000000000000
    Ag     O
     4     6
Direct
   0.67920000000000    0.32080000000000   -0.41420000000000   Ag   (16b)
   0.32080000000000    0.67920000000000   -0.58580000000000   Ag   (16b)
   0.66420000000000    0.83580000000000   -0.42920000000000   Ag   (16b)
   0.83580000000000    0.66420000000000   -0.07080000000000   Ag   (16b)
   0.84020000000000    0.84020000000000   -0.84020000000000   O    (8a)
   1.09020000000000    1.09020000000000   -0.59020000000000   O    (8a)
   0.76560000000000    0.46500000000000   -0.29900000000000   O   (16b)
   0.46500000000000    0.76560000000000   -0.93160000000000   O   (16b)
   0.54900000000000    1.18160000000000   -0.51560000000000   O   (16b)
   1.18160000000000    0.54900000000000   -0.21500000000000   O   (16b)
```

Natrolite ($Na_2Al_2Si_3O_{10}{\cdot}2H_2O$, $S6_{10}$): A2B4C2D12E3_oF184_43_b_2b_b_6b_ab - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Natrolite'
_chemical_formula_sum 'Al2 H4 Na2 O12 Si3'

loop_
_publ_author_name
 'A. Kirfel'
 'M. Orthen'
 'G. Will'
_journal_name_full_name
;
 Zeolites
;
_journal_volume 4
```

```
_journal_year 1984
_journal_page_first 140
_journal_page_last 146
_publ_Section_title
;
 Natrolite: refinement of the crystal structure of two samples from
      ↪ Marienberg (Usti nad Labem, CSSR)
;

_aflow_title 'Natrolite (Na$_{2}$Al$_{2}$Si$_{3}$O$_{10}$$\cdot$2H$_{2}
      ↪ $O, $S6_{10}$) Structure'
_aflow_proto 'A2B4C2D12E3_oF184_43_b_2b_b_6b_ab'
_aflow_params 'a,b/a,c/a,z_{1},x_{2},y_{2},z_{2},x_{3},y_{3},z_{3},x_{4
      ↪ },y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{6},x_{7},y_{7},
      ↪ z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{10},y_{10},z_{10},
      ↪ x_{11},y_{11},z_{11},x_{12},y_{12},z_{12}'
_aflow_params_values '18.296,1.01918452121,0.359914735461,0.0,0.0374,
      ↪ 0.0937,0.6152,0.0533,0.1604,0.0616,0.0848,0.1932,0.1731,0.2208,
      ↪ 0.0307,0.6179,0.0226,0.0686,0.866,0.0701,0.1818,0.6099,0.0983,
      ↪ 0.0351,0.5003,0.2065,0.1529,0.7261,0.1804,0.2273,0.3903,0.0561,
      ↪ 0.1896,0.1112,0.1533,0.2113,0.6231'
_aflow_Strukturbericht '$S6_{10}$'
_aflow_Pearson 'oF184'

_symmetry_space_group_name_H-M "F d d 2"
_symmetry_Int_Tables_number 43

_cell_length_a      18.29600
_cell_length_b      18.64700
_cell_length_c      6.58500
_cell_angle_alpha   90.00000
_cell_angle_beta    90.00000
_cell_angle_gamma   90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,z
3 -x+1/4,y+1/4,z+1/4
4 x+1/4,-y+1/4,z+1/4
5 x,y+1/2,z+1/2
6 -x,-y+1/2,z+1/2
7 -x+1/4,y+3/4,z+3/4
8 x+1/4,-y+3/4,z+3/4
9 x+1/2,y,z+1/2
10 -x+1/2,-y,z+1/2
11 -x+3/4,y+1/4,z+3/4
12 x+3/4,-y+1/4,z+3/4
13 x+1/2,y+1/2,z
14 -x+1/2,-y+1/2,z
15 -x+3/4,y+3/4,z+1/4
16 x+3/4,-y+3/4,z+1/4

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Si1 Si 8 a 0.00000 0.00000 0.00000 1.00000
Al1 Al 16 b 0.03740 0.09370 0.61520 1.00000
H1 H 16 b 0.05330 0.16040 0.06160 1.00000
H2 H 16 b 0.08480 0.19320 0.17310 1.00000
Na1 Na 16 b 0.22080 0.03070 0.61790 1.00000
O1 O 16 b 0.02260 0.06860 0.86600 1.00000
O2 O 16 b 0.07010 0.18180 0.60990 1.00000
O3 O 16 b 0.09830 0.03510 0.50030 1.00000
O4 O 16 b 0.20650 0.15290 0.72610 1.00000
O5 O 16 b 0.18040 0.22730 0.39030 1.00000
O6 O 16 b 0.05610 0.18960 0.11120 1.00000
Si2 Si 16 b 0.15330 0.21130 0.62310 1.00000
```

Natrolite ($Na_2Al_2Si_3O_{10}{\cdot}2H_2O$, $S6_{10}$): A2B4C2D12E3_oF184_43_b_2b_b_6b_ab - POSCAR

```
A2B4C2D12E3_oF184_43_b_2b_b_6b_ab & a,b/a,c/a,z1,x2,y2,z2,x3,y3,z3,x4,y4
      ↪ ,z4,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10,
      ↪ x11,y11,z11,x12,y12,z12 --params=18.296,1.01918452121,
      ↪ 0.359914735461,0.0,0.0374,0.0937,0.6152,0.0533,0.1604,0.0616,
      ↪ 0.0848,0.1932,0.1731,0.2208,0.0307,0.6179,0.0226,0.0686,0.866,
      ↪ 0.0701,0.1818,0.6099,0.0983,0.0351,0.5003,0.2065,0.1529,0.7261,
      ↪ 0.1804,0.2273,0.3903,0.0561,0.1896,0.1112,0.1533,0.2113,0.6231
      ↪ & Fdd2 C_{2v}^{19} #43 (ab^11) & oF184 & $S6_{10}$ &
      ↪ Al2H4Na2O12Si3 & Natrolite & A. Kirfel and M. Orthen and G.
      ↪ Will, Zeolites 4, 140-146 (1984)
   1.00000000000000
     0.00000000000000     9.32350000000000     3.29250000000000
     9.14800000000000     0.00000000000000     3.29250000000000
     9.14800000000000     9.32350000000000     0.00000000000000
      Al     H     Na     O     Si
       4     8      4    24      6
Direct
     0.67150000000000     0.55890000000000    -0.48410000000000    Al   (16b)
     0.55890000000000     0.67150000000000    -0.74630000000000    Al   (16b)
     0.73410000000000     0.99630000000000    -0.42150000000000    Al   (16b)
     0.99630000000000     0.73410000000000    -0.30890000000000    Al   (16b)
     0.16870000000000    -0.04550000000000     0.15210000000000    H    (16b)
    -0.04550000000000     0.16870000000000    -0.27530000000000    H    (16b)
     0.09790000000000     0.52530000000000     0.08130000000000    H    (16b)
     0.52530000000000     0.09790000000000     0.29550000000000    H    (16b)
     0.28150000000000     0.06470000000000     0.10490000000000    H    (16b)
     0.06470000000000     0.28150000000000    -0.45110000000000    H    (16b)
     0.14510000000000     0.70110000000000    -0.03150000000000    H    (16b)
     0.70110000000000     0.14510000000000     0.18530000000000    H    (16b)
     0.42780000000000     0.80800000000000    -0.36640000000000    Na   (16b)
     0.80800000000000     0.42780000000000    -0.86940000000000    Na   (16b)
     0.61640000000000     1.11940000000000    -0.17780000000000    Na   (16b)
     1.11940000000000     0.61640000000000    -0.55800000000000    Na   (16b)
     0.91200000000000     0.82000000000000    -0.77480000000000    O    (16b)
     0.82000000000000     0.91200000000000    -0.95720000000000    O    (16b)
     1.02480000000000     1.20720000000000    -0.66200000000000    O    (16b)
     1.20720000000000     1.02480000000000    -0.57000000000000    O    (16b)
     0.72160000000000     0.49820000000000    -0.35800000000000    O    (16b)
     0.49820000000000     0.72160000000000    -0.86180000000000    O    (16b)
     0.60800000000000     1.11180000000000    -0.47160000000000    O    (16b)
     1.11180000000000     0.60800000000000    -0.24820000000000    O    (16b)
     0.43710000000000     0.56350000000000    -0.36690000000000    O    (16b)
     0.56350000000000     0.43710000000000    -0.63370000000000    O    (16b)
     0.61690000000000     0.88370000000000    -0.18710000000000    O    (16b)
     0.88370000000000     0.61690000000000    -0.31350000000000    O    (16b)
     0.67250000000000     0.77970000000000    -0.36670000000000    O    (16b)
     0.77970000000000     0.67250000000000    -1.08550000000000    O    (16b)
     0.61670000000000     1.33550000000000    -0.42250000000000    O    (16b)
     1.33550000000000     0.61670000000000    -0.52970000000000    O    (16b)
     0.43720000000000     0.34340000000000     0.01740000000000    O    (16b)
     0.34340000000000     0.43720000000000    -0.79800000000000    O    (16b)
     0.23260000000000     1.04800000000000    -0.18720000000000    O    (16b)
     1.04800000000000     0.23260000000000    -0.09340000000000    O    (16b)
     0.24470000000000    -0.02230000000000     0.13450000000000    O    (16b)
    -0.02230000000000     0.24470000000000    -0.35690000000000    O    (16b)
     0.11550000000000     0.60690000000000     0.00530000000000    O    (16b)
     0.60690000000000     0.11550000000000     0.27230000000000    O    (16b)
     0.00000000000000     0.00000000000000     0.00000000000000    Si   (8a)
     0.25000000000000     0.25000000000000     0.25000000000000    Si   (8a)
     0.68110000000000     0.56510000000000    -0.25850000000000    Si   (16b)
     0.56510000000000     0.68110000000000    -0.98770000000000    Si   (16b)
     0.50850000000000     1.23770000000000    -0.43110000000000    Si   (16b)
     1.23770000000000     0.50850000000000    -0.31510000000000    Si   (16b)
```

Blossite (α-$Cu_2V_2O_7$): A2B7C2_oF88_43_b_a3b_b - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Blossite'
_chemical_formula_sum 'Cu2 O7 V2'

loop_
_publ_author_name
 'C. Calvo'
 'R. Faggiani'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 31
_journal_year 1975
_journal_page_first 603
_journal_page_last 605
_publ_Section_title
;
 $\alpha$ Cupric Divanadate
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'Blossite ($\alpha$-Cu$_{2}$V$_{2}$O$_{7}$) Structure'
_aflow_proto 'A2B7C2_oF88_43_b_a3b_b'
_aflow_params 'a,b/a,c/a,z_{1},x_{2},y_{2},z_{2},x_{3},y_{3},z_{3},x_{4
      ↪ },y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{6}'
_aflow_params_values '20.645,0.406054734803,0.312036812788,-0.0908,
      ↪ 0.1658,0.3646,0.75,0.2453,0.5622,0.2774,0.1446,0.4368,0.0332,
      ↪ 0.1619,0.3477,0.4592,0.19906,0.4054,0.2343'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oF88'

_symmetry_space_group_name_H-M "F d d 2"
_symmetry_Int_Tables_number 43

_cell_length_a      20.64500
_cell_length_b      8.38300
_cell_length_c      6.44200
_cell_angle_alpha   90.00000
_cell_angle_beta    90.00000
_cell_angle_gamma   90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,z
3 -x+1/4,y+1/4,z+1/4
4 x+1/4,-y+1/4,z+1/4
5 x,y+1/2,z+1/2
6 -x,-y+1/2,z+1/2
7 -x+1/4,y+3/4,z+3/4
8 x+1/4,-y+3/4,z+3/4
9 x+1/2,y,z+1/2
10 -x+1/2,-y,z+1/2
11 -x+3/4,y+1/4,z+3/4
12 x+3/4,-y+1/4,z+3/4
13 x+1/2,y+1/2,z
14 -x+1/2,-y+1/2,z
15 -x+3/4,y+3/4,z+1/4
16 x+3/4,-y+3/4,z+1/4
```

```
loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 8 a 0.00000 0.00000 -0.09080 1.00000
Cu1 Cu 16 b 0.16580 0.36460 0.75000 1.00000
O2 O 16 b 0.24530 0.56220 0.27740 1.00000
O3 O 16 b 0.14460 0.43680 0.03320 1.00000
O4 O 16 b 0.16190 0.34770 0.45920 1.00000
V1 V 16 b 0.19906 0.40540 0.23430 1.00000
```

Blossite (α-$Cu_2V_2O_7$): A2B7C2_oF88_43_b_a3b_b - POSCAR

```
A2B7C2_oF88_43_b_a3b_b & a,b/a,c/a,z1,x2,y2,z2,x3,y3,z3,x4,y4,z4,x5,y5,
    ↪ z5,x6,y6,z6 --params=20.645,0.406054734803,0.312036812788,-
    ↪ 0.0908,0.1658,0.3646,0.75,0.2453,0.5622,0.2774,0.1446,0.4368,
    ↪ 0.0332,0.1619,0.3477,0.4592,0.19906,0.4054,0.2343 & Fdd2 C_{2v
    ↪ }^{19} #43 (ab^5) & oF88 & None & Cu2O7V2 & Blossite & C. Calvo
    ↪  and R. Faggiani, Acta Crystallogr. Sect. B Struct. Sci. 31,
    ↪ 603-605 (1975)
   1.00000000000000
   0.00000000000000    4.19150000000000    3.22100000000000
  10.32250000000000    0.00000000000000    3.22100000000000
  10.32250000000000    4.19150000000000    0.00000000000000
     Cu      O      V
      4     14      4
Direct
   0.94880000000000    0.55120000000000   -0.21960000000000    Cu   (16b)
   0.55120000000000    0.94880000000000   -1.28040000000000    Cu   (16b)
   0.46960000000000    1.53040000000000   -0.69880000000000    Cu   (16b)
   1.53040000000000    0.46960000000000   -0.30120000000000    Cu   (16b)
  -0.09080000000000   -0.09080000000000    0.09080000000000    O    (8a)
   0.15920000000000    0.15920000000000    0.34080000000000    O    (8a)
   0.59430000000000   -0.03950000000000    0.53010000000000    O   (16b)
  -0.03950000000000    0.59430000000000   -1.08490000000000    O   (16b)
  -0.28010000000000    1.33490000000000   -0.34430000000000    O   (16b)
   1.33490000000000   -0.28010000000000    0.28950000000000    O   (16b)
   0.32540000000000   -0.25900000000000    0.54820000000000    O   (16b)
  -0.25900000000000    0.32540000000000   -0.61460000000000    O   (16b)
  -0.29820000000000    0.86460000000000   -0.07540000000000    O   (16b)
   0.86460000000000   -0.29820000000000    0.50900000000000    O   (16b)
   0.64500000000000    0.27340000000000    0.05040000000000    O   (16b)
   0.27340000000000    0.64500000000000   -0.96880000000000    O   (16b)
   0.19960000000000    1.21880000000000   -0.39500000000000    O   (16b)
   1.21880000000000    0.19960000000000   -0.02340000000000    O   (16b)
   0.44064000000000    0.02796000000000    0.37016000000000    V   (16b)
   0.02796000000000    0.44064000000000   -0.83876000000000    V   (16b)
  -0.12016000000000    1.08876000000000   -0.19064000000000    V   (16b)
   1.08876000000000   -0.12016000000000    0.22204000000000    V   (16b)
```

Archerite (KH_2PO_4): A2BC4D_oF64_43_b_a_2b_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Archerite'
_chemical_formula_sum 'H2 K O4 P'

loop_
_publ_author_name
 'H. A. Levy'
 'S. W. Peterson'
 'S. H. Simonsen'
_journal_name_full_name
;
 Physical Review
;
_journal_volume 93
_journal_year 1954
_journal_page_first 1120
_journal_page_last 1121
_publ_Section_title
;
 Neutron Diffraction Study of the Ferroelectric Modification of
    ↪ Potassium Dihydrogen Phosphate
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'Archerite (KH$_{2}$PO$_{4}$) Structure'
_aflow_proto 'A2BC4D_oF64_43_b_a_2b_a'
_aflow_params 'a,b/a,c/a,z_{1},z_{2},x_{3},y_{3},z_{3},x_{4},y_{4},z_{4
    ↪ },x_{5},y_{5},z_{5}'
_aflow_params_values '10.53,0.991452991453,0.655270655271,0.512,0.0,
    ↪ 0.188,-0.0375,0.1355,0.116,0.0345,0.131,-0.0345,0.116,0.8765'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oF64'

_symmetry_space_group_name_H-M "F d d 2"
_symmetry_Int_Tables_number 43

_cell_length_a     10.53000
_cell_length_b     10.44000
_cell_length_c     6.90000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,z
3 -x+1/4,y+1/4,z+1/4
4 x+1/4,-y+1/4,z+1/4
5 x,y+1/2,z+1/2
6 -x,-y+1/2,z+1/2
7 -x+1/4,y+3/4,z+3/4
8 x+1/4,-y+3/4,z+3/4
9 x+1/2,y,z+1/2
10 -x+1/2,-y,z+1/2
11 -x+3/4,y+1/4,z+3/4
12 x+3/4,-y+1/4,z+3/4
13 x+1/2,y+1/2,z
14 -x+1/2,-y+1/2,z
15 -x+3/4,y+3/4,z+1/4
16 x+3/4,-y+3/4,z+1/4

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
K1 K 8 a 0.00000 0.00000 0.51200 1.00000
P1 P 8 a 0.00000 0.00000 0.00000 1.00000
H1 H 16 b 0.18800 -0.03750 0.13550 1.00000
O1 O 16 b 0.11600 0.03450 0.13100 1.00000
O2 O 16 b -0.03450 0.11600 0.87650 1.00000
```

Archerite (KH_2PO_4): A2BC4D_oF64_43_b_a_2b_a - POSCAR

```
A2BC4D_oF64_43_b_a_2b_a & a,b/a,c/a,z1,z2,x3,y3,z3,x4,y4,z4,x5,y5,z5 --
    ↪ params=10.53,0.991452991453,0.655270655271,0.512,0.0,0.188,-
    ↪ 0.0375,0.1355,0.116,0.0345,0.131,-0.0345,0.116,0.8765 & Fdd2 C_
    ↪ {2v}^{19} #43 (a^2b^3) & oF64 & None & H2KO4P & Archerite & H.
    ↪ A. Levy and S. W. Peterson and S. H. Simonsen, Phys. Rev. 93,
    ↪ 1120-1121 (1954)
   1.00000000000000
   0.00000000000000    5.22000000000000    3.45000000000000
   5.26500000000000    0.00000000000000    3.45000000000000
   5.26500000000000    5.22000000000000    0.00000000000000
     H      K      O      P
     4      2      8      2
Direct
  -0.09000000000000    0.36100000000000    0.01500000000000    H   (16b)
   0.36100000000000   -0.09000000000000   -0.28600000000000    H   (16b)
   0.23500000000000    0.53600000000000    0.34000000000000    H   (16b)
   0.53600000000000    0.23500000000000   -0.11100000000000    H   (16b)
   0.51200000000000    0.51200000000000   -0.51200000000000    K    (8a)
   0.76200000000000    0.76200000000000   -0.26200000000000    K    (8a)
   0.04950000000000    0.21250000000000    0.01950000000000    O   (16b)
   0.21250000000000    0.04950000000000   -0.28150000000000    O   (16b)
   0.23050000000000    0.53150000000000    0.20050000000000    O   (16b)
   0.53150000000000    0.23050000000000    0.03750000000000    O   (16b)
   1.02700000000000    0.72600000000000   -0.79500000000000    O   (16b)
   0.72600000000000    1.02700000000000   -0.95800000000000    O   (16b)
   1.04500000000000    1.20800000000000   -0.77700000000000    O   (16b)
   1.20800000000000    1.04500000000000   -0.47600000000000    O   (16b)
   0.00000000000000    0.00000000000000    0.00000000000000    P    (8a)
   0.25000000000000    0.25000000000000    0.25000000000000    P    (8a)
```

Cs_2Se: A2B_oF24_43_b_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Cs2Se'
_chemical_formula_sum 'Cs2 Se'

loop_
_publ_author_name
 'P. B\"{o}ttcher'
_journal_name_full_name
;
 Journal of the Less-Common Metals
;
_journal_volume 76
_journal_year 1980
_journal_page_first 271
_journal_page_last 277
_publ_Section_title
;
 Zur Kenntnis von Cs$_{2}$Se
;

# Found in Pearson's Handbook of Crystallographic Data, 1991

_aflow_title 'Cs$_{2}$Se Structure'
_aflow_proto 'A2B_oF24_43_b_a'
_aflow_params 'a,b/a,c/a,z_{1},x_{2},y_{2},z_{2}'
_aflow_params_values '16.49,0.712553062462,0.410855063675,0.0,0.0749,
    ↪ 0.202,0.8118'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oF24'

_symmetry_space_group_name_H-M "F d d 2"
_symmetry_Int_Tables_number 43

_cell_length_a     16.49000
```

```
_cell_length_b      11.75000
_cell_length_c      6.77500
_cell_angle_alpha   90.00000
_cell_angle_beta    90.00000
_cell_angle_gamma   90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,z
3 -x+1/4,y+1/4,z+1/4
4 x+1/4,-y+1/4,z+1/4
5 x,y+1/2,z+1/2
6 -x,-y+1/2,z+1/2
7 -x+1/4,y+3/4,z+3/4
8 x+1/4,-y+3/4,z+3/4
9 x+1/2,y,z+1/2
10 -x+1/2,-y,z+1/2
11 -x+3/4,y+1/4,z+3/4
12 x+3/4,-y+1/4,z+3/4
13 x+1/2,y+1/2,z
14 -x+1/2,-y+1/2,z
15 -x+3/4,y+3/4,z+1/4
16 x+3/4,-y+3/4,z+1/4

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Se1 Se 8 a 0.00000 0.00000 0.00000 1.00000
Cs1 Cs 16 b 0.07490 0.20200 0.81180 1.00000
```

Cs_2Se: A2B_oF24_43_b_a - POSCAR

```
A2B_oF24_43_b_a & a,b/a,c/a,z1,x2,y2,z2 --params=16.49,0.712553062462,
      ↪ 0.410855063675,0.0,0.0749,0.202,0.8118 & Fdd2 C_{2v}^{19} #43 (
      ↪ ab) & oF24 & None & Cs2Se & Cs2Se & P. B\"{o}ttcher, J.
      ↪ Less-Common Met. 76, 271-277 (1980)
   1.00000000000000
   0.00000000000000    5.87500000000000    3.38750000000000
   8.24500000000000    0.00000000000000    3.38750000000000
   8.24500000000000    5.87500000000000    0.00000000000000
    Cs     Se
     4      2
Direct
   0.93890000000000    0.68470000000000   -0.53490000000000    Cs    (16b)
   0.68470000000000    0.93890000000000   -1.08870000000000    Cs    (16b)
   0.78490000000000    1.33870000000000   -0.68890000000000    Cs    (16b)
   1.33870000000000    0.78490000000000   -0.43470000000000    Cs    (16b)
   0.00000000000000    0.00000000000000    0.00000000000000    Se     (8a)
   0.25000000000000    0.25000000000000    0.25000000000000    Se     (8a)
```

Zr_2Al_3: A3B2_oF40_43_ab_b - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Al3Zr2'
_chemical_formula_sum 'Al3 Zr2'

loop_
_publ_author_name
 'T. J. Renouf'
 'C. A. Beevers'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 14
_journal_year 1961
_journal_page_first 469
_journal_page_last 472
_publ_Section_title
;
 The Crystal Structure of Zr${2}$Al${3}$
;

# Found in Crystal Structure Investigations on the Zr-Al and Hf-Al
      ↪ Systems, 1962

_aflow_title 'Zr${2}$Al${3}$ Structure'
_aflow_proto 'A3B2_oF40_43_ab_b'
_aflow_params 'a,b/a,c/a,z_{1},x_{2},y_{2},z_{2},x_{3},y_{3},z_{3}'
_aflow_params_values '9.601,1.44839079263,0.58014790126,0.125,0.065,
      ↪ 0.116,0.5,0.068,0.196,0.0'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oF40'

_symmetry_space_group_name_H-M "F d d 2"
_symmetry_Int_Tables_number 43

_cell_length_a      9.60100
_cell_length_b      13.90600
_cell_length_c      5.57000
_cell_angle_alpha   90.00000
_cell_angle_beta    90.00000
_cell_angle_gamma   90.00000
```

```
loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,z
3 -x+1/4,y+1/4,z+1/4
4 x+1/4,-y+1/4,z+1/4
5 x,y+1/2,z+1/2
6 -x,-y+1/2,z+1/2
7 -x+1/4,y+3/4,z+3/4
8 x+1/4,-y+3/4,z+3/4
9 x+1/2,y,z+1/2
10 -x+1/2,-y,z+1/2
11 -x+3/4,y+1/4,z+3/4
12 x+3/4,-y+1/4,z+3/4
13 x+1/2,y+1/2,z
14 -x+1/2,-y+1/2,z
15 -x+3/4,y+3/4,z+1/4
16 x+3/4,-y+3/4,z+1/4

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Al1 Al 8 a 0.00000 0.00000 0.12500 1.00000
Al2 Al 16 b 0.06500 0.11600 0.50000 1.00000
Zr1 Zr 16 b 0.06800 0.19600 0.00000 1.00000
```

Zr_2Al_3: A3B2_oF40_43_ab_b - POSCAR

```
A3B2_oF40_43_ab_b & a,b/a,c/a,z1,x2,y2,z2,x3,y3,z3 --params=9.601,
      ↪ 1.44839079263,0.58014790126,0.125,0.065,0.116,0.5,0.068,0.196,
      ↪ 0.0 & Fdd2 C_{2v}^{19} #43 (ab^2) & oF40 & None & Al3Zr2 &
      ↪ Al3Zr2 & T. J. Renouf and C. A. Beevers, Acta Cryst. 14,
      ↪ 469-472 (1961)
   1.00000000000000
   0.00000000000000    6.95300000000000    2.78500000000000
   4.80050000000000    0.00000000000000    2.78500000000000
   4.80050000000000    6.95300000000000    0.00000000000000
    Al     Zr
     6      4
Direct
   0.12500000000000    0.12500000000000   -0.12500000000000    Al    (8a)
   0.37500000000000    0.37500000000000    0.12500000000000    Al    (8a)
   0.55100000000000    0.44900000000000   -0.31900000000000    Al   (16b)
   0.44900000000000    0.55100000000000   -0.68100000000000    Al   (16b)
   0.56900000000000    0.93100000000000   -0.30100000000000    Al   (16b)
   0.93100000000000    0.56900000000000   -0.19900000000000    Al   (16b)
   0.12800000000000   -0.12800000000000    0.26400000000000    Zr   (16b)
  -0.12800000000000    0.12800000000000   -0.26400000000000    Zr   (16b)
  -0.01400000000000    0.51400000000000    0.12200000000000    Zr   (16b)
   0.51400000000000   -0.01400000000000    0.37800000000000    Zr   (16b)
```

Hemimorphite ($Zn_4Si_2O_7(OH)_2{\cdot}H_2O$, $S2_2$): A2B5CD2_oI40_44_2c_abcde_d_e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Hemimorphite'
_chemical_formula_sum 'H2 O5 Si Zn2'

loop_
_publ_author_name
 'R. J. Hill'
 'G. V. Gibbs'
 'J. R. Craig'
 'F. K. Ross'
 'J. M. Williams'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 146
_journal_year 1977
_journal_page_first 241
_journal_page_last 259
_publ_Section_title
;
 A neutron-diffraction study of hemimorphite
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'Hemimorphite (Zn${4}$Si${2}$O${7}$(OH)${2}\cdot$H${2}
      ↪ $O, $S2_{2}$) Structure'
_aflow_proto 'A2B5CD2_oI40_44_2c_abcde_d_e'
_aflow_params 'a,b/a,c/a,z_{1},z_{2},x_{3},z_{3},x_{4},z_{4},x_{5},z_{5
      ↪ },y_{6},z_{6},y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9}'
_aflow_params_values '8.367,1.28241902713,0.611330225887,0.5912,0.0195,
      ↪ 0.374,0.19,0.4256,0.643,0.305,0.041,0.1669,0.1938,0.1465,0.5076
      ↪ ,0.1602,0.2055,0.6362,0.2047,0.1613,0.0'
_aflow_Strukturbericht '$S2_{2}$'
_aflow_Pearson 'oI40'

_symmetry_space_group_name_H-M "I m m 2"
_symmetry_Int_Tables_number 44

_cell_length_a      8.36700
_cell_length_b      10.73000
```

```
_cell_length_c      5.11500
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,z
3 -x,y,z
4 x,-y,z
5 x+1/2,y+1/2,z+1/2
6 -x+1/2,-y+1/2,z+1/2
7 -x+1/2,y+1/2,z+1/2
8 x+1/2,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 2 a 0.00000 0.00000 0.59120 1.00000
O2 O 2 b 0.00000 0.50000 0.01950 1.00000
H1 H 4 c 0.37400 0.00000 0.19000 1.00000
H2 H 4 c 0.42560 0.00000 0.64300 1.00000
O3 O 4 c 0.30500 0.00000 0.04100 1.00000
O4 O 4 d 0.00000 0.16690 0.19380 1.00000
Si1 Si 4 d 0.00000 0.14650 0.50760 1.00000
O5 O 8 e 0.16020 0.20550 0.63620 1.00000
Zn1 Zn 8 e 0.20470 0.16130 0.00000 1.00000
```

Hemimorphite ($Zn_4Si_2O_7(OH)_2{\cdot}H_2O$, $S2_2$): A2B5CD2_oI40_44_2c_abcde_d_e - POSCAR

```
A2B5CD2_oI40_44_2c_abcde_d_e & a,b/a,c/a,z1,z2,x3,z3,x4,z4,x5,z5,y6,z6,
    ↪ y7,z7,x8,y8,z8,x9,y9,z9 --params=8.367,1.28241902713,
    ↪ 0.611330225887,0.5912,0.0195,0.374,0.19,0.4256,0.643,0.305,
    ↪ 0.041,0.1669,0.1938,0.1465,0.5076,0.1602,0.2055,0.6362,0.2047,
    ↪ 0.1613,0.0 & Imm2 C_{2v}^{20} #44 (abc^3d^2e^2) & oI40 & $S2_{2
    ↪ }$ & H2O5SiZn2 & Hemimorphite & R. J. Hill et al., Zeitschrift
    ↪ f{"u}r Kristallographie - Crystalline Materials 146, 241-259 (
    ↪ 1977)
   1.00000000000000
  -4.18350000000000    5.36500000000000    2.55750000000000
   4.18350000000000   -5.36500000000000    2.55750000000000
   4.18350000000000    5.36500000000000   -2.55750000000000
     H     O     Si     Zn
     4    10     2      4
Direct
   0.19000000000000    0.56400000000000    0.37400000000000    H    (4c)
   0.19000000000000   -0.18400000000000   -0.37400000000000    H    (4c)
   0.64300000000000    1.06860000000000    0.42560000000000    H    (4c)
   0.64300000000000    0.21740000000000   -0.42560000000000    H    (4c)
   0.59120000000000    0.59120000000000    0.00000000000000    O    (2a)
   0.51950000000000    0.01950000000000    0.50000000000000    O    (2b)
   0.04100000000000    0.34600000000000    0.30500000000000    O    (4c)
   0.04100000000000   -0.26400000000000   -0.30500000000000    O    (4c)
   0.36070000000000    0.19380000000000    0.16690000000000    O    (4d)
   0.02690000000000    0.19380000000000   -0.16690000000000    O    (4d)
   0.84170000000000    0.79640000000000    0.36570000000000    O    (8e)
   0.43070000000000    0.47600000000000   -0.36570000000000    O    (8e)
   0.43070000000000    0.79640000000000   -0.04530000000000    O    (8e)
   0.84170000000000    0.47600000000000    0.04530000000000    O    (8e)
   0.65410000000000    0.50760000000000    0.14650000000000    Si   (4d)
   0.36110000000000    0.50760000000000   -0.14650000000000    Si   (4d)
   0.16130000000000    0.20470000000000    0.36600000000000    Zn   (8e)
  -0.16130000000000   -0.20470000000000   -0.36600000000000    Zn   (8e)
  -0.16130000000000    0.20470000000000    0.04340000000000    Zn   (8e)
   0.16130000000000   -0.20470000000000   -0.04340000000000    Zn   (8e)
```

Ferroelectric $NaNO_2$ ($F5_5$): ABC2_oI8_44_a_a_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Sodium nitrite'
_chemical_formula_sum 'N Na O2'

loop_
_publ_author_name
 'M. I. Kay'
 'B. C. Frazier'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 14
_journal_year 1961
_journal_page_first 56
_journal_page_last 57
_publ_Section_title
;
 A neutron diffraction refinement of the low temperature phase of NaNO$_
    ↪ {2}$
;

_aflow_title 'Ferroelectric NaNO$_{2}$ ($F5_{5}$) Structure'
_aflow_proto 'ABC2_oI8_44_a_a_c'
_aflow_params 'a,b/a,c/a,z_{1},z_{2},x_{3},z_{3}'
_aflow_params_values '5.384,0.661218424963,1.03324665676,0.88,0.4147,
    ↪ 0.8059,0.0'
_aflow_Strukturbericht '$F5_{5}$'
_aflow_Pearson 'oI8'

_symmetry_space_group_name_H-M "I m m 2"
_symmetry_Int_Tables_number 44

_cell_length_a      5.38400
_cell_length_b      3.56000
_cell_length_c      5.56300
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,z
3 -x,y,z
4 x,-y,z
5 x+1/2,y+1/2,z+1/2
6 -x+1/2,-y+1/2,z+1/2
7 -x+1/2,y+1/2,z+1/2
8 x+1/2,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
N1 N 2 a 0.00000 0.00000 0.88000 1.00000
Na1 Na 2 a 0.00000 0.00000 0.41470 1.00000
O1 O 4 c 0.80590 0.00000 0.00000 1.00000
```

Ferroelectric $NaNO_2$ ($F5_5$): ABC2_oI8_44_a_a_c - POSCAR

```
ABC2_oI8_44_a_a_c & a,b/a,c/a,z1,z2,x3,z3 --params=5.384,0.661218424963,
    ↪ 1.03324665676,0.88,0.4147,0.8059,0.0 & Imm2 C_{2v}^{20} #44 (a^
    ↪ 2c) & oI8 & $F5_{5}$ & NaNO2 & Sodium nitrite & M. I. Kay and
    ↪ B. C. Frazier, Acta Cryst. 14, 56-57 (1961)
   1.00000000000000
  -2.69200000000000    1.78000000000000    2.78150000000000
   2.69200000000000   -1.78000000000000    2.78150000000000
   2.69200000000000    1.78000000000000   -2.78150000000000
     N     Na     O
     1     1      2
Direct
   0.88000000000000    0.88000000000000    0.00000000000000    N    (2a)
   0.41470000000000    0.41470000000000    0.00000000000000    Na   (2a)
   0.00000000000000    0.80590000000000    0.80590000000000    O    (4c)
   0.00000000000000   -0.80590000000000   -0.80590000000000    O    (4c)
```

$AgNO_2$ ($F5_{12}$): ABC2_oI8_44_a_a_d - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Silver nitrite'
_chemical_formula_sum 'Ag N O2'

loop_
_publ_author_name
 'S. Ohba'
 'Y. Saito'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 37
_journal_year 1981
_journal_page_first 1911
_journal_page_last 1913
_publ_Section_title
;
 Structure of silver(I) nitrite, a redetermination
;

_aflow_title 'AgNO$_{2}$ ($F5_{12}$) Structure'
_aflow_proto 'ABC2_oI8_44_a_a_d'
_aflow_params 'a,b/a,c/a,z_{1},z_{2},y_{3},z_{3}'
_aflow_params_values '3.528,1.74943310658,1.46853741497,0.0,0.4446,
    ↪ 0.1701,0.5747'
_aflow_Strukturbericht '$F5_{12}$'
_aflow_Pearson 'oI8'

_symmetry_space_group_name_H-M "I m m 2"
_symmetry_Int_Tables_number 44

_cell_length_a      3.52800
_cell_length_b      6.17200
_cell_length_c      5.18100
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
```

```
1 x,y,z
2 -x,-y,z
3 -x,y,z
4 x,-y,z
5 x+1/2,y+1/2,z+1/2
6 -x+1/2,-y+1/2,z+1/2
7 -x+1/2,y+1/2,z+1/2
8 x+1/2,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ag1 Ag 2 a 0.00000 0.00000 0.00000 1.00000
N1 N 2 a 0.00000 0.00000 0.44460 1.00000
O1 O 4 d 0.00000 0.17010 0.57470 1.00000
```

$AgNO_2$ ($F5_{12}$): ABC2_oI8_44_a_a_d - POSCAR

```
ABC2_oI8_44_a_a_d & a,b/a,c/a,z1,z2,y3,z3 --params=3.528,1.74943310658,
    ↪ 1.46853741497,0.0,0.4446,0.1701,0.5747 & Imm2 C_{2v}^{20} #44 (
    ↪ a^2d) & oI8 & $F5_{12}$ & AgNO2 & Silver nitrite & S. Ohba and
    ↪ Y. Saito, Acta Crystallogr. Sect. B Struct. Sci. 37, 1911-1913
    ↪ (1981)
   1.00000000000000
  -1.76400000000000    3.08600000000000    2.59050000000000
   1.76400000000000   -3.08600000000000    2.59050000000000
   1.76400000000000    3.08600000000000   -2.59050000000000
    Ag     N     O
     1     1     2
Direct
   0.00000000000000    0.00000000000000    0.00000000000000   Ag   (2a)
   0.44460000000000    0.44460000000000    0.00000000000000   N    (2a)
   0.74480000000000    0.57470000000000    0.17010000000000   O    (4d)
   0.40460000000000    0.57470000000000   -0.17010000000000   O    (4d)
```

*B*30 (MgZn?): AB_oI48_44_6d_ab2cde - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'MgZn'
_chemical_formula_sum 'Mg Zn'

loop_
_publ_author_name
 'C. Hermann'
 'O. Lohrmann'
 'H. Philipp'
_journal_year 1937
_publ_Section_title
;
 Strukturbericht Band II 1928-1932
;

_aflow_title '$B30$ (MgZn?) Structure'
_aflow_proto 'AB_oI48_44_6d_ab2cde'
_aflow_params 'a,b/a,c/a,z_{1},z_{2},x_{3},z_{3},x_{4},z_{4},y_{5},z_{5
    ↪ },y_{6},z_{6},y_{7},z_{7},y_{8},z_{8},y_{9},z_{9},y_{10},z_{10
    ↪ },y_{11},z_{11},x_{12},y_{12},z_{12}'
_aflow_params_values '7.53776,2.27653838806,1.0,-0.04167,-0.04167,0.25,
    ↪ 0.70833,0.75,0.20833,0.09375,0.45833,0.59375,0.45833,0.84375,
    ↪ 0.79167,0.34375,0.79167,0.59375,0.125,0.15625,0.125,0.75,
    ↪ 0.29167,0.75,0.25,0.04167'
_aflow_Strukturbericht '$B30$'
_aflow_Pearson 'oI48'

_symmetry_space_group_name_H-M "I m m 2"
_symmetry_Int_Tables_number 44

_cell_length_a    7.53776
_cell_length_b    17.16000
_cell_length_c    7.53776
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,z
3 -x,y,z
4 x,-y,z
5 x+1/2,y+1/2,z+1/2
6 -x+1/2,-y+1/2,z+1/2
7 -x+1/2,y+1/2,z+1/2
8 x+1/2,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Zn1 Zn 2 a 0.00000 0.00000 -0.04167 1.00000
Zn2 Zn 2 b 0.00000 0.50000 -0.04167 1.00000
Zn3 Zn 4 c 0.25000 0.00000 0.70833 1.00000
Zn4 Zn 4 c 0.75000 0.00000 0.20833 1.00000
Mg1 Mg 4 d 0.00000 0.09375 0.45833 1.00000
Mg2 Mg 4 d 0.00000 0.59375 0.45833 1.00000
Mg3 Mg 4 d 0.00000 0.84375 0.79167 1.00000
Mg4 Mg 4 d 0.00000 0.34375 0.79167 1.00000
Mg5 Mg 4 d 0.00000 0.59375 0.12500 1.00000
Mg6 Mg 4 d 0.00000 0.15625 0.12500 1.00000
Zn5 Zn 4 d 0.00000 0.75000 0.29167 1.00000
Zn6 Zn 8 e 0.75000 0.25000 0.04167 1.00000
```

*B*30 (MgZn?): AB_oI48_44_6d_ab2cde - POSCAR

```
AB_oI48_44_6d_ab2cde & a,b/a,c/a,z1,z2,x3,z3,x4,z4,y5,z5,y6,z6,y7,z7,y8,
    ↪ z8,y9,z9,y10,z10,y11,z11,x12,y12,z12 --params=7.53776,
    ↪ 2.27653838806,1.0,-0.04167,-0.04167,0.25,0.70833,0.75,0.20833,
    ↪ 0.09375,0.45833,0.59375,0.45833,0.84375,0.79167,0.34375,0.79167
    ↪ ,0.59375,0.125,0.15625,0.125,0.75,0.29167,0.75,0.25,0.04167 &
    ↪ Imm2 C_{2v}^{20} #44 (abc^2d^7e) & oI48 & $B30$ & MgZn & MgZn &
    ↪ C. Hermann and O. Lohrmann and H. Philipp, (1937)
   1.00000000000000
  -3.76888000000000    8.58000000000000    3.76888000000000
   3.76888000000000   -8.58000000000000    3.76888000000000
   3.76888000000000    8.58000000000000   -3.76888000000000
    Mg     Zn
    12     12
Direct
   0.55208000000000    0.45833000000000    0.09375000000000   Mg   (4d)
   0.36458000000000    0.45833000000000   -0.09375000000000   Mg   (4d)
   1.05208000000000    0.45833000000000    0.59375000000000   Mg   (4d)
  -0.13542000000000    0.45833000000000   -0.59375000000000   Mg   (4d)
   1.63542000000000    0.79167000000000    0.84375000000000   Mg   (4d)
  -0.05208000000000    0.79167000000000   -0.84375000000000   Mg   (4d)
   1.13542000000000    0.79167000000000    0.34375000000000   Mg   (4d)
   0.44792000000000    0.79167000000000   -0.34375000000000   Mg   (4d)
   0.71875000000000    0.12500000000000    0.59375000000000   Mg   (4d)
  -0.46875000000000    0.12500000000000   -0.59375000000000   Mg   (4d)
   0.28125000000000    0.12500000000000    0.15625000000000   Mg   (4d)
  -0.03125000000000    0.12500000000000   -0.15625000000000   Mg   (4d)
  -0.04167000000000   -0.04167000000000    0.00000000000000   Zn   (2a)
   0.45833000000000   -0.04167000000000    0.50000000000000   Zn   (2b)
   0.70833000000000    0.95833000000000    0.25000000000000   Zn   (4c)
   0.70833000000000    0.45833000000000   -0.25000000000000   Zn   (4c)
   0.20833000000000    0.95833000000000    0.75000000000000   Zn   (4c)
   0.20833000000000   -0.54167000000000   -0.75000000000000   Zn   (4c)
   1.04167000000000    0.29167000000000    0.75000000000000   Zn   (4d)
  -0.45833000000000    0.29167000000000   -0.75000000000000   Zn   (4d)
   0.29167000000000    0.79167000000000    1.00000000000000   Zn   (8e)
  -0.20833000000000   -0.70833000000000   -1.00000000000000   Zn   (8e)
  -0.20833000000000    0.79167000000000    0.50000000000000   Zn   (8e)
   0.29167000000000   -0.70833000000000   -0.50000000000000   Zn   (8e)
```

$Nb_2Zr_6O_{17}$: A2B17C6_oI100_46_ab_b8c_3c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Nb2O17Zr6'
_chemical_formula_sum 'Nb2 O17 Zr6'

loop_
_publ_author_name
 'J. Galy'
 'R. S. Roth'
_journal_name_full_name
;
 Journal of Solid State Chemistry
;
_journal_volume 7
_journal_year 1973
_journal_page_first 277
_journal_page_last 285
_publ_Section_title
;
 The Crystal Structure of Nb$_{2}$Zr$_{6}$O$_{17}$
;

# Found in Crystal structure solution for the $A_{6}$$B_{2}$O$_{17}$ (
    ↪ $A$ = Zr, Hf; $B$ = Nb, Ta) superstructure, 2019

_aflow_title 'Nb$_{2}$Zr$_{6}$O$_{17}$ Structure'
_aflow_proto 'A2B17C6_oI100_46_ab_b8c_3c'
_aflow_params 'a,b/a,c/a,z_{1},y_{2},z_{2},y_{3},z_{3},x_{4},y_{4},z_{4
    ↪ },x_{5},y_{5},z_{5},x_{6},y_{6},z_{6},x_{7},y_{7},z_{7},x_{8},
    ↪ y_{8},z_{8},x_{9},y_{9},z_{9},x_{10},y_{10},z_{10},x_{11},y_{11
    ↪ },z_{11},x_{12},y_{12},z_{12},x_{13},y_{13},z_{13},x_{14},y_{14
    ↪ },z_{14}'
_aflow_params_values '40.92,0.120478983382,0.128787878788,0.0,0.5151,
    ↪ 0.5087,0.3564,0.1537,0.1976,0.3485,0.4426,0.2195,0.8082,0.311,
    ↪ 0.1589,0.7925,0.3251,0.0956,0.7915,0.2785,0.0362,0.7505,0.3797,
    ↪ 0.0827,0.2994,0.4251,0.0249,0.2973,0.1839,0.1419,0.3421,0.1826,
    ↪ 0.065,-0.0152,0.5894,0.1278,0.5138,0.5136,0.18965,-0.0151,
    ↪ 0.5861'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oI100'

_symmetry_space_group_name_H-M "I m a 2"
_symmetry_Int_Tables_number 46

_cell_length_a    40.92000
_cell_length_b    4.93000
_cell_length_c    5.27000
_cell_angle_alpha 90.00000
```

```
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,z
3 -x+1/2,y,z
4 x+1/2,-y,z
5 x+1/2,y+1/2,z+1/2
6 -x+1/2,-y+1/2,z+1/2
7 -x,y+1/2,z+1/2
8 x,-y+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Nb1 Nb 4 a 0.00000 0.00000 0.00000 1.00000
Nb2 Nb 4 b 0.25000 0.51510 0.50870 1.00000
O1 O 4 b 0.25000 0.35640 0.15370 1.00000
O2 O 8 c 0.19760 0.34850 0.44260 1.00000
O3 O 8 c 0.21950 0.80820 0.31100 1.00000
O4 O 8 c 0.15890 0.79250 0.32510 1.00000
O5 O 8 c 0.09560 0.79150 0.27850 1.00000
O6 O 8 c 0.03620 0.75050 0.37970 1.00000
O7 O 8 c 0.08270 0.29940 0.42510 1.00000
O8 O 8 c 0.02490 0.29730 0.18390 1.00000
O9 O 8 c 0.14190 0.34210 0.18260 1.00000
Zr1 Zr 8 c 0.06500 -0.01520 0.58940 1.00000
Zr2 Zr 8 c 0.12780 0.51380 0.51360 1.00000
Zr3 Zr 8 c 0.18965 -0.01510 0.58610 1.00000
```

$Nb_2Zr_6O_{17}$: A2B17C6_oI100_46_ab_b8c_3c - POSCAR

```
A2B17C6_oI100_46_ab_b8c_3c & a,b/a,c/a,z1,y2,z2,y3,z3,x4,y4,z4,x5,y5,z5,
    ↪ x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10,x11,y11,z11,x12
    ↪ ,y12,z12,x13,y13,z13,x14,y14,z14 --params=40.92,0.120478983382,
    ↪ 0.128787878788,0.0,0.5151,0.5087,0.3564,0.1537,0.1976,0.3485,
    ↪ 0.4426,0.2195,0.8082,0.311,0.1589,0.7925,0.3251,0.0956,0.7915,
    ↪ 0.2785,0.0362,0.7505,0.3797,0.0827,0.2994,0.4251,0.0249,0.2973,
    ↪ 0.1839,0.1419,0.3421,0.1826,0.065,-0.0152,0.5894,0.1278,0.5138,
    ↪ 0.5136,0.18965,-0.0151,0.5861 & Ima2 C_{2v}^{22} #46 (ab^2c^11)
    ↪ & oI100 & None & Nb2O17Zr6 & Nb2O17Zr6 & J. Galy and R. S.
    ↪ Roth, J. Solid State Chem. 7, 277-285 (1973)
   1.00000000000000
 -20.46000000000000    2.46500000000000    2.63500000000000
  20.46000000000000   -2.46500000000000    2.63500000000000
  20.46000000000000    2.46500000000000   -2.63500000000000
     Nb      O      Zr
      4     34      12
Direct
    0.00000000000000    0.00000000000000    0.00000000000000   Nb   (4a)
    0.00000000000000    0.50000000000000    0.50000000000000   Nb   (4a)
    1.02380000000000    0.75870000000000    0.76510000000000   Nb   (4b)
   -0.00640000000000    1.25870000000000    0.23490000000000   Nb   (4b)
    0.51010000000000    0.40370000000000    0.60640000000000    O   (4b)
   -0.20270000000000    0.90370000000000    0.39360000000000    O   (4b)
    0.79110000000000    0.64020000000000    0.54610000000000    O   (8c)
    0.09410000000000    0.24500000000000   -0.54610000000000    O   (8c)
    0.09410000000000    1.14020000000000    0.34910000000000    O   (8c)
    0.79110000000000    0.74500000000000    0.65090000000000    O   (8c)
    1.11920000000000    0.53050000000000    1.02770000000000    O   (8c)
   -0.49720000000000    0.09150000000000   -1.02770000000000    O   (8c)
   -0.49720000000000    1.03050000000000   -0.08870000000000    O   (8c)
    1.11920000000000    0.59150000000000    1.08870000000000    O   (8c)
    1.11760000000000    0.48400000000000    0.95140000000000    O   (8c)
   -0.46740000000000    0.16620000000000   -0.95140000000000    O   (8c)
   -0.46740000000000    0.98400000000000   -0.13360000000000    O   (8c)
    1.11760000000000    0.66620000000000    1.13360000000000    O   (8c)
    1.07000000000000    0.37410000000000    0.88710000000000    O   (8c)
   -0.51300000000000    0.18290000000000   -0.88710000000000    O   (8c)
   -0.51300000000000    0.87410000000000   -0.19590000000000    O   (8c)
    1.07000000000000    0.68290000000000    1.19590000000000    O   (8c)
    1.13020000000000    0.41590000000000    0.78670000000000    O   (8c)
   -0.37080000000000    0.34350000000000   -0.78670000000000    O   (8c)
   -0.37080000000000    0.91590000000000   -0.21430000000000    O   (8c)
    1.13020000000000    0.84350000000000    1.21430000000000    O   (8c)
    0.72450000000000    0.50780000000000    0.38210000000000    O   (8c)
    0.12570000000000    0.34240000000000   -0.38210000000000    O   (8c)
    0.12570000000000    1.00780000000000    0.28330000000000    O   (8c)
    0.72450000000000    0.84240000000000    0.71670000000000    O   (8c)
    0.48120000000000    0.20880000000000    0.32220000000000    O   (8c)
   -0.11340000000000    0.15900000000000   -0.32220000000000    O   (8c)
   -0.11340000000000    0.70880000000000    0.22760000000000    O   (8c)
    0.48120000000000    0.65900000000000    0.77240000000000    O   (8c)
    0.52470000000000    0.32450000000000    0.48400000000000    O   (8c)
   -0.15950000000000    0.04070000000000   -0.48400000000000    O   (8c)
   -0.15950000000000    0.82450000000000    0.29980000000000    O   (8c)
    0.52470000000000    0.54070000000000    0.70020000000000    O   (8c)
    0.57420000000000    0.65440000000000    0.04980000000000   Zr   (8c)
    0.60460000000000    0.52440000000000   -0.04980000000000   Zr   (8c)
    0.60460000000000    1.15440000000000    0.58020000000000   Zr   (8c)
    0.57420000000000    1.02440000000000    0.41980000000000   Zr   (8c)
    1.02740000000000    0.64140000000000    0.64160000000000   Zr   (8c)
   -0.00020000000000    0.38580000000000   -0.64160000000000   Zr   (8c)
   -0.00020000000000    1.14140000000000    0.11400000000000   Zr   (8c)
    1.02740000000000    0.88580000000000    0.88600000000000   Zr   (8c)
    0.57100000000000    0.77575000000000    0.17455000000000   Zr   (8c)
    0.60120000000000    0.39645000000000   -0.17455000000000   Zr   (8c)
    0.60120000000000    1.27575000000000    0.70475000000000   Zr   (8c)
    0.57100000000000    0.89645000000000    0.29525000000000   Zr   (8c)
```

Parkerite ($Ni_3Bi_2S_2$): AB2C_oP8_51_e_be_f - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Parkerite'
_chemical_formula_sum 'Bi Ni2 S'

loop_
_publ_author_name
 'M. E. Fleet'
_journal_name_full_name
;
 American Mineralogist
;
_journal_volume 58
_journal_year 1973
_journal_page_first 435
_journal_page_last 439
_publ_Section_title
;
 The Crystal Structure of Parkerite (Ni$_{3}$Bi$_{2}$S$_{2}$)
;

_aflow_title 'Parkerite (Ni$_{3}$Bi$_{2}$S$_{2}$) Structure'
_aflow_proto 'AB2C_oP8_51_e_be_f'
_aflow_params 'a,b/a,c/a,z_{2},z_{3},z_{4}'
_aflow_params_values '5.545,0.730748422002,1.03354373309,0.7423,0.204,
    ↪ 0.256'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP8'

_symmetry_space_group_name_H-M "P 21/m 2/m 2/a"
_symmetry_Int_Tables_number 51

_cell_length_a     5.54500
_cell_length_b     4.05200
_cell_length_c     5.73100
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y,-z
3 -x,y,-z
4 -x+1/2,-y,z
5 -x,-y,-z
6 -x+1/2,y,z
7 x,-y,z
8 x+1/2,y,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ni1 Ni 2 b 0.00000 0.50000 0.00000 1.00000
Bi1 Bi 2 e 0.25000 0.00000 0.74230 1.00000
Ni2 Ni 2 e 0.25000 0.00000 0.20400 0.50000
S1 S 2 f 0.25000 0.50000 0.25600 1.00000
```

Parkerite ($Ni_3Bi_2S_2$): AB2C_oP8_51_e_be_f - POSCAR

```
AB2C_oP8_51_e_be_f & a,b/a,c/a,z2,z3,z4 --params=5.545,0.730748422002,
    ↪ 1.03354373309,0.7423,0.204,0.256 & Pmma D_{2h}^{5} #51 (be^2f)
    ↪ & oP8 & None & Bi2Ni3S2 & Parkerite & M. E. Fleet, Am. Mineral.
    ↪ 58, 435-439 (1973)
   1.00000000000000
   5.54500000000000    0.00000000000000    0.00000000000000
   0.00000000000000    4.05200000000000    0.00000000000000
   0.00000000000000    0.00000000000000    5.73100000000000
     Bi      Ni       S
      2       4       2
Direct
    0.25000000000000    0.00000000000000    0.74230000000000   Bi   (2e)
    0.75000000000000    0.00000000000000   -0.74230000000000   Bi   (2e)
    0.00000000000000    0.50000000000000    0.00000000000000   Ni   (2b)
    0.50000000000000    0.50000000000000    0.00000000000000   Ni   (2b)
    0.25000000000000    0.00000000000000    0.20400000000000   Ni   (2e)
    0.75000000000000    0.00000000000000   -0.20400000000000   Ni   (2e)
    0.25000000000000    0.50000000000000    0.25600000000000    S   (2f)
    0.75000000000000    0.50000000000000   -0.25600000000000    S   (2f)
```

$LiNb_6O_{15}F$: ABC6D15_oP46_51_f_d_2e2i_aef4i2j - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'FLiNb6O15'
_chemical_formula_sum 'F Li Nb6 O15'
```

```
loop_
_publ_author_name
 'M. Lundberg'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 19
_journal_year 1965
_journal_page_first 2274
_journal_page_last 2284
_publ_Section_title
;
 The Crystal Structure of LiNb$_{6}$O$_{15}$F
;

_aflow_title 'LiNb$_{6}$O$_{15}$F Structure'
_aflow_proto 'ABC6D15_oP46_51_f_d_2e2i_aef4i2j'
_aflow_params 'a,b/a,c/a,z_{3},z_{4},z_{5},z_{6},z_{7},x_{8},z_{8},x_{9
      ↪ },z_{9},x_{10},z_{10},x_{11},z_{11},x_{12},z_{12},x_{13},z_{13
      ↪ },x_{14},z_{14},x_{15},z_{15}'
_aflow_params_values '16.635,0.238292756237,0.534295160806,0.0578,0.6874
      ↪ ,0.4582,0.6922,0.0524,0.1305,0.3973,0.0593,0.8146,0.1309,0.6149
      ↪ ,0.1744,0.8823,0.1636,0.1963,0.0206,0.3386,0.1286,0.3998,0.0583
      ↪ ,0.817'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP46'

_symmetry_space_group_name_H-M "P 21/m 2/m 2/a"
_symmetry_Int_Tables_number 51

_cell_length_a      16.63500
_cell_length_b      3.96400
_cell_length_c      8.88800
_cell_angle_alpha   90.00000
_cell_angle_beta    90.00000
_cell_angle_gamma   90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y,-z
3 -x,y,-z
4 -x+1/2,-y,z
5 -x,-y,-z
6 -x+1/2,y,z
7 x,-y,z
8 x+1/2,y,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 2 a 0.00000 0.00000 0.00000 1.00000
Li1 Li 2 d 0.00000 0.50000 0.50000 1.00000
Nb1 Nb 2 e 0.25000 0.00000 0.05780 1.00000
Nb2 Nb 2 e 0.25000 0.00000 0.68740 1.00000
O2 O 2 e 0.25000 0.00000 0.45820 1.00000
F1 F 2 f 0.25000 0.50000 0.69220 1.00000
O3 O 2 f 0.25000 0.50000 0.05240 1.00000
Nb3 Nb 4 i 0.13050 0.00000 0.39730 1.00000
Nb4 Nb 4 i 0.05930 0.00000 0.81460 1.00000
O4 O 4 i 0.13090 0.00000 0.61490 1.00000
O5 O 4 i 0.17440 0.00000 0.88230 1.00000
O6 O 4 i 0.16360 0.00000 0.19630 1.00000
O7 O 4 i 0.02060 0.00000 0.33860 1.00000
O8 O 4 j 0.12860 0.50000 0.39980 1.00000
O9 O 4 j 0.05830 0.50000 0.81700 1.00000
```

LiNb$_6$O$_{15}$F: ABC6D15_oP46_51_f_d_2e2i_aef4i2j - POSCAR

```
ABC6D15_oP46_51_f_d_2e2i_aef4i2j & a,b/a,c/a,z3,z4,z5,z6,z7,x8,z8,x9,z9,
      ↪ x10,z10,x11,z11,x12,z12,x13,z13,x14,z14,x15,z15 --params=16.635
      ↪ ,0.238292756237,0.534295160806,0.0578,0.6874,0.4582,0.6922,
      ↪ 0.0524,0.1305,0.3973,0.0593,0.8146,0.1309,0.6149,0.1744,0.8823,
      ↪ 0.1636,0.1963,0.0206,0.3386,0.1286,0.3998,0.0583,0.817 & Pmma
      ↪ D_{2h}^{5} #51 (ade^3f^2i^6j^2) & oP46 & None & FLiNb6O15 &
      ↪ FLiNb6O15 & M. Lundberg, Acta Chem. Scand. 19, 2274-2284 (1965)
   1.00000000000000
  16.63500000000000     0.00000000000000     0.00000000000000
   0.00000000000000     3.96400000000000     0.00000000000000
   0.00000000000000     0.00000000000000     8.88800000000000
     F    Li    Nb     O
     2     2    12    30
Direct
   0.25000000000000     0.50000000000000     0.69220000000000     F    (2f)
   0.75000000000000     0.50000000000000    -0.69220000000000     F    (2f)
   0.00000000000000     0.50000000000000     0.50000000000000    Li    (2d)
   0.50000000000000     0.50000000000000     0.50000000000000    Li    (2d)
   0.25000000000000     0.00000000000000     0.05780000000000    Nb    (2e)
   0.75000000000000     0.00000000000000    -0.05780000000000    Nb    (2e)
   0.25000000000000     0.00000000000000     0.68740000000000    Nb    (2e)
   0.75000000000000     0.00000000000000    -0.68740000000000    Nb    (2e)
   0.13050000000000     0.00000000000000     0.39730000000000    Nb    (4i)
   0.36950000000000     0.00000000000000     0.39730000000000    Nb    (4i)
  -0.13050000000000     0.00000000000000    -0.39730000000000    Nb    (4i)
   0.63050000000000     0.00000000000000    -0.39730000000000    Nb    (4i)
   0.05930000000000     0.00000000000000     0.81460000000000    Nb    (4i)
   0.44070000000000     0.00000000000000     0.81460000000000    Nb    (4i)
  -0.05930000000000     0.00000000000000    -0.81460000000000    Nb    (4i)
   0.55930000000000     0.00000000000000    -0.81460000000000    Nb    (4i)
   0.00000000000000     0.00000000000000     0.00000000000000     O    (2a)
   0.50000000000000     0.00000000000000     0.00000000000000     O    (2a)
   0.25000000000000     0.00000000000000     0.45820000000000     O    (2e)
   0.75000000000000     0.00000000000000    -0.45820000000000     O    (2e)
   0.25000000000000     0.50000000000000     0.05240000000000     O    (2f)
   0.75000000000000     0.50000000000000    -0.05240000000000     O    (2f)
   0.13090000000000     0.00000000000000     0.61490000000000     O    (4i)
   0.36910000000000     0.00000000000000     0.61490000000000     O    (4i)
  -0.13090000000000     0.00000000000000    -0.61490000000000     O    (4i)
   0.63090000000000     0.00000000000000    -0.61490000000000     O    (4i)
   0.17440000000000     0.00000000000000     0.88230000000000     O    (4i)
   0.32560000000000     0.00000000000000     0.88230000000000     O    (4i)
  -0.17440000000000     0.00000000000000    -0.88230000000000     O    (4i)
   0.67440000000000     0.00000000000000    -0.88230000000000     O    (4i)
   0.16360000000000     0.00000000000000     0.19630000000000     O    (4i)
   0.33640000000000     0.00000000000000     0.19630000000000     O    (4i)
  -0.16360000000000     0.00000000000000    -0.19630000000000     O    (4i)
   0.66360000000000     0.00000000000000    -0.19630000000000     O    (4i)
   0.02060000000000     0.00000000000000     0.33860000000000     O    (4i)
   0.47940000000000     0.00000000000000     0.33860000000000     O    (4i)
  -0.02060000000000     0.00000000000000    -0.33860000000000     O    (4i)
   0.52060000000000     0.00000000000000    -0.33860000000000     O    (4i)
   0.12860000000000     0.50000000000000     0.39980000000000     O    (4j)
   0.37140000000000     0.50000000000000     0.39980000000000     O    (4j)
  -0.12860000000000     0.50000000000000    -0.39980000000000     O    (4j)
   0.62860000000000     0.50000000000000    -0.39980000000000     O    (4j)
   0.05830000000000     0.50000000000000     0.81700000000000     O    (4j)
   0.44170000000000     0.50000000000000     0.81700000000000     O    (4j)
  -0.05830000000000     0.50000000000000    -0.81700000000000     O    (4j)
   0.55830000000000     0.50000000000000    -0.81700000000000     O    (4j)
```

Carnallite [Mg(H$_2$O)$_6$KCl$_3$]: A3B12CDE6_oP276_52_d4e_18e_ce_de_2d8e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Carnallite'
_chemical_formula_sum 'Cl3 H12 K Mg O6'

loop_
_publ_author_name
 'E. O. Schlemper'
 'P. K. {Sen Gupta}'
 'T. Zoltai'
_journal_name_full_name
;
 American Mineralogist
;
_journal_volume 70
_journal_year 1985
_journal_page_first 1309
_journal_page_last 1313
_publ_Section_title
;
 Refinement of the structure of carnallite, Mg(H$_{2}$O)$_{6}$KCl$_{3}$
;

_aflow_title 'Carnallite [Mg(H$_{2}$O)$_{6}$KCl$_{3}$] Structure'
_aflow_proto 'A3B12CDE6_oP276_52_d4e_18e_ce_de_2d8e'
_aflow_params 'a,b/a,c/a,z_{1},x_{2},x_{3},x_{4},x_{5},x_{6},y_{6},z_{6
      ↪ },x_{7},y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{10},
      ↪ y_{10},z_{10},x_{11},y_{11},z_{11},x_{12},y_{12},z_{12},x_{13},
      ↪ y_{13},z_{13},x_{14},y_{14},z_{14},x_{15},y_{15},z_{15},x_{16},
      ↪ y_{16},z_{16},x_{17},y_{17},z_{17},x_{18},y_{18},z_{18},x_{19},
      ↪ y_{19},z_{19},x_{20},y_{20},z_{20},x_{21},y_{21},z_{21},x_{22},
      ↪ y_{22},z_{22},x_{23},y_{23},z_{23},x_{24},y_{24},z_{24},x_{25},
      ↪ y_{25},z_{25},x_{26},y_{26},z_{26},x_{27},y_{27},z_{27},x_{28},
      ↪ y_{28},z_{28},x_{29},y_{29},z_{29},x_{30},y_{30},z_{30},x_{31},
      ↪ y_{31},z_{31},x_{32},y_{32},z_{32},x_{33},y_{33},z_{33},x_{34},
      ↪ y_{34},z_{34},x_{35},y_{35},z_{35},x_{36},y_{36},z_{36},x_{37},
      ↪ y_{37},z_{37}'
_aflow_params_values '16.119,1.39413114958,0.592530554005,0.2514,0.76061
      ↪ ,0.24343,0.3692,0.1162,0.16583,0.57525,-0.01474,0.16928,0.58113
      ↪ ,0.48834,0.41462,0.58176,0.25235,0.01987,0.74957,-0.02314,0.604
      ↪ ,0.75,0.579,0.412,0.751,0.313,0.737,0.673,0.543,0.25,0.722,
      ↪ 0.505,0.734,0.848,0.587,0.303,0.845,0.335,0.514,0.509,0.326,
      ↪ 0.417,0.494,0.606,0.768,0.572,0.345,0.264,-0.066,0.682,0.399,-
      ↪ 0.01,0.568,0.422,-0.058,0.5,0.359,0.817,0.673,0.429,0.811,0.623
      ↪ ,0.565,-0.099,0.679,0.573,0.604,0.695,0.438,0.66,0.561,0.411,
      ↪ 0.615,0.475,0.0886,0.65668,0.74995,0.42094,0.58981,0.7477,
      ↪ 0.2554,0.70704,0.4384,0.2673,0.66936,0.1506,0.4466,0.51041,0.65
      ↪ ,0.2967,0.57117,0.7349,0.4259,0.54662,-0.0646,0.3956,0.66912,
      ↪ 0.8444,0.5448,0.60769,0.7631,0.4179,0.63155,0.558'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP276'

_symmetry_space_group_name_H-M "P 2/n 21/n 2/a"
_symmetry_Int_Tables_number 52

_cell_length_a      16.11900
_cell_length_b      22.47200
_cell_length_c      9.55100
_cell_angle_alpha   90.00000
_cell_angle_beta    90.00000
_cell_angle_gamma   90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y+1/2,-z+1/2
3 -x+1/2,y+1/2,-z+1/2
```

```
4 -x+1/2,-y,z
5 -x,-y,-z
6 -x,y+1/2,z+1/2
7 x+1/2,-y+1/2,z+1/2
8 x+1/2,y,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
K1 K 4 c 0.25000 0.00000 0.25140 1.00000
Cl1 Cl 4 d 0.76061 0.25000 0.25000 1.00000
Mg1 Mg 4 d 0.24343 0.25000 0.25000 1.00000
O1 O 4 d 0.36920 0.25000 0.25000 1.00000
O2 O 4 d 0.11620 0.25000 0.25000 1.00000
Cl2 Cl 8 e 0.16583 0.57525 -0.01474 1.00000
Cl3 Cl 8 e 0.16928 0.58113 0.48834 1.00000
Cl4 Cl 8 e 0.41462 0.58176 0.25235 1.00000
Cl5 Cl 8 e 0.01987 0.74957 -0.02314 1.00000
H1 H 8 e 0.60400 0.75000 0.57900 1.00000
H2 H 8 e 0.41200 0.75100 0.31300 1.00000
H3 H 8 e 0.73700 0.67300 0.54300 1.00000
H4 H 8 e 0.25000 0.72200 0.50500 1.00000
H5 H 8 e 0.73400 0.84800 0.58700 1.00000
H6 H 8 e 0.30300 0.84500 0.33500 1.00000
H7 H 8 e 0.51400 0.50900 0.32600 1.00000
H8 H 8 e 0.41700 0.49400 0.60600 1.00000
H9 H 8 e 0.76800 0.57200 0.34500 1.00000
H10 H 8 e 0.26400 -0.06600 0.68200 1.00000
H11 H 8 e 0.39900 -0.01000 0.56800 1.00000
H12 H 8 e 0.42200 -0.05800 0.50000 1.00000
H13 H 8 e 0.35900 0.81700 0.67300 1.00000
H14 H 8 e 0.42900 0.81100 0.62300 1.00000
H15 H 8 e 0.56500 -0.09900 0.67900 1.00000
H16 H 8 e 0.57300 0.60400 0.69500 1.00000
H17 H 8 e 0.43800 0.66000 0.56100 1.00000
H18 H 8 e 0.41100 0.61500 0.47500 1.00000
K2 K 8 e 0.08860 0.65668 0.74995 1.00000
Mg2 Mg 8 e 0.42094 0.58981 0.74770 1.00000
O3 O 8 e 0.25540 0.70704 0.43840 1.00000
O4 O 8 e 0.26730 0.66936 0.15060 1.00000
O5 O 8 e 0.44660 0.51041 0.65000 1.00000
O6 O 8 e 0.29670 0.57117 0.73490 1.00000
O7 O 8 e 0.42590 0.54662 -0.06460 1.00000
O8 O 8 e 0.39560 0.66912 0.84440 1.00000
O9 O 8 e 0.54480 0.60769 0.76310 1.00000
O10 O 8 e 0.41790 0.63155 0.55800 1.00000
```

Carnallite [$Mg(H_2O)_6KCl_3$]: A3B12CDE6_oP276_52_d4e_18e_ce_de_2d8e - POSCAR

```
A3B12CDE6_oP276_52_d4e_18e_ce_de_2d8e & a,b/a,c/a,z1,x2,x3,x4,x5,x6,y6,
    ↪ z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10,x11,y11,z11,x12,y12,
    ↪ z12,x13,y13,z13,x14,y14,z14,x15,y15,z15,x16,y16,z16,x17,y17,z17
    ↪ ,x18,y18,z18,x19,y19,z19,x20,y20,z20,x21,y21,z21,x22,y22,z22,
    ↪ x23,y23,z23,x24,y24,z24,x25,y25,z25,x26,y26,z26,x27,y27,z27,x28
    ↪ ,y28,z28,x29,y29,z29,x30,y30,z30,x31,y31,z31,x32,y32,z32,x33,
    ↪ y33,z33,x34,y34,z34,x35,y35,z35,x36,y36,z36,x37,y37,z37 --
    ↪ params=16.119,1.39413114958,0.592530554005,0.2514,0.76061,
    ↪ 0.24343,0.3692,0.1162,0.16583,0.57525,-0.01474,0.16928,0.58113,
    ↪ 0.48834,0.41462,0.58176,0.25235,0.01987,0.74957,-0.02314,0.604,
    ↪ 0.75,0.579,0.412,0.751,0.313,0.737,0.673,0.543,0.25,0.722,0.505
    ↪ ,0.734,0.848,0.587,0.303,0.845,0.335,0.514,0.509,0.326,0.417,
    ↪ 0.494,0.606,0.768,0.572,0.345,0.264,-0.066,0.682,0.399,-0.01,
    ↪ 0.568,0.422,-0.058,0.5,0.359,0.817,0.673,0.429,0.811,0.623,
    ↪ 0.565,-0.099,0.679,0.573,0.604,0.695,0.438,0.66,0.561,0.411,
    ↪ 0.615,0.475,0.0886,0.65668,0.74995,0.42094,0.58981,0.7477,
    ↪ 0.2554,0.70704,0.4384,0.2673,0.66936,0.1506,0.4466,0.51041,0.65
    ↪ ,0.2967,0.57117,0.7349,0.4259,0.54662,-0.0646,0.3956,0.66912,
    ↪ 0.8444,0.5448,0.60769,0.7631,0.4179,0.63155,0.558 & Pnna D_{2h
    ↪ }^{6} #52 (cd^4e^32) & oP276 & None & Cl3H12KMgO6 & Carnallite
    ↪ & E. O. Schlemper and P. K. {Sen Gupta} and T. Zoltai, Am.
    ↪ Mineral. 70, 1309-1313 (1985)
   1.00000000000000
  16.11900000000000    0.00000000000000    0.00000000000000
   0.00000000000000   22.47200000000000    0.00000000000000
   0.00000000000000    0.00000000000000    9.55100000000000
     Cl     H     K    Mg     O
     36   144    12    12    72
Direct
   0.76061000000000    0.25000000000000    0.25000000000000   Cl   (4d)
  -0.26061000000000    0.75000000000000    0.25000000000000   Cl   (4d)
  -0.76061000000000    0.75000000000000    0.75000000000000   Cl   (4d)
   1.26061000000000    0.25000000000000    0.75000000000000   Cl   (4d)
   0.16583000000000    0.57525000000000   -0.01474000000000   Cl   (8e)
   0.33417000000000   -0.57525000000000   -0.01474000000000   Cl   (8e)
   0.33417000000000    1.07525000000000    0.51474000000000   Cl   (8e)
   0.16583000000000   -0.07525000000000    0.51474000000000   Cl   (8e)
  -0.16583000000000   -0.57525000000000    0.01474000000000   Cl   (8e)
   0.66583000000000    0.57525000000000    0.01474000000000   Cl   (8e)
   0.66583000000000   -0.07525000000000    0.48526000000000   Cl   (8e)
  -0.16583000000000    1.07525000000000    0.48526000000000   Cl   (8e)
   0.16928000000000    0.58113000000000    0.48834000000000   Cl   (8e)
   0.33072000000000   -0.58113000000000    0.48834000000000   Cl   (8e)
   0.33072000000000    1.08113000000000    0.01166000000000   Cl   (8e)
   0.16928000000000   -0.08113000000000    0.01166000000000   Cl   (8e)
  -0.16928000000000   -0.58113000000000   -0.48834000000000   Cl   (8e)
   0.66928000000000    0.58113000000000   -0.48834000000000   Cl   (8e)
   0.66928000000000   -0.08113000000000    0.98834000000000   Cl   (8e)
  -0.16928000000000    1.08113000000000    0.98834000000000   Cl   (8e)
   0.41462000000000    0.58176000000000    0.25235000000000   Cl   (8e)
   0.08538000000000   -0.58176000000000    0.25235000000000   Cl   (8e)
   0.08538000000000    1.08176000000000    0.24765000000000   Cl   (8e)
   0.41462000000000   -0.08176000000000    0.24765000000000   Cl   (8e)
  -0.41462000000000   -0.58176000000000   -0.25235000000000   Cl   (8e)
   0.91462000000000    0.58176000000000   -0.25235000000000   Cl   (8e)
   0.91462000000000   -0.08176000000000    0.75235000000000   Cl   (8e)
  -0.41462000000000    1.08176000000000    0.75235000000000   Cl   (8e)
   0.01987000000000    0.74957000000000   -0.02314000000000   Cl   (8e)
   0.48013000000000   -0.74957000000000   -0.02314000000000   Cl   (8e)
   0.48013000000000    1.24957000000000    0.52314000000000   Cl   (8e)
   0.01987000000000   -0.24957000000000    0.52314000000000   Cl   (8e)
  -0.01987000000000   -0.74957000000000    0.02314000000000   Cl   (8e)
   0.51987000000000    0.74957000000000    0.02314000000000   Cl   (8e)
   0.51987000000000   -0.24957000000000    0.47686000000000   Cl   (8e)
  -0.01987000000000    1.24957000000000    0.47686000000000   Cl   (8e)
   0.60400000000000    0.75000000000000    0.57900000000000   H    (8e)
  -0.10400000000000   -0.75000000000000    0.57900000000000   H    (8e)
  -0.10400000000000    1.25000000000000   -0.07900000000000   H    (8e)
   0.60400000000000   -0.25000000000000   -0.07900000000000   H    (8e)
  -0.60400000000000   -0.75000000000000   -0.57900000000000   H    (8e)
   1.10400000000000    0.75000000000000   -0.57900000000000   H    (8e)
   1.10400000000000   -0.25000000000000    1.07900000000000   H    (8e)
  -0.60400000000000    1.25000000000000    1.07900000000000   H    (8e)
   0.41200000000000    0.75100000000000    0.31300000000000   H    (8e)
   0.08800000000000   -0.75100000000000    0.31300000000000   H    (8e)
   0.08800000000000    1.25100000000000    0.18700000000000   H    (8e)
   0.41200000000000   -0.25100000000000    0.18700000000000   H    (8e)
  -0.41200000000000   -0.75100000000000   -0.31300000000000   H    (8e)
   0.91200000000000    0.75100000000000   -0.31300000000000   H    (8e)
   0.91200000000000   -0.25100000000000    0.81300000000000   H    (8e)
  -0.41200000000000    1.25100000000000    0.81300000000000   H    (8e)
   0.73700000000000    0.67300000000000    0.54300000000000   H    (8e)
  -0.23700000000000   -0.67300000000000    0.54300000000000   H    (8e)
  -0.23700000000000    1.17300000000000   -0.04300000000000   H    (8e)
   0.73700000000000   -0.17300000000000   -0.04300000000000   H    (8e)
  -0.73700000000000   -0.67300000000000   -0.54300000000000   H    (8e)
   1.23700000000000    0.67300000000000   -0.54300000000000   H    (8e)
   1.23700000000000   -0.17300000000000    1.04300000000000   H    (8e)
  -0.73700000000000    1.17300000000000    1.04300000000000   H    (8e)
   0.25000000000000    0.72200000000000    0.50500000000000   H    (8e)
   0.25000000000000   -0.72200000000000    0.50500000000000   H    (8e)
   0.25000000000000    1.22200000000000   -0.00500000000000   H    (8e)
   0.25000000000000   -0.22200000000000   -0.00500000000000   H    (8e)
  -0.25000000000000   -0.72200000000000   -0.50500000000000   H    (8e)
   0.75000000000000    0.72200000000000   -0.50500000000000   H    (8e)
   0.75000000000000   -0.22200000000000    1.00500000000000   H    (8e)
  -0.25000000000000    1.22200000000000    1.00500000000000   H    (8e)
   0.73400000000000    0.84800000000000    0.58700000000000   H    (8e)
  -0.23400000000000   -0.84800000000000    0.58700000000000   H    (8e)
  -0.23400000000000    1.34800000000000   -0.08700000000000   H    (8e)
   0.73400000000000   -0.34800000000000   -0.08700000000000   H    (8e)
  -0.73400000000000   -0.84800000000000   -0.58700000000000   H    (8e)
   1.23400000000000    0.84800000000000   -0.58700000000000   H    (8e)
   1.23400000000000   -0.34800000000000    1.08700000000000   H    (8e)
  -0.73400000000000    1.34800000000000    1.08700000000000   H    (8e)
   0.30300000000000    0.84500000000000    0.33500000000000   H    (8e)
   0.19700000000000   -0.84500000000000    0.33500000000000   H    (8e)
   0.19700000000000    1.34500000000000    0.16500000000000   H    (8e)
   0.30300000000000   -0.34500000000000    0.16500000000000   H    (8e)
  -0.30300000000000   -0.84500000000000   -0.33500000000000   H    (8e)
   0.80300000000000    0.84500000000000   -0.33500000000000   H    (8e)
   0.80300000000000   -0.34500000000000    0.83500000000000   H    (8e)
  -0.30300000000000    1.34500000000000    0.83500000000000   H    (8e)
   0.51400000000000    0.50900000000000    0.32600000000000   H    (8e)
  -0.01400000000000   -0.50900000000000    0.32600000000000   H    (8e)
  -0.01400000000000    1.00900000000000    0.17400000000000   H    (8e)
   0.51400000000000   -0.00900000000000    0.17400000000000   H    (8e)
  -0.51400000000000   -0.50900000000000   -0.32600000000000   H    (8e)
   1.01400000000000    0.50900000000000   -0.32600000000000   H    (8e)
   1.01400000000000   -0.00900000000000    0.82600000000000   H    (8e)
  -0.51400000000000    1.00900000000000    0.82600000000000   H    (8e)
   0.41700000000000    0.49400000000000    0.60600000000000   H    (8e)
   0.08300000000000   -0.49400000000000    0.60600000000000   H    (8e)
   0.08300000000000    0.99400000000000   -0.10600000000000   H    (8e)
   0.41700000000000    0.00600000000000   -0.10600000000000   H    (8e)
  -0.41700000000000   -0.49400000000000   -0.60600000000000   H    (8e)
   0.91700000000000    0.49400000000000   -0.60600000000000   H    (8e)
   0.91700000000000    0.00600000000000    1.10600000000000   H    (8e)
  -0.41700000000000    0.99400000000000    1.10600000000000   H    (8e)
   0.76800000000000    0.57200000000000    0.34500000000000   H    (8e)
  -0.26800000000000   -0.57200000000000    0.34500000000000   H    (8e)
  -0.26800000000000    1.07200000000000    0.15500000000000   H    (8e)
   0.76800000000000   -0.07200000000000    0.15500000000000   H    (8e)
  -0.76800000000000   -0.57200000000000   -0.34500000000000   H    (8e)
   1.26800000000000    0.57200000000000   -0.34500000000000   H    (8e)
   1.26800000000000   -0.07200000000000    0.84500000000000   H    (8e)
  -0.76800000000000    1.07200000000000    0.84500000000000   H    (8e)
   0.26400000000000   -0.06600000000000    0.68200000000000   H    (8e)
   0.23600000000000    0.06600000000000    0.68200000000000   H    (8e)
   0.23600000000000    0.43400000000000   -0.18200000000000   H    (8e)
   0.26400000000000    0.56600000000000   -0.18200000000000   H    (8e)
  -0.26400000000000    0.06600000000000   -0.68200000000000   H    (8e)
   0.76400000000000   -0.06600000000000   -0.68200000000000   H    (8e)
   0.76400000000000    0.56600000000000    1.18200000000000   H    (8e)
  -0.26400000000000    0.43400000000000    1.18200000000000   H    (8e)
   0.39900000000000   -0.01000000000000    0.56800000000000   H    (8e)
   0.10100000000000    0.01000000000000    0.56800000000000   H    (8e)
   0.10100000000000    0.49000000000000   -0.06800000000000   H    (8e)
   0.39900000000000    0.51000000000000   -0.06800000000000   H    (8e)
  -0.39900000000000    0.01000000000000   -0.56800000000000   H    (8e)
   0.89900000000000   -0.01000000000000   -0.56800000000000   H    (8e)
   0.89900000000000    0.51000000000000    1.06800000000000   H    (8e)
  -0.39900000000000    0.49000000000000    1.06800000000000   H    (8e)
   0.42200000000000   -0.05800000000000    0.50000000000000   H    (8e)
   0.07800000000000    0.05800000000000    0.50000000000000   H    (8e)
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0.07800000000000	0.44200000000000	0.00000000000000	H	(8e)
0.42200000000000	0.55800000000000	0.00000000000000	H	(8e)
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0.92200000000000	−0.05800000000000	−0.50000000000000	H	(8e)
0.92200000000000	0.55800000000000	1.00000000000000	H	(8e)
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0.35900000000000	0.81700000000000	0.67300000000000	H	(8e)
0.14100000000000	−0.81700000000000	0.67300000000000	H	(8e)
0.14100000000000	1.31700000000000	−0.17300000000000	H	(8e)
0.35900000000000	−0.31700000000000	−0.17300000000000	H	(8e)
−0.35900000000000	−0.81700000000000	−0.67300000000000	H	(8e)
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0.85900000000000	−0.31700000000000	1.17300000000000	H	(8e)
−0.35900000000000	1.31700000000000	1.17300000000000	H	(8e)
0.42900000000000	0.81100000000000	0.62300000000000	H	(8e)
0.07100000000000	−0.81100000000000	0.62300000000000	H	(8e)
0.07100000000000	1.31100000000000	−0.12300000000000	H	(8e)
0.42900000000000	−0.31100000000000	−0.12300000000000	H	(8e)
−0.42900000000000	−0.81100000000000	−0.62300000000000	H	(8e)
0.92900000000000	0.81100000000000	−0.62300000000000	H	(8e)
0.92900000000000	−0.31100000000000	1.12300000000000	H	(8e)
−0.42900000000000	1.31100000000000	1.12300000000000	H	(8e)
0.56500000000000	−0.09900000000000	0.67900000000000	H	(8e)
−0.06500000000000	0.09900000000000	0.67900000000000	H	(8e)
−0.06500000000000	0.40100000000000	−0.17900000000000	H	(8e)
0.56500000000000	0.59900000000000	−0.17900000000000	H	(8e)
−0.56500000000000	0.09900000000000	−0.67900000000000	H	(8e)
1.06500000000000	−0.09900000000000	−0.67900000000000	H	(8e)
1.06500000000000	0.59900000000000	1.17900000000000	H	(8e)
−0.56500000000000	0.40100000000000	1.17900000000000	H	(8e)
0.57300000000000	0.60400000000000	0.69500000000000	H	(8e)
−0.07300000000000	−0.60400000000000	0.69500000000000	H	(8e)
−0.07300000000000	1.10400000000000	−0.19500000000000	H	(8e)
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−0.57300000000000	−0.60400000000000	−0.69500000000000	H	(8e)
1.07300000000000	0.60400000000000	−0.69500000000000	H	(8e)
1.07300000000000	−0.10400000000000	1.19500000000000	H	(8e)
−0.57300000000000	1.10400000000000	1.19500000000000	H	(8e)
0.43800000000000	0.66000000000000	0.56100000000000	H	(8e)
0.06200000000000	−0.66000000000000	0.56100000000000	H	(8e)
0.06200000000000	1.16000000000000	−0.06100000000000	H	(8e)
0.43800000000000	−0.16000000000000	−0.06100000000000	H	(8e)
−0.43800000000000	−0.66000000000000	−0.56100000000000	H	(8e)
0.93800000000000	0.66000000000000	−0.56100000000000	H	(8e)
0.93800000000000	−0.16000000000000	1.06100000000000	H	(8e)
−0.43800000000000	1.16000000000000	1.06100000000000	H	(8e)
0.41100000000000	0.61500000000000	0.47500000000000	H	(8e)
0.08900000000000	−0.61500000000000	0.47500000000000	H	(8e)
0.08900000000000	1.11500000000000	0.02500000000000	H	(8e)
0.41100000000000	−0.11500000000000	0.02500000000000	H	(8e)
−0.41100000000000	−0.61500000000000	−0.47500000000000	H	(8e)
0.91100000000000	0.61500000000000	−0.47500000000000	H	(8e)
0.91100000000000	−0.11500000000000	0.97500000000000	H	(8e)
−0.41100000000000	1.11500000000000	0.97500000000000	H	(8e)
0.25000000000000	0.00000000000000	0.25140000000000	K	(4c)
0.25000000000000	0.50000000000000	0.24860000000000	K	(4c)
0.75000000000000	0.00000000000000	−0.25140000000000	K	(4c)
0.75000000000000	0.50000000000000	0.75140000000000	K	(4c)
0.08860000000000	0.65668000000000	0.74995000000000	K	(8e)
0.41140000000000	−0.65668000000000	0.74995000000000	K	(8e)
0.41140000000000	1.15668000000000	−0.24995000000000	K	(8e)
0.08860000000000	−0.15668000000000	−0.24995000000000	K	(8e)
−0.08860000000000	−0.65668000000000	−0.74995000000000	K	(8e)
0.58860000000000	0.65668000000000	−0.74995000000000	K	(8e)
0.58860000000000	−0.15668000000000	1.24995000000000	K	(8e)
−0.08860000000000	1.15668000000000	1.24995000000000	K	(8e)
0.24343000000000	0.25000000000000	0.25000000000000	Mg	(4d)
0.25657000000000	0.75000000000000	0.25000000000000	Mg	(4d)
−0.24343000000000	0.75000000000000	0.75000000000000	Mg	(4d)
0.74343000000000	0.25000000000000	0.75000000000000	Mg	(4d)
0.42094000000000	0.58981000000000	0.74770000000000	Mg	(8e)
0.07906000000000	−0.58981000000000	0.74770000000000	Mg	(8e)
0.07906000000000	1.08981000000000	−0.24770000000000	Mg	(8e)
0.42094000000000	−0.08981000000000	−0.24770000000000	Mg	(8e)
−0.42094000000000	−0.58981000000000	−0.74770000000000	Mg	(8e)
0.92094000000000	0.58981000000000	−0.74770000000000	Mg	(8e)
0.92094000000000	−0.08981000000000	1.24770000000000	Mg	(8e)
−0.42094000000000	1.08981000000000	1.24770000000000	Mg	(8e)
0.36920000000000	0.25000000000000	0.25000000000000	O	(4d)
0.13080000000000	0.75000000000000	0.25000000000000	O	(4d)
−0.36920000000000	0.75000000000000	0.75000000000000	O	(4d)
0.86920000000000	0.25000000000000	0.75000000000000	O	(4d)
0.11620000000000	0.25000000000000	0.25000000000000	O	(4d)
0.38380000000000	0.75000000000000	0.25000000000000	O	(4d)
−0.11620000000000	0.75000000000000	0.75000000000000	O	(4d)
0.61620000000000	0.25000000000000	0.75000000000000	O	(4d)
0.25540000000000	0.70704000000000	0.43840000000000	O	(8e)
0.24460000000000	−0.70704000000000	0.43840000000000	O	(8e)
0.24460000000000	1.20704000000000	0.06160000000000	O	(8e)
0.25540000000000	−0.20704000000000	0.06160000000000	O	(8e)
−0.25540000000000	−0.70704000000000	−0.43840000000000	O	(8e)
0.75540000000000	0.70704000000000	−0.43840000000000	O	(8e)
0.75540000000000	−0.20704000000000	0.93840000000000	O	(8e)
−0.25540000000000	1.20704000000000	0.93840000000000	O	(8e)
0.26730000000000	0.66936000000000	0.15060000000000	O	(8e)
0.23270000000000	−0.66936000000000	0.15060000000000	O	(8e)
0.23270000000000	1.16936000000000	0.34940000000000	O	(8e)
0.26730000000000	−0.16936000000000	0.34940000000000	O	(8e)
−0.26730000000000	−0.66936000000000	−0.15060000000000	O	(8e)
0.76730000000000	0.66936000000000	−0.15060000000000	O	(8e)
0.76730000000000	−0.16936000000000	0.65060000000000	O	(8e)
−0.26730000000000	1.16936000000000	0.65060000000000	O	(8e)
0.44660000000000	0.51041000000000	0.65000000000000	O	(8e)
0.05340000000000	−0.51041000000000	0.65000000000000	O	(8e)
0.05340000000000	1.01041000000000	−0.15000000000000	O	(8e)
0.44660000000000	−0.01041000000000	−0.15000000000000	O	(8e)
−0.44660000000000	−0.51041000000000	−0.65000000000000	O	(8e)
0.94660000000000	0.51041000000000	−0.65000000000000	O	(8e)
0.94660000000000	−0.01041000000000	1.15000000000000	O	(8e)
−0.44660000000000	1.01041000000000	1.15000000000000	O	(8e)
0.29670000000000	0.57117000000000	0.73490000000000	O	(8e)
0.20330000000000	−0.57117000000000	0.73490000000000	O	(8e)
0.20330000000000	1.07117000000000	−0.23490000000000	O	(8e)
0.29670000000000	−0.07117000000000	−0.23490000000000	O	(8e)
−0.29670000000000	−0.57117000000000	−0.73490000000000	O	(8e)
0.79670000000000	0.57117000000000	−0.73490000000000	O	(8e)
0.79670000000000	−0.07117000000000	1.23490000000000	O	(8e)
−0.29670000000000	1.07117000000000	1.23490000000000	O	(8e)
0.42590000000000	0.54662000000000	−0.06460000000000	O	(8e)
0.07410000000000	−0.54662000000000	−0.06460000000000	O	(8e)
0.07410000000000	1.04662000000000	0.56460000000000	O	(8e)
0.42590000000000	−0.04662000000000	0.56460000000000	O	(8e)
−0.42590000000000	−0.54662000000000	0.06460000000000	O	(8e)
0.92590000000000	0.54662000000000	0.06460000000000	O	(8e)
0.92590000000000	−0.04662000000000	0.43540000000000	O	(8e)
−0.42590000000000	1.04662000000000	0.43540000000000	O	(8e)
0.39560000000000	0.66912000000000	0.84440000000000	O	(8e)
0.10440000000000	−0.66912000000000	0.84440000000000	O	(8e)
0.10440000000000	1.16912000000000	−0.34440000000000	O	(8e)
0.39560000000000	−0.16912000000000	−0.34440000000000	O	(8e)
−0.39560000000000	−0.66912000000000	−0.84440000000000	O	(8e)
0.89560000000000	0.66912000000000	−0.84440000000000	O	(8e)
0.89560000000000	−0.16912000000000	1.34440000000000	O	(8e)
−0.39560000000000	1.16912000000000	1.34440000000000	O	(8e)
0.54480000000000	0.60769000000000	0.76310000000000	O	(8e)
−0.04480000000000	−0.60769000000000	0.76310000000000	O	(8e)
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0.54480000000000	−0.10769000000000	−0.26310000000000	O	(8e)
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1.04480000000000	0.60769000000000	−0.76310000000000	O	(8e)
1.04480000000000	−0.10769000000000	1.26310000000000	O	(8e)
−0.54480000000000	1.10769000000000	1.26310000000000	O	(8e)
0.41790000000000	0.63155000000000	0.55800000000000	O	(8e)
0.08210000000000	−0.63155000000000	0.55800000000000	O	(8e)
0.08210000000000	1.13155000000000	−0.05800000000000	O	(8e)
0.41790000000000	−0.13155000000000	−0.05800000000000	O	(8e)
−0.41790000000000	−0.63155000000000	−0.55800000000000	O	(8e)
0.91790000000000	0.63155000000000	−0.55800000000000	O	(8e)
0.91790000000000	−0.13155000000000	1.05800000000000	O	(8e)
−0.41790000000000	1.13155000000000	1.05800000000000	O	(8e)

Eriochalcite ($CuCl_2{\cdot}2H_2O$, *C*45): A2BC4D2_oP18_53_h_a_i_e - CIF

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A redetermination of the crystal structure of cupric chloride dihydrate
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2 x,-y,-z
3 -x+1/2,y,-z+1/2
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x,y,z
7 x+1/2,-y,z+1/2
8 x+1/2,y,-z+1/2
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O1 O 4 e 0.24020 0.00000 0.00000 1.00000
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H1 H 8 i 0.27500 0.06400 0.09900 1.00000
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Eriochalcite ($CuCl_2{\cdot}2H_2O$, *C*45): A2BC4D2_oP18_53_h_a_i_e - POSCAR

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      ↪ 0.099 & Pmna D_{2h}^{7} #53 (aehi) & oP18 & $C45$ & Cl2CuH4O2 &
      ↪  Eriochalcite & S. Brownstein et al., Zeitschrift f{"u}r
      ↪ Kristallographie - Crystalline Materials 189, 13-15 (1989)
   1.00000000000000
   8.08860000000000    0.00000000000000    0.00000000000000
   0.00000000000000    3.74580000000000    0.00000000000000
   0.00000000000000    0.00000000000000    7.41410000000000
    Cl     Cu     H     O
     4      2     8     4
Direct
   0.00000000000000    0.37980000000000    0.23998000000000   Cl    (4h)
   0.50000000000000   -0.37980000000000    0.73998000000000   Cl    (4h)
   0.50000000000000    0.37980000000000    0.26002000000000   Cl    (4h)
   0.00000000000000   -0.37980000000000   -0.23998000000000   Cl    (4h)
   0.00000000000000    0.00000000000000    0.00000000000000   Cu    (2a)
   0.50000000000000    0.00000000000000    0.50000000000000   Cu    (2a)
   0.27500000000000    0.06400000000000    0.09900000000000   H     (8i)
   0.22500000000000   -0.06400000000000    0.59900000000000   H     (8i)
   0.22500000000000    0.06400000000000    0.40100000000000   H     (8i)
   0.27500000000000   -0.06400000000000   -0.09900000000000   H     (8i)
  -0.27500000000000   -0.06400000000000   -0.09900000000000   H     (8i)
   0.77500000000000    0.06400000000000    0.40100000000000   H     (8i)
   0.77500000000000   -0.06400000000000    0.59900000000000   H     (8i)
  -0.27500000000000    0.06400000000000    0.09900000000000   H     (8i)
   0.24020000000000    0.00000000000000    0.00000000000000   O     (4e)
   0.25980000000000    0.00000000000000    0.50000000000000   O     (4e)
  -0.24020000000000    0.00000000000000    0.00000000000000   O     (4e)
   0.74020000000000    0.00000000000000    0.50000000000000   O     (4e)
```

NH_4HF_2 ($F5_8$): A2BC_oP16_53_eh_ab_g - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'F2H5N'
_chemical_formula_sum 'F2 H (NH4)'

loop_
_publ_author_name
 'M. T. Rogers'
 'L. Helmholz'
_journal_name_full_name
;
 Journal of the American Chemical Society
;
_journal_volume 62
_journal_year 1940
_journal_page_first 1533
_journal_page_last 1536
_publ_Section_title
;
 A Redetermination of the Parameters in Ammonium Bifluoride
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'NH$_{4}$HF$_{2}$ ($F5_{8}$) Structure'
_aflow_proto 'A2BC_oP16_53_eh_ab_g'
_aflow_params 'a,b/a,c/a,x_{3},y_{4},y_{5},z_{5}'
_aflow_params_values '8.426,0.437930215998,0.970804652267,0.863,0.46,
      ↪ 0.869,0.371'
_aflow_Strukturbericht '$F5_{8}$'
_aflow_Pearson 'oP16'

_symmetry_space_group_name_H-M "P 2/m 2/n 21/a"
_symmetry_Int_Tables_number 53

_cell_length_a     8.42600
_cell_length_b     3.69000
_cell_length_c     8.18000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x+1/2,y,-z+1/2
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x,y,z
7 x+1/2,-y,z+1/2
8 x+1/2,y,-z+1/2
```

```
loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
H1 H 2 a 0.00000 0.00000 0.00000 1.00000
H2 H 2 b 0.50000 0.00000 0.00000 1.00000
F1 F 4 e 0.86300 0.00000 0.00000 1.00000
NH41 NH4 4 g 0.25000 0.46000 0.25000 1.00000
F2 F 4 h 0.00000 0.86900 0.37100 1.00000
```

NH_4HF_2 ($F5_8$): A2BC_oP16_53_eh_ab_g - POSCAR

```
A2BC_oP16_53_eh_ab_g & a,b/a,c/a,x3,y4,y5,z5 --params=8.426,
      ↪ 0.437930215998,0.970804652267,0.863,0.46,0.869,0.371 & Pmna D_{
      ↪ 2h}^{7} #53 (abegh) & oP16 & $F5_{8}$ & F2H5N & F2H5N & M. T.
      ↪ Rogers and L. Helmholz, J. Am. Chem. Soc. 62, 1533-1536 (1940)
   1.00000000000000
   8.42600000000000    0.00000000000000    0.00000000000000
   0.00000000000000    3.69000000000000    0.00000000000000
   0.00000000000000    0.00000000000000    8.18000000000000
     F     H   NH4
     8     4     4
Direct
   0.86300000000000    0.00000000000000    0.00000000000000   F    (4e)
  -0.36300000000000    0.00000000000000    0.50000000000000   F    (4e)
  -0.86300000000000    0.00000000000000    0.00000000000000   F    (4e)
   1.36300000000000    0.00000000000000    0.50000000000000   F    (4e)
   0.00000000000000    0.86900000000000    0.37100000000000   F    (4h)
   0.50000000000000   -0.86900000000000    0.87100000000000   F    (4h)
   0.50000000000000    0.86900000000000    0.12900000000000   F    (4h)
   0.00000000000000   -0.86900000000000   -0.37100000000000   F    (4h)
   0.00000000000000    0.00000000000000    0.00000000000000   H    (2a)
   0.50000000000000    0.00000000000000    0.50000000000000   H    (2a)
   0.50000000000000    0.00000000000000    0.00000000000000   H    (2b)
   0.00000000000000    0.00000000000000    0.50000000000000   H    (2b)
   0.25000000000000    0.46000000000000    0.25000000000000   NH4  (4g)
   0.25000000000000   -0.46000000000000    0.75000000000000   NH4  (4g)
   0.75000000000000   -0.46000000000000    0.75000000000000   NH4  (4g)
   0.75000000000000    0.46000000000000    0.25000000000000   NH4  (4g)
```

Orthorhombic $Sr_4Ru_3O_{10}$: A10B3C4_oP68_55_2e2fgh2i_adef_2e2f - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'O10Ru3Sr4'
_chemical_formula_sum 'O10 Ru3 Sr4'

loop_
_publ_author_name
 'M. K. Crawford'
 'R. L. Harlow'
 'W. Marshall'
 'Z. Li'
 'G. Cao'
 'R. L. Lindstrom'
 'Q. Huang'
 'J. W. Lynn'
_journal_name_full_name
;
 Physical Review B
;
_journal_volume 65
_journal_year 2002
_journal_page_first 214412
_journal_page_last 214412
_publ_Section_title
;
 Structure and magnetism of single crystal Sr$_{4}$Ru$_{3}$O$_{10}$: A
      ↪ ferromagnetic triple-layer ruthenate
;

_aflow_title 'Orthorhombic Sr$_{4}$Ru$_{3}$O$_{10}$ Structure'
_aflow_proto 'A10B3C4_oP68_55_2e2fgh2i_adef_2e2f'
_aflow_params 'a,b/a,c/a,z_{3},z_{4},z_{5},z_{6},z_{7},z_{8},z_{9},z_{10
      ↪ },z_{11},z_{12},x_{13},y_{13},x_{14},y_{14},x_{15},y_{15},z_{15
      ↪ },x_{16},y_{16},z_{16}'
_aflow_params_values '3.9001,1.0,7.32622240455,0.0695,0.213,0.1402,
      ↪ 0.2961,0.4301,0.2871,0.4303,0.3598,0.0699,0.2038,0.2028,0.2971,
      ↪ 0.2966,0.2964,0.2721,0.2271,0.1392,0.2266,0.2275,0.3608'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP68'

_symmetry_space_group_name_H-M "P 21/b 21/a 2/m"
_symmetry_Int_Tables_number 55

_cell_length_a     3.90010
_cell_length_b     3.90010
_cell_length_c     28.57300
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z
3 -x+1/2,y+1/2,-z
```

```
4 -x,-y,z
5 -x,-y,-z
6 -x+1/2,y+1/2,z
7 x+1/2,-y+1/2,z
8 x,y,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ru1 Ru 2 a 0.00000 0.00000 0.00000 1.00000
Ru2 Ru 2 d 0.00000 0.50000 0.50000 1.00000
O1 O 4 e 0.00000 0.00000 0.06950 1.00000
O2 O 4 e 0.00000 0.00000 0.21300 1.00000
Ru3 Ru 4 e 0.00000 0.00000 0.14020 1.00000
Sr1 Sr 4 e 0.00000 0.00000 0.29610 1.00000
Sr2 Sr 4 e 0.00000 0.00000 0.43010 1.00000
O3 O 4 f 0.00000 0.50000 0.28710 1.00000
O4 O 4 f 0.00000 0.50000 0.43030 1.00000
Ru4 Ru 4 f 0.00000 0.50000 0.35980 1.00000
Sr3 Sr 4 f 0.00000 0.50000 0.06990 1.00000
Sr4 Sr 4 f 0.00000 0.50000 0.20380 1.00000
O5 O 4 g 0.20280 0.29710 0.00000 1.00000
O6 O 4 h 0.29660 0.29640 0.50000 1.00000
O7 O 8 i 0.27210 0.22710 0.13920 1.00000
O8 O 8 i 0.22660 0.22750 0.36080 1.00000
```

Orthorhombic $Sr_4Ru_3O_{10}$: A10B3C4_oP68_55_2e2fgh2i_adef_2e2f - POSCAR

```
A10B3C4_oP68_55_2e2fgh2i_adef_2e2f & a,b/a,c/a,z3,z4,z5,z6,z7,z8,z9,z10,
    ↪ z11,z12,x13,y13,x14,y14,x15,y15,z15,x16,y16,z16 --params=3.9001
    ↪ ,1.0,7.32622240455,0.0695,0.213,0.1402,0.2961,0.4301,0.2871,
    ↪ 0.4303,0.3598,0.0699,0.2038,0.2028,0.2971,0.2966,0.2964,0.2721,
    ↪ 0.2271,0.1392,0.2266,0.2275,0.3608 & Pbam D_{2h}^{9} #55 (ade^
    ↪ 5f^5ghi^2) & oP68 & None & O10Ru3Sr4 & O10Ru3Sr4 & M. K.
    ↪ Crawford et al., Phys. Rev. B 65, 214412(2002)
   1.00000000000000
   3.90010000000000   0.00000000000000   0.00000000000000
   0.00000000000000   3.90010000000000   0.00000000000000
   0.00000000000000   0.00000000000000  28.57300000000000
     O    Ru    Sr
    40    12    16
Direct
   0.00000000000000   0.00000000000000   0.06950000000000   O   (4e)
   0.50000000000000   0.50000000000000  -0.06950000000000   O   (4e)
   0.00000000000000   0.00000000000000  -0.06950000000000   O   (4e)
   0.50000000000000   0.50000000000000   0.06950000000000   O   (4e)
   0.00000000000000   0.00000000000000   0.21300000000000   O   (4e)
   0.50000000000000   0.50000000000000  -0.21300000000000   O   (4e)
   0.00000000000000   0.00000000000000  -0.21300000000000   O   (4e)
   0.50000000000000   0.50000000000000   0.21300000000000   O   (4e)
   0.00000000000000   0.50000000000000   0.28710000000000   O   (4f)
   0.50000000000000   0.00000000000000  -0.28710000000000   O   (4f)
   0.00000000000000   0.50000000000000  -0.28710000000000   O   (4f)
   0.50000000000000   0.00000000000000   0.28710000000000   O   (4f)
   0.00000000000000   0.50000000000000   0.43030000000000   O   (4f)
   0.50000000000000   0.00000000000000  -0.43030000000000   O   (4f)
   0.00000000000000   0.50000000000000  -0.43030000000000   O   (4f)
   0.50000000000000   0.00000000000000   0.43030000000000   O   (4f)
   0.20280000000000   0.29710000000000   0.00000000000000   O   (4g)
  -0.20280000000000  -0.29710000000000   0.00000000000000   O   (4g)
   0.29720000000000   0.79710000000000   0.00000000000000   O   (4g)
   0.70280000000000   0.20290000000000   0.00000000000000   O   (4g)
   0.29660000000000   0.29640000000000   0.50000000000000   O   (4h)
  -0.29660000000000  -0.29640000000000   0.50000000000000   O   (4h)
   0.20340000000000   0.79640000000000   0.50000000000000   O   (4h)
   0.79660000000000   0.20360000000000   0.50000000000000   O   (4h)
   0.27210000000000   0.22710000000000   0.13920000000000   O   (8i)
  -0.27210000000000  -0.22710000000000   0.13920000000000   O   (8i)
   0.22790000000000   0.72710000000000  -0.13920000000000   O   (8i)
   0.77210000000000   0.27290000000000  -0.13920000000000   O   (8i)
  -0.27210000000000  -0.22710000000000  -0.13920000000000   O   (8i)
   0.27210000000000   0.22710000000000  -0.13920000000000   O   (8i)
   0.77210000000000   0.27290000000000   0.13920000000000   O   (8i)
   0.22790000000000   0.72710000000000   0.13920000000000   O   (8i)
   0.22660000000000   0.22750000000000   0.36080000000000   O   (8i)
  -0.22660000000000  -0.22750000000000   0.36080000000000   O   (8i)
   0.27340000000000   0.72750000000000  -0.36080000000000   O   (8i)
   0.72660000000000   0.27250000000000  -0.36080000000000   O   (8i)
  -0.22660000000000  -0.22750000000000  -0.36080000000000   O   (8i)
   0.22660000000000   0.22750000000000  -0.36080000000000   O   (8i)
   0.72660000000000   0.27250000000000   0.36080000000000   O   (8i)
   0.27340000000000   0.72750000000000   0.36080000000000   O   (8i)
   0.00000000000000   0.00000000000000   0.00000000000000   Ru  (2a)
   0.50000000000000   0.50000000000000   0.00000000000000   Ru  (2a)
   0.00000000000000   0.50000000000000   0.50000000000000   Ru  (2d)
   0.50000000000000   0.00000000000000   0.50000000000000   Ru  (2d)
   0.00000000000000   0.00000000000000   0.14020000000000   Ru  (4e)
   0.50000000000000   0.50000000000000  -0.14020000000000   Ru  (4e)
   0.00000000000000   0.00000000000000  -0.14020000000000   Ru  (4e)
   0.50000000000000   0.50000000000000   0.14020000000000   Ru  (4e)
   0.00000000000000   0.50000000000000   0.35980000000000   Ru  (4f)
   0.50000000000000   0.00000000000000  -0.35980000000000   Ru  (4f)
   0.00000000000000   0.50000000000000  -0.35980000000000   Ru  (4f)
   0.50000000000000   0.00000000000000   0.35980000000000   Ru  (4f)
   0.00000000000000   0.00000000000000   0.29610000000000   Sr  (4e)
   0.50000000000000   0.50000000000000  -0.29610000000000   Sr  (4e)
   0.00000000000000   0.00000000000000  -0.29610000000000   Sr  (4e)
   0.50000000000000   0.50000000000000   0.29610000000000   Sr  (4e)
   0.00000000000000   0.00000000000000   0.43010000000000   Sr  (4e)
   0.50000000000000   0.50000000000000  -0.43010000000000   Sr  (4e)
   0.00000000000000   0.00000000000000  -0.43010000000000   Sr  (4e)
   0.50000000000000   0.50000000000000   0.43010000000000   Sr  (4e)
   0.00000000000000   0.50000000000000   0.06990000000000   Sr  (4f)
   0.50000000000000   0.00000000000000  -0.06990000000000   Sr  (4f)
   0.00000000000000   0.50000000000000  -0.06990000000000   Sr  (4f)
   0.50000000000000   0.00000000000000   0.06990000000000   Sr  (4f)
   0.00000000000000   0.50000000000000   0.20380000000000   Sr  (4f)
   0.50000000000000   0.00000000000000  -0.20380000000000   Sr  (4f)
   0.00000000000000   0.50000000000000  -0.20380000000000   Sr  (4f)
   0.50000000000000   0.00000000000000   0.20380000000000   Sr  (4f)
```

$Nb_2Pd_3Se_8$: A2B3C8_oP26_55_h_ag_2g2h - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Nb2Pd3Se8'
_chemical_formula_sum 'Nb2 Pd3 Se8'

loop_
_publ_author_name
 'D. A. Keszler'
 'J. A. Ibers'
_journal_name_full_name
;
 Journal of Solid State Chemistry
;
_journal_volume 52
_journal_year 1984
_journal_page_first 73
_journal_page_last 79
_publ_Section_title
;
 A new structural type in ternary chalcogenide chemistry: Structure and
    ↪ properties of Nb$_{2}$Pd$_{3}$Se$_{8}$
;

_aflow_title 'Nb$_{2}$Pd$_{3}$Se$_{8}$ Structure'
_aflow_proto 'A2B3C8_oP26_55_h_ag_2g2h'
_aflow_params 'a,b/a,c/a,x_{2},y_{2},x_{3},y_{3},x_{4},y_{4},x_{5},y_{5
    ↪ },x_{6},y_{6},x_{7},y_{7}'
_aflow_params_values '15.074,0.701406395117,0.235305824599,0.21616,
    ↪ 0.3812,-0.01109,0.23185,0.15734,0.04391,0.11591,0.21532,0.28335
    ↪ ,0.25032,0.11562,0.45742'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP26'

_symmetry_space_group_name_H-M "P 21/b 21/a 2/m"
_symmetry_Int_Tables_number 55

_cell_length_a    15.07400
_cell_length_b    10.57300
_cell_length_c    3.54700
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z
3 -x+1/2,y+1/2,-z
4 -x,-y,z
5 -x,-y,-z
6 -x+1/2,y+1/2,z
7 x+1/2,-y+1/2,z
8 x,y,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Pd1 Pd 2 a 0.00000 0.00000 0.00000 1.00000
Pd2 Pd 4 g 0.21616 0.38120 0.00000 1.00000
Se1 Se 4 g -0.01109 0.23185 0.00000 1.00000
Se2 Se 4 g 0.15734 0.04391 0.00000 1.00000
Nb1 Nb 4 h 0.11591 0.21532 0.50000 1.00000
Se3 Se 4 h 0.28335 0.25032 0.50000 1.00000
Se4 Se 4 h 0.11562 0.45742 0.50000 1.00000
```

$Nb_2Pd_3Se_8$: A2B3C8_oP26_55_h_ag_2g2h - POSCAR

```
A2B3C8_oP26_55_h_ag_2g2h & a,b/a,c/a,x2,y2,x3,y3,x4,y4,x5,y5,x6,y6,x7,y7
    ↪ --params=15.074,0.701406395117,0.235305824599,0.21616,0.3812,-
    ↪ 0.01109,0.23185,0.15734,0.04391,0.11591,0.21532,0.28335,0.25032
    ↪ ,0.11562,0.45742 & Pbam D_{2h}^{9} #55 (ag^3h^3) & oP26 & None
    ↪ & Nb2Pd3Se8 & Nb2Pd3Se8 & D. A. Keszler and J. A. Ibers, J.
    ↪ Solid State Chem. 52, 73-79 (1984)
   1.00000000000000
  15.07400000000000   0.00000000000000   0.00000000000000
   0.00000000000000  10.57300000000000   0.00000000000000
   0.00000000000000   0.00000000000000   3.54700000000000
     Nb    Pd    Se
      4     6    16
Direct
   0.11591000000000   0.21532000000000   0.50000000000000   Nb   (4h)
  -0.11591000000000  -0.21532000000000   0.50000000000000   Nb   (4h)
```

```
  0.38409000000000   0.71532000000000   0.50000000000000   Nb   (4h)
  0.61591000000000   0.28468000000000   0.50000000000000   Nb   (4h)
  0.00000000000000   0.00000000000000   0.00000000000000   Pd   (2a)
  0.50000000000000   0.50000000000000   0.00000000000000   Pd   (2a)
  0.21616000000000   0.38120000000000   0.00000000000000   Pd   (4g)
 -0.21616000000000  -0.38120000000000   0.00000000000000   Pd   (4g)
  0.28384000000000   0.88120000000000   0.00000000000000   Pd   (4g)
  0.71616000000000   0.11880000000000   0.00000000000000   Pd   (4g)
 -0.01109000000000   0.23185000000000   0.00000000000000   Se   (4g)
  0.01109000000000  -0.23185000000000   0.00000000000000   Se   (4g)
  0.51109000000000   0.73185000000000   0.00000000000000   Se   (4g)
  0.48891000000000   0.26815000000000   0.00000000000000   Se   (4g)
  0.15734000000000   0.04391000000000   0.00000000000000   Se   (4g)
 -0.15734000000000  -0.04391000000000   0.00000000000000   Se   (4g)
  0.34266000000000   0.54391000000000   0.00000000000000   Se   (4g)
  0.65734000000000   0.45609000000000   0.00000000000000   Se   (4g)
  0.28335000000000   0.25032000000000   0.50000000000000   Se   (4h)
 -0.28335000000000  -0.25032000000000   0.50000000000000   Se   (4h)
  0.21665000000000   0.75032000000000   0.50000000000000   Se   (4h)
  0.78335000000000   0.24968000000000   0.50000000000000   Se   (4h)
  0.11562000000000   0.45742000000000   0.50000000000000   Se   (4h)
 -0.11562000000000  -0.45742000000000   0.50000000000000   Se   (4h)
  0.38438000000000   0.95742000000000   0.50000000000000   Se   (4h)
  0.61562000000000   0.04258000000000   0.50000000000000   Se   (4h)
```

$K_2HgCl_4{\cdot}H_2O$ ($E3_4$): A4BCD2_oP32_55_ghi_f_e_gh - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Cl4(H2O)HgK2'
_chemical_formula_sum 'Cl4 (H2O) Hg K2'

loop_
_publ_author_name
 'K. Aurivillius'
 'C. St{\aa}lhandske'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 27
_journal_year 1973
_journal_page_first 1086
_journal_page_last 1088
_publ_Section_title
;
 An X-Ray Single Crystal Study of K$_{2}$HgCl$_{4}$$\cdot$H$_{2}$O
;

_aflow_title 'K$_{2}$HgCl$_{4}$$\cdot$H$_{2}$O ($E3_{4}$) Structure'
_aflow_proto 'A4BCD2_oP32_55_ghi_f_e_gh'
_aflow_params 'a,b/a,c/a,z_{1},z_{2},x_{3},y_{3},x_{4},y_{4},x_{5},y_{5
      ↪ },x_{6},y_{6},x_{7},y_{7},z_{7}'
_aflow_params_values '8.258,1.41220634536,1.08077016227,0.22925,0.2312,
      ↪ 0.2042,0.0763,0.0819,0.341,0.2489,0.0601,0.1047,0.3047,0.8831,
      ↪ 0.186,0.252'
_aflow_Strukturbericht '$E3_{4}$'
_aflow_Pearson 'oP32'

_symmetry_space_group_name_H-M "P 21/b 21/a 2/m"
_symmetry_Int_Tables_number 55

_cell_length_a    8.25800
_cell_length_b    11.66200
_cell_length_c    8.92500
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z
3 -x+1/2,y+1/2,-z
4 -x,-y,z
5 -x,-y,-z
6 -x+1/2,y+1/2,z
7 x+1/2,-y+1/2,z
8 x,y,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Hg1 Hg 4 e 0.00000 0.00000 0.22925 1.00000
H2O1 H2O 4 f 0.00000 0.50000 0.23120 1.00000
Cl1 Cl 4 g 0.20420 0.07630 0.00000 1.00000
K1 K 4 g 0.08190 0.34100 0.00000 1.00000
Cl2 Cl 4 h 0.24890 0.06010 0.50000 1.00000
K2 K 4 h 0.10470 0.30470 0.50000 1.00000
Cl3 Cl 8 i 0.88310 0.18600 0.25200 1.00000
```

$K_2HgCl_4{\cdot}H_2O$ ($E3_4$): A4BCD2_oP32_55_ghi_f_e_gh - POSCAR

```
A4BCD2_oP32_55_ghi_f_e_gh & a,b/a,c/a,z1,z2,x3,y3,x4,y4,x5,y5,x6,y6,x7,
      ↪ y7,z7 --params=8.258,1.41220634536,1.08077016227,0.22925,0.2312
      ↪ ,0.2042,0.0763,0.0819,0.341,0.2489,0.0601,0.1047,0.3047,0.8831,
      ↪ 0.186,0.252 & Pbam D_{2h}^{9} #55 (efg^2h^2i) & oP32 & $E3_{4}$
      ↪ & Cl4(H2O)HgK2 & Cl4(H2O)HgK2 & K. Aurivillius and C. St{\aa}
      ↪ lhandske, Acta Chem. Scand. 27, 1086-1088 (1973)
   1.00000000000000
   8.25800000000000   0.00000000000000   0.00000000000000
   0.00000000000000  11.66200000000000   0.00000000000000
   0.00000000000000   0.00000000000000   8.92500000000000
     Cl    H2O     Hg      K
     16      4      4      8
Direct
   0.20420000000000   0.07630000000000   0.00000000000000   Cl   (4g)
  -0.20420000000000  -0.07630000000000   0.00000000000000   Cl   (4g)
   0.29580000000000   0.57630000000000   0.00000000000000   Cl   (4g)
   0.70420000000000   0.42370000000000   0.00000000000000   Cl   (4g)
   0.24890000000000   0.06010000000000   0.50000000000000   Cl   (4h)
  -0.24890000000000  -0.06010000000000   0.50000000000000   Cl   (4h)
   0.25110000000000   0.56010000000000   0.50000000000000   Cl   (4h)
   0.74890000000000   0.43990000000000   0.50000000000000   Cl   (4h)
   0.88310000000000   0.18600000000000   0.25200000000000   Cl   (8i)
  -0.88310000000000  -0.18600000000000   0.25200000000000   Cl   (8i)
  -0.38310000000000   0.68600000000000  -0.25200000000000   Cl   (8i)
   1.38310000000000   0.31400000000000  -0.25200000000000   Cl   (8i)
  -0.88310000000000  -0.18600000000000  -0.25200000000000   Cl   (8i)
   0.88310000000000   0.18600000000000  -0.25200000000000   Cl   (8i)
   1.38310000000000   0.31400000000000   0.25200000000000   Cl   (8i)
  -0.38310000000000   0.68600000000000   0.25200000000000   Cl   (8i)
   0.00000000000000   0.50000000000000   0.23120000000000   H2O  (4f)
   0.50000000000000   0.00000000000000  -0.23120000000000   H2O  (4f)
   0.00000000000000   0.50000000000000  -0.23120000000000   H2O  (4f)
   0.50000000000000   0.00000000000000   0.23120000000000   H2O  (4f)
   0.00000000000000   0.00000000000000   0.22925000000000   Hg   (4e)
   0.50000000000000   0.50000000000000  -0.22925000000000   Hg   (4e)
   0.00000000000000   0.00000000000000  -0.22925000000000   Hg   (4e)
   0.50000000000000   0.50000000000000   0.22925000000000   Hg   (4e)
   0.08190000000000   0.34100000000000   0.00000000000000   K    (4g)
  -0.08190000000000  -0.34100000000000   0.00000000000000   K    (4g)
   0.41810000000000   0.84100000000000   0.00000000000000   K    (4g)
   0.58190000000000   0.15900000000000   0.00000000000000   K    (4g)
   0.10470000000000   0.30470000000000   0.50000000000000   K    (4h)
  -0.10470000000000  -0.30470000000000   0.50000000000000   K    (4h)
   0.39530000000000   0.80470000000000   0.50000000000000   K    (4h)
   0.60470000000000   0.19530000000000   0.50000000000000   K    (4h)
```

$Ru_{11}B_8$: A8B11_oP38_55_g3h_a3g2h - CIF

```
# CIF file
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_audit_creation_method FINDSYM

_chemical_name_mineral 'B8Ru11'
_chemical_formula_sum 'B8 Ru11'

loop_
_publ_author_name
 'J. {\AA}selius'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 14
_journal_year 1960
_journal_page_first 2169
_journal_page_last 2176
_publ_Section_title
;
 The Crystal Structure of Ru$_{11}$B$_{8}$
;

_aflow_title 'Ru$_{11}$B$_{8}$ Structure'
_aflow_proto 'A8B11_oP38_55_g3h_a3g2h'
_aflow_params 'a,b/a,c/a,x_{2},y_{2},x_{3},y_{3},x_{4},y_{4},x_{5},y_{5
      ↪ },x_{6},y_{6},x_{7},y_{7},x_{8},y_{8},x_{9},y_{9},x_{10},y_{10
      ↪ }'
_aflow_params_values '11.609,0.97700060298,0.244293220777,0.34,0.21,
      ↪ 0.2844,0.3913,0.0429,0.3952,0.1686,0.174,0.13,0.01,0.15,0.32,
      ↪ 0.27,0.25,0.4636,0.2962,0.3404,0.0616'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP38'

_symmetry_space_group_name_H-M "P 21/b 21/a 2/m"
_symmetry_Int_Tables_number 55

_cell_length_a    11.60900
_cell_length_b    11.34200
_cell_length_c    2.83600
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z
3 -x+1/2,y+1/2,-z
4 -x,-y,z
5 -x,-y,-z
6 -x+1/2,y+1/2,z
7 x+1/2,-y+1/2,z
8 x,y,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
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```
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ru1 Ru 2 a 0.00000 0.00000 0.00000 1.00000
B1 B 4 g 0.34000 0.21000 0.00000 1.00000
Ru2 Ru 4 g 0.28440 0.39130 0.00000 1.00000
Ru3 Ru 4 g 0.04290 0.39520 0.00000 1.00000
Ru4 Ru 4 g 0.16860 0.17400 0.00000 1.00000
B2 B 4 h 0.13000 0.01000 0.50000 1.00000
B3 B 4 h 0.15000 0.32000 0.50000 1.00000
B4 B 4 h 0.27000 0.25000 0.50000 1.00000
Ru5 Ru 4 h 0.46360 0.29620 0.50000 1.00000
Ru6 Ru 4 h 0.34040 0.06160 0.50000 1.00000
```

$Ru_{11}B_8$: A8B11_oP38_55_g3h_a3g2h - POSCAR

```
A8B11_oP38_55_g3h_a3g2h & a,b/a,c/a,x2,y2,x3,y3,x4,y4,x5,y5,x6,y6,x7,y7,
    ↪ x8,y8,x9,y9,x10,y10 --params=11.609,0.97700060298,
    ↪ 0.244293220777,0.34,0.21,0.2844,0.3913,0.0429,0.3952,0.1686,
    ↪ 0.174,0.13,0.01,0.15,0.32,0.27,0.25,0.4636,0.2962,0.3404,0.0616
    ↪ & Pbam D_{2h}^{9} #55 (ag^4h^5) & oP38 & None & B8Ru11 &
    ↪ B8Ru11 & J. {\AA}selius, Acta Chem. Scand. 14, 2169-2176 (1960)
  1.00000000000000
 11.60900000000000   0.00000000000000   0.00000000000000
  0.00000000000000  11.34200000000000   0.00000000000000
  0.00000000000000   0.00000000000000   2.83600000000000
    B    Ru
   16    22
Direct
   0.34000000000000   0.21000000000000   0.00000000000000   B   (4g)
  -0.34000000000000  -0.21000000000000   0.00000000000000   B   (4g)
   0.16000000000000   0.71000000000000   0.00000000000000   B   (4g)
   0.84000000000000   0.29000000000000   0.00000000000000   B   (4g)
   0.13000000000000   0.01000000000000   0.50000000000000   B   (4h)
  -0.13000000000000  -0.01000000000000   0.50000000000000   B   (4h)
   0.37000000000000   0.51000000000000   0.50000000000000   B   (4h)
   0.63000000000000   0.49000000000000   0.50000000000000   B   (4h)
   0.15000000000000   0.32000000000000   0.50000000000000   B   (4h)
  -0.15000000000000  -0.32000000000000   0.50000000000000   B   (4h)
   0.35000000000000   0.82000000000000   0.50000000000000   B   (4h)
   0.65000000000000   0.18000000000000   0.50000000000000   B   (4h)
   0.27000000000000   0.25000000000000   0.50000000000000   B   (4h)
  -0.27000000000000  -0.25000000000000   0.50000000000000   B   (4h)
   0.23000000000000   0.75000000000000   0.50000000000000   B   (4h)
   0.77000000000000   0.25000000000000   0.50000000000000   B   (4h)
   0.00000000000000   0.00000000000000   0.00000000000000   Ru  (2a)
   0.50000000000000   0.50000000000000   0.00000000000000   Ru  (2a)
   0.28440000000000   0.39130000000000   0.00000000000000   Ru  (4g)
  -0.28440000000000  -0.39130000000000   0.00000000000000   Ru  (4g)
   0.21560000000000   0.89130000000000   0.00000000000000   Ru  (4g)
   0.78440000000000   0.10870000000000   0.00000000000000   Ru  (4g)
   0.04290000000000   0.39520000000000   0.00000000000000   Ru  (4g)
  -0.04290000000000  -0.39520000000000   0.00000000000000   Ru  (4g)
   0.45710000000000   0.89520000000000   0.00000000000000   Ru  (4g)
   0.54290000000000   0.10480000000000   0.00000000000000   Ru  (4g)
   0.16860000000000   0.17400000000000   0.00000000000000   Ru  (4g)
  -0.16860000000000  -0.17400000000000   0.00000000000000   Ru  (4g)
   0.33140000000000   0.67400000000000   0.00000000000000   Ru  (4g)
   0.66860000000000   0.32600000000000   0.00000000000000   Ru  (4g)
   0.46360000000000   0.29620000000000   0.50000000000000   Ru  (4h)
  -0.46360000000000  -0.29620000000000   0.50000000000000   Ru  (4h)
   0.03640000000000   0.79620000000000   0.50000000000000   Ru  (4h)
   0.96360000000000   0.20380000000000   0.50000000000000   Ru  (4h)
   0.34040000000000   0.06160000000000   0.50000000000000   Ru  (4h)
  -0.34040000000000  -0.06160000000000   0.50000000000000   Ru  (4h)
   0.15960000000000   0.56160000000000   0.50000000000000   Ru  (4h)
   0.84040000000000   0.43840000000000   0.50000000000000   Ru  (4h)
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$HoMn_2O_5$: AB2C5_oP32_55_g_fh_eghi - CIF

```
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_audit_creation_method FINDSYM

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_chemical_formula_sum 'Ho Mn2 O5'

loop_
_publ_author_name
 'S. {Quezel-Ambrunaz}'
 'F. Bertaut'
 'G. Buisson'
_journal_name_full_name
;
 Comptes rendus de l'{A}cad\'{e}mie des Sciences
;
_journal_volume 258
_journal_year 1964
_journal_page_first 3025
_journal_page_last 3027
_publ_Section_title
;
 Structure des compos\\'{e}s d\'oxydes de terres rares et de mangan\'{e}
    ↪ se de formule TMn$_{2}$O$_{5}$
;

# Found in Structure of NdMn$_{2}$O$_{5}$, 1993

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_aflow_proto 'AB2C5_oP32_55_g_fh_eghi'
_aflow_params 'a,b/a,c/a,z_{1},z_{2},x_{3},y_{3},x_{4},y_{4},x_{5},y_{5
    ↪ },x_{6},y_{6},x_{7},y_{7},z_{7}'
_aflow_params_values '7.36,1.1535326087,0.773097826087,0.25,0.25,0.143,
    ↪ 0.172,0.14,0.44,0.09,0.848,0.14,0.44,0.1,0.72,0.25'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP32'

_symmetry_space_group_name_H-M "P 21/b 21/a 2/m"
_symmetry_Int_Tables_number 55

_cell_length_a    7.36000
_cell_length_b    8.49000
_cell_length_c    5.69000
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z
3 -x+1/2,y+1/2,-z
4 -x,-y,z
5 -x,-y,-z
6 -x+1/2,y+1/2,z
7 x+1/2,-y+1/2,z
8 x,y,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 4 e 0.00000 0.00000 0.25000 1.00000
Mn1 Mn 4 f 0.00000 0.50000 0.25000 1.00000
Ho1 Ho 4 g 0.14300 0.17200 0.00000 1.00000
O2 O 4 g 0.14000 0.44000 0.00000 1.00000
Mn2 Mn 4 h 0.09000 0.84800 0.50000 1.00000
O3 O 4 h 0.14000 0.44000 0.50000 1.00000
O4 O 8 i 0.10000 0.72000 0.25000 1.00000
```

$HoMn_2O_5$: AB2C5_oP32_55_g_fh_eghi - POSCAR

```
AB2C5_oP32_55_g_fh_eghi & a,b/a,c/a,z1,z2,x3,y3,x4,y4,x5,y5,x6,y6,x7,y7,
    ↪ z7 --params=7.36,1.1535326087,0.773097826087,0.25,0.25,0.143,
    ↪ 0.172,0.14,0.44,0.09,0.848,0.14,0.44,0.1,0.72,0.25 & Pbam D_{2h
    ↪ }^{9} #55 (efg^2h^2i) & oP32 & None & HoMn2O5 & HoMn2O5 & S. {
    ↪ Quezel-Ambrunaz} and F. Bertaut and G. Buisson, {C. R. Acad.
    ↪ Sci. 258, 3025-3027 (1964)
  1.00000000000000
  7.36000000000000   0.00000000000000   0.00000000000000
  0.00000000000000   8.49000000000000   0.00000000000000
  0.00000000000000   0.00000000000000   5.69000000000000
   Ho    Mn    O
    4     8    20
Direct
   0.14300000000000   0.17200000000000   0.00000000000000   Ho   (4g)
  -0.14300000000000  -0.17200000000000   0.00000000000000   Ho   (4g)
   0.35700000000000   0.67200000000000   0.00000000000000   Ho   (4g)
   0.64300000000000   0.32800000000000   0.00000000000000   Ho   (4g)
   0.00000000000000   0.50000000000000   0.25000000000000   Mn   (4f)
   0.50000000000000   0.00000000000000  -0.25000000000000   Mn   (4f)
   0.00000000000000   0.50000000000000  -0.25000000000000   Mn   (4f)
   0.50000000000000   0.00000000000000   0.25000000000000   Mn   (4f)
   0.09000000000000   0.84800000000000   0.50000000000000   Mn   (4h)
  -0.09000000000000  -0.84800000000000   0.50000000000000   Mn   (4h)
   0.41000000000000   1.34800000000000   0.50000000000000   Mn   (4h)
   0.59000000000000  -0.34800000000000   0.50000000000000   Mn   (4h)
   0.00000000000000   0.00000000000000   0.25000000000000   O    (4e)
   0.50000000000000   0.50000000000000  -0.25000000000000   O    (4e)
   0.00000000000000   0.00000000000000  -0.25000000000000   O    (4e)
   0.50000000000000   0.50000000000000   0.25000000000000   O    (4e)
   0.14000000000000   0.44000000000000   0.00000000000000   O    (4g)
  -0.14000000000000  -0.44000000000000   0.00000000000000   O    (4g)
   0.36000000000000   0.94000000000000   0.00000000000000   O    (4g)
   0.64000000000000   0.06000000000000   0.00000000000000   O    (4g)
   0.14000000000000   0.44000000000000   0.50000000000000   O    (4h)
  -0.14000000000000  -0.44000000000000   0.50000000000000   O    (4h)
   0.36000000000000   0.94000000000000   0.50000000000000   O    (4h)
   0.64000000000000   0.06000000000000   0.50000000000000   O    (4h)
   0.10000000000000   0.72000000000000   0.25000000000000   O    (8i)
  -0.10000000000000  -0.72000000000000   0.25000000000000   O    (8i)
   0.40000000000000   1.22000000000000  -0.25000000000000   O    (8i)
   0.60000000000000  -0.22000000000000  -0.25000000000000   O    (8i)
  -0.10000000000000  -0.72000000000000  -0.25000000000000   O    (8i)
   0.10000000000000   0.72000000000000  -0.25000000000000   O    (8i)
   0.60000000000000  -0.22000000000000   0.25000000000000   O    (8i)
   0.40000000000000   1.22000000000000   0.25000000000000   O    (8i)
```

Calciborite (CaB_2O_4 II): A2BC4_oP56_56_2e_e_4e - CIF

```
# CIF file
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_audit_creation_method FINDSYM

_chemical_name_mineral 'Calciborite'
_chemical_formula_sum 'B2 Ca O4'

loop_
_publ_author_name
 'D. N. Shashkin'
 'M. A. Simonov'
```

```
 'N. V. Belov'
_journal_name_full_name
;
 Doklady Akademii Nauk SSSR
;
_journal_volume 195
_journal_year 1970
_journal_page_first 345
_journal_page_last 348
_publ_Section_title
;
 Crystal structure of calciborite CaB$_{2}$O$_{4}$=Ca$_{2}$[BO$_{3}$BO]
      ↪ $_{2}$
;

# Found in The American Mineralogist Crystal Structure Database, 2003

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_aflow_proto 'A2BC4_oP56_56_2e_e_4e'
_aflow_params 'a,b/a,c/a,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3},y_{3
      ↪ },z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{6},
      ↪ x_{7},y_{7},z_{7}'
_aflow_params_values '8.38,1.6491646778,0.597374701671,0.537,0.139,0.624
      ↪ ,0.742,0.052,0.365,0.386,0.143,0.123,0.391,0.185,0.633,0.742,-
      ↪ 0.009,0.114,0.596,0.112,0.365,0.885,0.112,0.378'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP56'

_symmetry_space_group_name_H-M "P 21/c 21/c 2/n"
_symmetry_Int_Tables_number 56

_cell_length_a     8.38000
_cell_length_b     13.82000
_cell_length_c     5.00600
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y,-z+1/2
3 -x,y+1/2,-z+1/2
4 -x+1/2,-y+1/2,z
5 -x,-y,-z
6 -x+1/2,y,z+1/2
7 x,-y+1/2,z+1/2
8 x+1/2,y+1/2,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
B1 B 8 e 0.53700 0.13900 0.62400 1.00000
B2 B 8 e 0.74200 0.05200 0.36500 1.00000
Ca1 Ca 8 e 0.38600 0.14300 0.12300 1.00000
O1 O 8 e 0.39100 0.18500 0.63300 1.00000
O2 O 8 e 0.74200 -0.00900 0.11400 1.00000
O3 O 8 e 0.59600 0.11200 0.36500 1.00000
O4 O 8 e 0.88500 0.11200 0.37800 1.00000
```

Calciborite (CaB_2O_4 II): A2BC4_oP56_56_2e_e_4e - POSCAR

```
A2BC4_oP56_56_2e_e_4e & a,b/a,c/a,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,z4,x5
      ↪ ,y5,z5,x6,y6,z6,x7,y7,z7 --params=8.38,1.6491646778,
      ↪ 0.597374701671,0.537,0.139,0.624,0.742,0.052,0.365,0.386,0.143,
      ↪ 0.123,0.391,0.185,0.633,0.742,-0.009,0.114,0.596,0.112,0.365,
      ↪ 0.885,0.112,0.378 & Pccn D_{2h}^{10} #56 (e^7) & oP56 & None &
      ↪ B2CaO4 & Calciborite & D. N. Shashkin and M. A. Simonov and N.
      ↪ V. Belov, Doklady Akademii Nauk SSSR 195, 345-348 (1970)
   1.00000000000000
     8.38000000000000     0.00000000000000     0.00000000000000
     0.00000000000000    13.82000000000000     0.00000000000000
     0.00000000000000     0.00000000000000     5.00600000000000
       B    Ca     O
      16     8    32
Direct
     0.53700000000000     0.13900000000000     0.62400000000000     B     (8e)
    -0.03700000000000     0.36100000000000     0.62400000000000     B     (8e)
    -0.53700000000000     0.63900000000000    -0.12400000000000     B     (8e)
     1.03700000000000    -0.13900000000000    -0.12400000000000     B     (8e)
    -0.53700000000000    -0.13900000000000    -0.62400000000000     B     (8e)
     1.03700000000000     0.63900000000000    -0.62400000000000     B     (8e)
     0.53700000000000     0.36100000000000     1.12400000000000     B     (8e)
    -0.03700000000000     0.13900000000000     1.12400000000000     B     (8e)
     0.74200000000000     0.05200000000000     0.36500000000000     B     (8e)
    -0.24200000000000     0.44800000000000     0.36500000000000     B     (8e)
    -0.74200000000000     0.55200000000000     0.13500000000000     B     (8e)
     1.24200000000000    -0.05200000000000     0.13500000000000     B     (8e)
    -0.74200000000000    -0.05200000000000    -0.36500000000000     B     (8e)
     1.24200000000000     0.55200000000000    -0.36500000000000     B     (8e)
     0.74200000000000     0.44800000000000     0.86500000000000     B     (8e)
    -0.24200000000000     0.05200000000000     0.86500000000000     B     (8e)
     0.38600000000000     0.14300000000000     0.12300000000000     Ca    (8e)
     0.11400000000000     0.35700000000000     0.12300000000000     Ca    (8e)
    -0.38600000000000     0.64300000000000     0.37700000000000     Ca    (8e)
     0.88600000000000    -0.14300000000000     0.37700000000000     Ca    (8e)
    -0.38600000000000    -0.14300000000000    -0.12300000000000     Ca    (8e)
     0.88600000000000     0.64300000000000    -0.12300000000000     Ca    (8e)
     0.38600000000000     0.35700000000000     0.62300000000000     Ca    (8e)
     0.11400000000000     0.14300000000000     0.62300000000000     Ca    (8e)
     0.39100000000000     0.18500000000000     0.63300000000000     O     (8e)
     0.10900000000000     0.31500000000000     0.63300000000000     O     (8e)
    -0.39100000000000     0.68500000000000    -0.13300000000000     O     (8e)
     0.89100000000000    -0.18500000000000    -0.13300000000000     O     (8e)
    -0.39100000000000    -0.18500000000000    -0.63300000000000     O     (8e)
     0.89100000000000     0.68500000000000    -0.63300000000000     O     (8e)
     0.39100000000000     0.31500000000000     1.13300000000000     O     (8e)
     0.10900000000000     0.18500000000000     1.13300000000000     O     (8e)
     0.74200000000000    -0.00900000000000     0.11400000000000     O     (8e)
    -0.24200000000000     0.50900000000000     0.11400000000000     O     (8e)
    -0.74200000000000     0.49100000000000     0.38600000000000     O     (8e)
     1.24200000000000     0.00900000000000     0.38600000000000     O     (8e)
    -0.74200000000000     0.00900000000000    -0.11400000000000     O     (8e)
     1.24200000000000     0.49100000000000    -0.11400000000000     O     (8e)
     0.74200000000000     0.50900000000000     0.61400000000000     O     (8e)
    -0.24200000000000    -0.00900000000000     0.61400000000000     O     (8e)
     0.59600000000000     0.11200000000000     0.36500000000000     O     (8e)
    -0.09600000000000     0.38800000000000     0.36500000000000     O     (8e)
    -0.59600000000000     0.61200000000000     0.13500000000000     O     (8e)
     1.09600000000000    -0.11200000000000     0.13500000000000     O     (8e)
    -0.59600000000000    -0.11200000000000    -0.36500000000000     O     (8e)
     1.09600000000000     0.61200000000000    -0.36500000000000     O     (8e)
     0.59600000000000     0.38800000000000     0.86500000000000     O     (8e)
    -0.09600000000000     0.11200000000000     0.86500000000000     O     (8e)
     0.88500000000000     0.11200000000000     0.37800000000000     O     (8e)
    -0.38500000000000     0.38800000000000     0.37800000000000     O     (8e)
    -0.88500000000000     0.61200000000000     0.12200000000000     O     (8e)
     1.38500000000000    -0.11200000000000     0.12200000000000     O     (8e)
    -0.88500000000000    -0.11200000000000    -0.37800000000000     O     (8e)
     1.38500000000000     0.61200000000000    -0.37800000000000     O     (8e)
     0.88500000000000     0.38800000000000     0.87800000000000     O     (8e)
    -0.38500000000000     0.11200000000000     0.87800000000000     O     (8e)
```

$D0_{10}$ (WO_3) (*obsolete*): A3B_oP16_57_a2d_d - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'O3W'
_chemical_formula_sum 'O3 W'

loop_
_publ_author_name
 'P. M. Woodward'
 'A. W. Sleight'
 'T. Vogt'
_journal_name_full_name
;
 Journal of Solid State Chemistry
;
_journal_volume 131
_journal_year 1997
_journal_page_first 9
_journal_page_last 17
_publ_Section_title
;
 Ferroelectric Tungsten Trioxide
;

_aflow_title '$D0_{10}$ (WO$_{3}$) ({\em obsolete}) Structure'
_aflow_proto 'A3B_oP16_57_a2d_d'
_aflow_params 'a,b/a,c/a,x_{2},y_{2},x_{3},y_{3},x_{4},y_{4}'
_aflow_params_values '3.8,1.94736842105,1.89473684211,0.0,0.78125,0.5625
      ↪ ,-0.03125,0.0625,-0.03125'
_aflow_Strukturbericht '$D0_{10}$'
_aflow_Pearson 'oP16'

_symmetry_space_group_name_H-M "P 2/b 21/c 21/m"
_symmetry_Int_Tables_number 57

_cell_length_a     3.80000
_cell_length_b     7.40000
_cell_length_c     7.20000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y+1/2,-z
3 -x,y+1/2,-z+1/2
4 -x,-y,z+1/2
5 -x,-y,-z
6 -x,y+1/2,z
7 x,-y+1/2,z+1/2
8 x,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 4 a 0.00000 0.00000 0.00000 1.00000
O2 O 4 d 0.00000 0.78125 0.25000 1.00000
O3 O 4 d 0.56250 -0.03125 0.25000 1.00000
W1 W 4 d 0.06250 -0.03125 0.25000 1.00000
```

$D0_{10}$ (WO_3) (*obsolete*): A3B_oP16_57_a2d_d - POSCAR

```
A3B_oP16_57_a2d_d & a,b/a,c/a,x2,y2,x3,y3,x4,y4 --params=3.8,
    ↪ 1.94736842105,1.89473684211,0.0,0.78125,0.5625,-0.03125,0.0625
    ↪ ,-0.03125 & Pbcm D_{2h}^{11} #57 (ad^3) & oP16 & $D0_{10}$ &
    ↪ O3W & O3W & P. M. Woodward and A. W. Sleight and T. Vogt, J.
    ↪ Solid State Chem. 131, 9-17 (1997)
   1.00000000000000
   3.80000000000000    0.00000000000000    0.00000000000000
   0.00000000000000    7.40000000000000    0.00000000000000
   0.00000000000000    0.00000000000000    7.20000000000000
     O     W
    12     4
Direct
   0.00000000000000    0.00000000000000    0.00000000000000  O  (4a)
   0.00000000000000    0.00000000000000    0.50000000000000  O  (4a)
   0.00000000000000    0.50000000000000    0.50000000000000  O  (4a)
   0.00000000000000    0.50000000000000    0.00000000000000  O  (4a)
   0.00000000000000    0.78125000000000    0.25000000000000  O  (4d)
   0.00000000000000   -0.78125000000000    0.75000000000000  O  (4d)
   0.00000000000000    1.28125000000000    0.25000000000000  O  (4d)
   0.00000000000000   -0.28125000000000    0.75000000000000  O  (4d)
   0.56250000000000   -0.03125000000000    0.25000000000000  O  (4d)
  -0.56250000000000    0.03125000000000    0.75000000000000  O  (4d)
  -0.56250000000000    0.46875000000000    0.25000000000000  O  (4d)
   0.56250000000000    0.53125000000000    0.75000000000000  O  (4d)
   0.06250000000000   -0.03125000000000    0.25000000000000  W  (4d)
  -0.06250000000000    0.03125000000000    0.75000000000000  W  (4d)
  -0.06250000000000    0.46875000000000    0.25000000000000  W  (4d)
   0.06250000000000    0.53125000000000    0.75000000000000  W  (4d)
```

$SrUO_4$: A4BC_oP24_57_cde_d_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'O4SrU'
_chemical_formula_sum 'O4 Sr U'

loop_
_publ_author_name
 'B. O. Loopstra'
 'H. M. Rietveld'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 25
_journal_year 1969
_journal_page_first 787
_journal_page_last 791
_publ_Section_title
;
 The structure of some alkaline-earth metal uranates
;

_aflow_title 'SrUO${4}$ Structure'
_aflow_proto 'A4BC_oP24_57_cde_d_a'
_aflow_params 'a,b/a,c/a,x_{2},x_{3},y_{3},x_{4},y_{4},x_{5},y_{5},z_{5
    ↪ }'
_aflow_params_values '5.4896,1.4531113378,1.48092757214,0.1726,0.8602,
    ↪ 0.03,0.471,0.2013,0.2998,-0.0807,0.081'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP24'

_symmetry_space_group_name_H-M "P 2/b 21/c 21/m"
_symmetry_Int_Tables_number 57

_cell_length_a    5.48960
_cell_length_b    7.97700
_cell_length_c    8.12970
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y+1/2,-z
3 -x,y+1/2,-z+1/2
4 -x,-y,z+1/2
5 -x,-y,-z
6 -x,y+1/2,z
7 x,-y+1/2,z+1/2
8 x,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
U1 U 4 a 0.00000 0.00000 0.00000 1.00000
O1 O 4 c 0.17260 0.25000 0.00000 1.00000
O2 O 4 d 0.86020 0.03000 0.25000 1.00000
Sr1 Sr 4 d 0.47100 0.20130 0.25000 1.00000
O3 O 8 e 0.29980 -0.08070 0.08100 1.00000
```

$SrUO_4$: A4BC_oP24_57_cde_d_a - POSCAR

```
A4BC_oP24_57_cde_d_a & a,b/a,c/a,x2,x3,y3,x4,y4,x5,y5,z5 --params=5.4896
    ↪ ,1.4531113378,1.48092757214,0.1726,0.8602,0.03,0.471,0.2013,
    ↪ 0.2998,-0.0807,0.081 & Pbcm D_{2h}^{11} #57 (acd^2e) & oP24 &
    ↪ None & O4SrU & O4SrU & B. O. Loopstra and H. M. Rietveld, Acta
    ↪ Crystallogr. Sect. B Struct. Sci. 25, 787-791 (1969)
   1.00000000000000
   5.48960000000000    0.00000000000000    0.00000000000000
   0.00000000000000    7.97700000000000    0.00000000000000
   0.00000000000000    0.00000000000000    8.12970000000000
     O     Sr     U
    16     4     4
Direct
   0.17260000000000    0.25000000000000    0.00000000000000  O  (4c)
  -0.17260000000000    0.75000000000000    0.50000000000000  O  (4c)
  -0.17260000000000    0.75000000000000    0.00000000000000  O  (4c)
   0.17260000000000    0.25000000000000    0.50000000000000  O  (4c)
   0.86020000000000    0.03000000000000    0.25000000000000  O  (4d)
  -0.86020000000000   -0.03000000000000    0.75000000000000  O  (4d)
  -0.86020000000000    0.53000000000000    0.25000000000000  O  (4d)
   0.86020000000000    0.47000000000000    0.75000000000000  O  (4d)
   0.29980000000000   -0.08070000000000    0.08100000000000  O  (8e)
  -0.29980000000000    0.08070000000000    0.58100000000000  O  (8e)
  -0.29980000000000    0.41930000000000    0.41900000000000  O  (8e)
   0.29980000000000    0.58070000000000   -0.08100000000000  O  (8e)
  -0.29980000000000    0.08070000000000   -0.08100000000000  O  (8e)
   0.29980000000000   -0.08070000000000    0.41900000000000  O  (8e)
   0.29980000000000    0.58070000000000    0.58100000000000  O  (8e)
  -0.29980000000000    0.41930000000000    0.08100000000000  O  (8e)
   0.47100000000000    0.20130000000000    0.25000000000000  Sr  (4d)
  -0.47100000000000   -0.20130000000000    0.75000000000000  Sr  (4d)
  -0.47100000000000    0.70130000000000    0.25000000000000  Sr  (4d)
   0.47100000000000    0.29870000000000    0.75000000000000  Sr  (4d)
   0.00000000000000    0.00000000000000    0.00000000000000  U  (4a)
   0.00000000000000    0.00000000000000    0.50000000000000  U  (4a)
   0.00000000000000    0.50000000000000    0.50000000000000  U  (4a)
   0.00000000000000    0.50000000000000    0.00000000000000  U  (4a)
```

Lueshite ($NaNbO_3$): ABC3_oP40_57_cd_e_cd2e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Lueshite'
_chemical_formula_sum 'Na Nb O3'

loop_
_publ_author_name
 'A. C. {Sakowski-Cowley}'
 'K. Lukaszewicz'
 'H. D. Megaw'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 25
_journal_year 1969
_journal_page_first 851
_journal_page_last 865
_publ_Section_title
;
 The structure of sodium niobate at room temperature, and the problem of
    ↪ reliability in pseudosymmetric structures
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'Lueshite (NaNbO${3}$) Structure'
_aflow_proto 'ABC3_oP40_57_cd_e_cd2e'
_aflow_params 'a,b/a,c/a,x_{1},x_{2},x_{3},y_{3},x_{4},y_{4},x_{5},y_{5
    ↪ },z_{5},x_{6},y_{6},z_{6},x_{7},y_{7},z_{7}'
_aflow_params_values '5.506,1.01089720305,2.81874318925,0.243,0.696,
    ↪ 0.239,0.218,0.191,0.767,0.2566,0.7722,0.6262,0.536,0.532,0.64,-
    ↪ 0.034,-0.033,0.61'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP40'

_symmetry_space_group_name_H-M "P 2/b 21/c 21/m"
_symmetry_Int_Tables_number 57

_cell_length_a    5.50600
_cell_length_b    5.56600
_cell_length_c    15.52000
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y+1/2,-z
3 -x,y+1/2,-z+1/2
4 -x,-y,z+1/2
5 -x,-y,-z
6 -x,y+1/2,z
7 x,-y+1/2,z+1/2
8 x,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
```

```
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Na1 Na 4 c 0.24300 0.25000 0.00000 1.00000
O1 O 4 c 0.69600 0.25000 0.00000 1.00000
Na2 Na 4 d 0.23900 0.21800 0.25000 1.00000
O2 O 4 d 0.19100 0.76700 0.25000 1.00000
Nb1 Nb 8 e 0.25660 0.77220 0.62620 1.00000
O3 O 8 e 0.53600 0.53200 0.64000 1.00000
O4 O 8 e -0.03400 -0.03300 0.61000 1.00000
```

Lueshite ($NaNbO_3$): ABC3_oP40_57_cd_e_cd2e - POSCAR

```
ABC3_oP40_57_cd_e_cd2e & a,b/a,c/a,x1,x2,x3,y3,x4,y4,x5,y5,z5,x6,y6,z6,
    ↪ x7,y7,z7 --params=5.506,1.01089720305,2.81874318925,0.243,0.696
    ↪ ,0.239,0.218,0.191,0.767,0.2566,0.7722,0.6262,0.536,0.532,0.64
    ↪ ,-0.034,-0.033,0.61 & Pbcm D_{2h}^{11} #57 (c^2d^2e^3) & oP40 &
    ↪  None & NaNbO3 & Lueshite & A. C. {Sakowski-Cowley} and K.
    ↪ Lukaszewicz and H. D. Megaw, Acta Crystallogr. Sect. B Struct.
    ↪ Sci. 25, 851-865 (1969)
   1.00000000000000
   5.50600000000000    0.00000000000000    0.00000000000000
   0.00000000000000    5.56600000000000    0.00000000000000
   0.00000000000000    0.00000000000000   15.52000000000000
     Na     Nb      O
      8      8     24
Direct
   0.24300000000000    0.25000000000000    0.00000000000000   Na   (4c)
  -0.24300000000000    0.75000000000000    0.50000000000000   Na   (4c)
  -0.24300000000000    0.75000000000000    0.00000000000000   Na   (4c)
   0.24300000000000    0.25000000000000    0.50000000000000   Na   (4c)
   0.23900000000000    0.21800000000000    0.25000000000000   Na   (4d)
  -0.23900000000000   -0.21800000000000    0.75000000000000   Na   (4d)
  -0.23900000000000    0.71800000000000    0.25000000000000   Na   (4d)
   0.23900000000000    0.28200000000000    0.75000000000000   Na   (4d)
   0.25660000000000    0.77220000000000    0.62620000000000   Nb   (8e)
  -0.25660000000000   -0.77220000000000    1.12620000000000   Nb   (8e)
  -0.25660000000000    1.27220000000000   -0.12620000000000   Nb   (8e)
   0.25660000000000   -0.27220000000000   -0.62620000000000   Nb   (8e)
  -0.25660000000000   -0.77220000000000   -0.62620000000000   Nb   (8e)
   0.25660000000000    0.77220000000000   -0.12620000000000   Nb   (8e)
   0.25660000000000   -0.27220000000000    1.12620000000000   Nb   (8e)
  -0.25660000000000    1.27220000000000    0.62620000000000   Nb   (8e)
   0.69600000000000    0.25000000000000    0.00000000000000   O    (4c)
  -0.69600000000000    0.75000000000000    0.50000000000000   O    (4c)
  -0.69600000000000    0.75000000000000    0.00000000000000   O    (4c)
   0.69600000000000    0.25000000000000    0.50000000000000   O    (4c)
   0.19100000000000    0.76700000000000    0.25000000000000   O    (4d)
  -0.19100000000000   -0.76700000000000    0.75000000000000   O    (4d)
  -0.19100000000000    1.26700000000000    0.25000000000000   O    (4d)
   0.19100000000000   -0.26700000000000    0.75000000000000   O    (4d)
   0.53600000000000    0.53200000000000    0.64000000000000   O    (8e)
  -0.53600000000000   -0.53200000000000    1.14000000000000   O    (8e)
  -0.53600000000000    1.03200000000000   -0.14000000000000   O    (8e)
   0.53600000000000   -0.03200000000000   -0.64000000000000   O    (8e)
  -0.53600000000000   -0.53200000000000   -0.64000000000000   O    (8e)
   0.53600000000000    0.53200000000000   -0.14000000000000   O    (8e)
   0.53600000000000   -0.03200000000000    1.14000000000000   O    (8e)
  -0.53600000000000    1.03200000000000    0.64000000000000   O    (8e)
  -0.03400000000000   -0.03300000000000    0.61000000000000   O    (8e)
   0.03400000000000    0.03300000000000    1.11000000000000   O    (8e)
   0.03400000000000    0.46700000000000   -0.11000000000000   O    (8e)
  -0.03400000000000    0.53300000000000   -0.61000000000000   O    (8e)
   0.03400000000000    0.03300000000000   -0.61000000000000   O    (8e)
  -0.03400000000000   -0.03300000000000   -0.11000000000000   O    (8e)
  -0.03400000000000    0.53300000000000    1.11000000000000   O    (8e)
   0.03400000000000    0.46700000000000    0.61000000000000   O    (8e)
```

Kotoite ($Mg_3(BO_3)_2$): A2B3C6_oP22_58_g_af_gh - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Kotoite'
_chemical_formula_sum 'B2 Mg3 O6'

loop_
_publ_author_name
 'S. V. Berger'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 3
_journal_year 1949
_journal_page_first 660
_journal_page_last 675
_publ_Section_title
;
 The Crystal Structure of the Isomorphous Orthoborates of Cobalt and
    ↪ Magnesium
;

_aflow_title 'Kotoite (Mg$_{3}$(BO$_{3}$)$_{2}$) Structure'
_aflow_proto 'A2B3C6_oP22_58_g_af_gh'
_aflow_params 'a,b/a,c/a,z_{2},x_{3},y_{3},x_{4},y_{4},x_{5},y_{5},z_{5
    ↪ }'
_aflow_params_values '4.497,1.20035579275,1.8714698688,0.179,0.56,0.25,
    ↪ 0.258,0.316,0.705,0.218,0.139'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP22'

_symmetry_space_group_name_H-M "P 21/n 21/n 2/m"
_symmetry_Int_Tables_number 58
_cell_length_a    4.49700
_cell_length_b    5.39800
_cell_length_c    8.41600
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x+1/2,y+1/2,-z+1/2
4 -x,-y,z
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x+1/2,-y+1/2,z+1/2
8 x,y,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Mg1 Mg 2 a 0.00000 0.00000 0.00000 1.00000
Mg2 Mg 4 f 0.00000 0.50000 0.17900 1.00000
B1 B 4 g 0.56000 0.25000 0.00000 1.00000
O1 O 4 g 0.25800 0.31600 0.00000 1.00000
O2 O 8 h 0.70500 0.21800 0.13900 1.00000
```

Kotoite ($Mg_3(BO_3)_2$): A2B3C6_oP22_58_g_af_gh - POSCAR

```
A2B3C6_oP22_58_g_af_gh & a,b/a,c/a,z2,x3,y3,x4,y4,x5,y5,z5 --params=
    ↪ 4.497,1.20035579275,1.8714698688,0.179,0.56,0.25,0.258,0.316,
    ↪ 0.705,0.218,0.139 & Pnnm D_{2h}^{12} #58 (afg^2h) & oP22 & None
    ↪  & B2Mg3O6 & Kotoite & S. V. Berger, Acta Chem. Scand. 3,
    ↪ 660-675 (1949)
   1.00000000000000
   4.49700000000000    0.00000000000000    0.00000000000000
   0.00000000000000    5.39800000000000    0.00000000000000
   0.00000000000000    0.00000000000000    8.41600000000000
     B     Mg      O
     4      6     12
Direct
   0.56000000000000    0.25000000000000    0.00000000000000   B    (4g)
  -0.56000000000000   -0.25000000000000    0.00000000000000   B    (4g)
  -0.06000000000000    0.75000000000000    0.50000000000000   B    (4g)
   1.06000000000000    0.25000000000000    0.50000000000000   B    (4g)
   0.00000000000000    0.00000000000000    0.00000000000000   Mg   (2a)
   0.50000000000000    0.50000000000000    0.50000000000000   Mg   (2a)
   0.00000000000000    0.50000000000000    0.17900000000000   Mg   (4f)
   0.50000000000000    0.00000000000000    0.32100000000000   Mg   (4f)
   0.00000000000000    0.50000000000000   -0.17900000000000   Mg   (4f)
   0.50000000000000    0.00000000000000    0.67900000000000   Mg   (4f)
   0.25800000000000    0.31600000000000    0.00000000000000   O    (4g)
  -0.25800000000000   -0.31600000000000    0.00000000000000   O    (4g)
   0.24200000000000    0.81600000000000    0.50000000000000   O    (4g)
   0.75800000000000    0.18400000000000    0.50000000000000   O    (4g)
   0.70500000000000    0.21800000000000    0.13900000000000   O    (8h)
  -0.70500000000000   -0.21800000000000    0.13900000000000   O    (8h)
  -0.20500000000000    0.71800000000000    0.36100000000000   O    (8h)
   1.20500000000000    0.28200000000000    0.36100000000000   O    (8h)
  -0.70500000000000   -0.21800000000000   -0.13900000000000   O    (8h)
   0.70500000000000    0.21800000000000   -0.13900000000000   O    (8h)
   1.20500000000000    0.28200000000000    0.63900000000000   O    (8h)
  -0.20500000000000    0.71800000000000    0.63900000000000   O    (8h)
```

Andalusite (Al_2SiO_5, $S0_2$): A2B5C_oP32_58_eg_3gh_g - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Andalusite'
_chemical_formula_sum 'Al2 O5 S'

loop_
_publ_author_name
 'J. K. Winter'
 'S. Ghose'
_journal_name_full_name
;
 American Mineralogist
;
_journal_volume 64
_journal_year 1979
_journal_page_first 573
_journal_page_last 586
_publ_Section_title
;
 Thermal expansion and high-temperaturce crystal chemistry of the Al$_{2
    ↪ }$SiO$_{4}$ polymorphs
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'Andalusite (Al$_{2}$SiO$_{5}$, $S0_{2}$) Structure'
_aflow_proto 'A2B5C_oP32_58_eg_3gh_g'
_aflow_params 'a,b/a,c/a,z_{1},x_{2},y_{2},x_{3},y_{3},x_{4},y_{4},x_{5
    ↪ },y_{5},x_{6},y_{6},x_{7},y_{7},z_{7}'
```

```
_aflow_params_values '7.798,1.01347781482,0.712567324955,0.2419,0.1295,
    ↪ 0.6391,0.0767,0.8629,0.4246,0.3629,0.103,0.4003,0.246,0.252,
    ↪ 0.2305,0.1339,0.2394'
_aflow_Strukturbericht '$S0_{2}$'
_aflow_Pearson 'oP32'

_symmetry_space_group_name_H-M "P 21/n 21/n 2/m"
_symmetry_Int_Tables_number 58

_cell_length_a 7.79800
_cell_length_b 7.90310
_cell_length_c 5.55660
_cell_angle_alpha 90.00000
_cell_angle_beta 90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x+1/2,y+1/2,-z+1/2
4 -x,-y,z
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x+1/2,-y+1/2,z+1/2
8 x,y,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Al1 Al 4 e 0.00000 0.00000 0.24190 1.00000
Al2 Al 4 g 0.12950 0.63910 0.00000 1.00000
O1 O 4 g 0.07670 0.86290 0.00000 1.00000
O2 O 4 g 0.42460 0.36290 0.00000 1.00000
O3 O 4 g 0.10300 0.40030 0.00000 1.00000
S1 S 4 g 0.24600 0.25200 0.00000 1.00000
O4 O 8 h 0.23050 0.13390 0.23940 1.00000
```

Andalusite (Al_2SiO_5, $S0_2$): A2B5C_oP32_58_eg_3gh_g - POSCAR

```
A2B5C_oP32_58_eg_3gh_g & a,b/a,c/a,z1,x2,y2,x3,y3,x4,y4,x5,y5,x6,y6,x7,
    ↪ y7,z7 --params=7.798,1.01347781482,0.712567324955,0.2419,0.1295
    ↪ ,0.6391,0.0767,0.8629,0.4246,0.3629,0.103,0.4003,0.246,0.252,
    ↪ 0.2305,0.1339,0.2394 & Pnnm D_{2h}^{12} #58 (eg^5h) & oP32 &
    ↪ $S0_{2}$ & Al2O5Si & Andalusite & J. K. Winter and S. Ghose,
    ↪ Am. Mineral. 64, 573-586 (1979)
   1.00000000000000
   7.79800000000000    0.00000000000000    0.00000000000000
   0.00000000000000    7.90310000000000    0.00000000000000
   0.00000000000000    0.00000000000000    5.55660000000000
     Al      O      S
      8     20      4
Direct
   0.00000000000000    0.00000000000000    0.24190000000000   Al   (4e)
   0.50000000000000    0.50000000000000    0.25810000000000   Al   (4e)
   0.00000000000000    0.00000000000000   -0.24190000000000   Al   (4e)
   0.50000000000000    0.50000000000000    0.74190000000000   Al   (4e)
   0.12950000000000    0.63910000000000    0.00000000000000   Al   (4g)
  -0.12950000000000   -0.63910000000000    0.00000000000000   Al   (4g)
   0.37050000000000    1.13910000000000    0.50000000000000   Al   (4g)
   0.62950000000000   -0.13910000000000    0.50000000000000   Al   (4g)
   0.07670000000000    0.86290000000000    0.00000000000000   O    (4g)
  -0.07670000000000   -0.86290000000000    0.00000000000000   O    (4g)
   0.42330000000000    1.36290000000000    0.50000000000000   O    (4g)
   0.57670000000000   -0.36290000000000    0.50000000000000   O    (4g)
   0.42460000000000    0.36290000000000    0.00000000000000   O    (4g)
  -0.42460000000000   -0.36290000000000    0.00000000000000   O    (4g)
   0.07540000000000    0.86290000000000    0.50000000000000   O    (4g)
   0.92460000000000    0.13710000000000    0.50000000000000   O    (4g)
   0.10300000000000    0.40030000000000    0.00000000000000   O    (4g)
  -0.10300000000000   -0.40030000000000    0.00000000000000   O    (4g)
   0.39700000000000    0.90030000000000    0.50000000000000   O    (4g)
   0.60300000000000    0.09970000000000    0.50000000000000   O    (4g)
   0.23050000000000    0.13390000000000    0.23940000000000   O    (8h)
  -0.23050000000000   -0.13390000000000    0.23940000000000   O    (8h)
   0.26950000000000    0.63390000000000    0.26060000000000   O    (8h)
   0.73050000000000    0.36610000000000    0.26060000000000   O    (8h)
  -0.23050000000000   -0.13390000000000   -0.23940000000000   O    (8h)
   0.23050000000000    0.13390000000000   -0.23940000000000   O    (8h)
   0.73050000000000    0.36610000000000    0.73940000000000   O    (8h)
   0.26950000000000    0.63390000000000    0.73940000000000   O    (8h)
   0.24600000000000    0.25200000000000    0.00000000000000   S    (4g)
  -0.24600000000000   -0.25200000000000    0.00000000000000   S    (4g)
   0.25400000000000    0.75200000000000    0.50000000000000   S    (4g)
   0.74600000000000    0.24800000000000    0.50000000000000   S    (4g)
```

Protoanthophyllite ($H_2Mg_7Si_8O_{24}$): A2B7C24D8_oP82_58_g_ae2f_2g5h_2h - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Protoanthophyllite'
_chemical_formula_sum 'H2 Mg7 O24 Si8'

loop_
_publ_author_name
 'H. Konishi'
 'T. L. Groy'
 'I. D\\'{o}dony'
 'R. Miyawaki'
 'S. Matsubara'
 'P. R. Buseck'
_journal_name_full_name
;
 American Mineralogist
;
_journal_volume 88
_journal_year 2003
_journal_page_first 1718
_journal_page_last 1723
_publ_Section_title
;
 Crystal structure of protoanthophyllite: A new mineral from the Takase
    ↪ ultramafic complex, Japan
;

_aflow_title 'Protoanthophyllite (H$_{2}$Mg$_{7}$Si$_{8}$O$_{24}$)
    ↪ Structure'
_aflow_proto 'A2B7C24D8_oP82_58_g_ae2f_2g5h_2h'
_aflow_params 'a,b/a,c/a,z_{2},z_{3},z_{4},x_{5},y_{5},x_{6},y_{6},x_{7
    ↪ },y_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{10},y_{10},z_{10
    ↪ },x_{11},y_{11},z_{11},x_{12},y_{12},z_{12},x_{13},y_{13},z_{13
    ↪ },x_{14},y_{14},z_{14}'
_aflow_params_values '5.3117,1.76126287253,3.37571775514,0.17748,0.41276
    ↪ ,0.24009,0.682,0.225,0.6656,0.1128,0.1502,0.3414,0.1666,0.11279
    ↪ ,0.08868,0.6699,0.12137,0.17464,0.1834,0.1211,0.25265,0.433,
    ↪ 0.3436,0.11987,-0.0626,0.34888,0.13244,0.17327,0.28464,0.08483,
    ↪ 0.67034,0.29399,0.17095'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP82'

_symmetry_space_group_name_H-M "P 21/n 21/n 2/m"
_symmetry_Int_Tables_number 58

_cell_length_a 5.31170
_cell_length_b 9.35530
_cell_length_c 17.93080
_cell_angle_alpha 90.00000
_cell_angle_beta 90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x+1/2,y+1/2,-z+1/2
4 -x,-y,z
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x+1/2,-y+1/2,z+1/2
8 x,y,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Mg1 Mg 2 a 0.00000 0.00000 0.00000 1.00000
Mg2 Mg 4 e 0.00000 0.00000 0.17748 1.00000
Mg3 Mg 4 f 0.00000 0.50000 0.41276 1.00000
Mg4 Mg 4 f 0.00000 0.50000 0.24009 1.00000
H1 H 4 g 0.68200 0.22500 0.00000 1.00000
O1 O 4 g 0.66560 0.11280 0.00000 1.00000
O2 O 4 g 0.15020 0.34140 0.00000 1.00000
O3 O 8 h 0.16660 0.11279 0.08868 1.00000
O4 O 8 h 0.66990 0.12137 0.17464 1.00000
O5 O 8 h 0.18340 0.12110 0.25265 1.00000
O6 O 8 h 0.43300 0.34360 0.11987 1.00000
O7 O 8 h -0.06260 0.34888 0.13244 1.00000
Si1 Si 8 h 0.17327 0.28464 0.08483 1.00000
Si2 Si 8 h 0.67034 0.29399 0.17095 1.00000
```

Protoanthophyllite ($H_2Mg_7Si_8O_{24}$): A2B7C24D8_oP82_58_g_ae2f_2g5h_2h - POSCAR

```
A2B7C24D8_oP82_58_g_ae2f_2g5h_2h & a,b/a,c/a,z2,z3,z4,x5,y5,x6,y6,x7,y7,
    ↪ x8,y8,z8,x9,y9,z9,x10,y10,z10,x11,y11,z11,x12,y12,z12,x13,y13,
    ↪ z13,x14,y14,z14 --params=5.3117,1.76126287253,3.37571775514,
    ↪ 0.17748,0.41276,0.24009,0.682,0.225,0.6656,0.1128,0.1502,0.3414
    ↪ ,0.1666,0.11279,0.08868,0.6699,0.12137,0.17464,0.1834,0.1211,
    ↪ 0.25265,0.433,0.3436,0.11987,-0.0626,0.34888,0.13244,0.17327,
    ↪ 0.28464,0.08483,0.67034,0.29399,0.17095 & Pnnm D_{2h}^{12} #58
    ↪ (aef^2g^3h^7) & oP82 & None & H2Mg7O24Si8 & Protoanthophyllite
    ↪ & H. Konishi et al., Am. Mineral. 88, 1718-1723 (2003)
   1.00000000000000
   5.31170000000000    0.00000000000000    0.00000000000000
   0.00000000000000    9.35530000000000    0.00000000000000
   0.00000000000000    0.00000000000000   17.93080000000000
     H      Mg      O      Si
     4      14      48     16
Direct
   0.68200000000000    0.22500000000000    0.00000000000000   H    (4g)
  -0.68200000000000   -0.22500000000000    0.00000000000000   H    (4g)
  -0.18200000000000    0.72500000000000    0.50000000000000   H    (4g)
   1.18200000000000    0.27500000000000    0.50000000000000   H    (4g)
   0.00000000000000    0.00000000000000    0.00000000000000   Mg   (2a)
   0.50000000000000    0.50000000000000    0.50000000000000   Mg   (2a)
   0.00000000000000    0.00000000000000    0.17748000000000   Mg   (4e)
```

```
  0.50000000000000   0.50000000000000   0.32252000000000   Mg   (4e)
  0.00000000000000   0.00000000000000  -0.17748000000000   Mg   (4e)
  0.50000000000000   0.50000000000000   0.67748000000000   Mg   (4e)
  0.00000000000000   0.50000000000000   0.41276000000000   Mg   (4f)
  0.50000000000000   0.00000000000000   0.08724000000000   Mg   (4f)
  0.00000000000000   0.50000000000000  -0.41276000000000   Mg   (4f)
  0.50000000000000   0.00000000000000   0.91276000000000   Mg   (4f)
  0.00000000000000   0.50000000000000   0.24009000000000   Mg   (4f)
  0.50000000000000   0.00000000000000   0.25991000000000   Mg   (4f)
  0.00000000000000   0.50000000000000  -0.24009000000000   Mg   (4f)
  0.50000000000000   0.00000000000000   0.74009000000000   Mg   (4f)
  0.66560000000000   0.11280000000000   0.00000000000000   O    (4g)
 -0.66560000000000  -0.11280000000000   0.00000000000000   O    (4g)
 -0.16560000000000   0.61280000000000   0.50000000000000   O    (4g)
  1.16560000000000   0.38720000000000   0.50000000000000   O    (4g)
  0.15020000000000   0.34140000000000   0.00000000000000   O    (4g)
 -0.15020000000000  -0.34140000000000   0.00000000000000   O    (4g)
  0.34980000000000   0.84140000000000   0.50000000000000   O    (4g)
  0.65020000000000   0.15860000000000   0.50000000000000   O    (4g)
  0.16660000000000   0.11279000000000   0.08868000000000   O    (8h)
 -0.16660000000000  -0.11279000000000   0.08868000000000   O    (8h)
  0.33340000000000   0.61279000000000   0.41132000000000   O    (8h)
  0.66660000000000   0.38721000000000   0.41132000000000   O    (8h)
 -0.16660000000000  -0.11279000000000  -0.08868000000000   O    (8h)
  0.16660000000000   0.11279000000000  -0.08868000000000   O    (8h)
  0.66660000000000   0.38721000000000   0.58868000000000   O    (8h)
  0.33340000000000   0.61279000000000   0.58868000000000   O    (8h)
  0.66990000000000   0.12137000000000   0.17464000000000   O    (8h)
 -0.66990000000000  -0.12137000000000   0.17464000000000   O    (8h)
 -0.16990000000000   0.62137000000000   0.32536000000000   O    (8h)
  1.16990000000000   0.37863000000000   0.32536000000000   O    (8h)
 -0.66990000000000  -0.12137000000000  -0.17464000000000   O    (8h)
  0.66990000000000   0.12137000000000  -0.17464000000000   O    (8h)
  1.16990000000000   0.37863000000000   0.67464000000000   O    (8h)
 -0.16990000000000   0.62137000000000   0.67464000000000   O    (8h)
  0.18340000000000   0.12110000000000   0.25265000000000   O    (8h)
 -0.18340000000000  -0.12110000000000   0.25265000000000   O    (8h)
  0.31660000000000   0.62110000000000   0.24735000000000   O    (8h)
  0.68340000000000   0.37890000000000   0.24735000000000   O    (8h)
 -0.18340000000000  -0.12110000000000  -0.25265000000000   O    (8h)
  0.18340000000000   0.12110000000000  -0.25265000000000   O    (8h)
  0.68340000000000   0.37890000000000   0.75265000000000   O    (8h)
  0.31660000000000   0.62110000000000   0.75265000000000   O    (8h)
  0.43300000000000   0.34360000000000   0.11987000000000   O    (8h)
 -0.43300000000000  -0.34360000000000   0.11987000000000   O    (8h)
  0.06700000000000   0.84360000000000   0.38013000000000   O    (8h)
  0.93300000000000   0.15640000000000   0.38013000000000   O    (8h)
 -0.43300000000000  -0.34360000000000  -0.11987000000000   O    (8h)
  0.43300000000000   0.34360000000000  -0.11987000000000   O    (8h)
  0.93300000000000   0.15640000000000   0.61987000000000   O    (8h)
  0.06700000000000   0.84360000000000   0.61987000000000   O    (8h)
 -0.06260000000000   0.34888000000000   0.13244000000000   O    (8h)
  0.06260000000000  -0.34888000000000   0.13244000000000   O    (8h)
  0.56260000000000   0.84888000000000   0.36756000000000   O    (8h)
  0.43740000000000   0.15112000000000   0.36756000000000   O    (8h)
  0.06260000000000  -0.34888000000000  -0.13244000000000   O    (8h)
 -0.06260000000000   0.34888000000000  -0.13244000000000   O    (8h)
  0.43740000000000   0.15112000000000   0.63244000000000   O    (8h)
  0.56260000000000   0.84888000000000   0.63244000000000   O    (8h)
  0.17327000000000   0.28464000000000   0.08483000000000   Si   (8h)
 -0.17327000000000  -0.28464000000000   0.08483000000000   Si   (8h)
  0.32673000000000   0.78464000000000   0.41517000000000   Si   (8h)
  0.67327000000000   0.21536000000000   0.41517000000000   Si   (8h)
 -0.17327000000000  -0.28464000000000  -0.08483000000000   Si   (8h)
  0.17327000000000   0.28464000000000  -0.08483000000000   Si   (8h)
  0.67327000000000   0.21536000000000   0.58483000000000   Si   (8h)
  0.32673000000000   0.78464000000000   0.58483000000000   Si   (8h)
  0.67034000000000   0.29399000000000   0.17095000000000   Si   (8h)
 -0.67034000000000  -0.29399000000000   0.17095000000000   Si   (8h)
 -0.17034000000000   0.79399000000000   0.32905000000000   Si   (8h)
  1.17034000000000   0.20601000000000   0.32905000000000   Si   (8h)
 -0.67034000000000  -0.29399000000000  -0.17095000000000   Si   (8h)
  0.67034000000000   0.29399000000000  -0.17095000000000   Si   (8h)
  1.17034000000000   0.20601000000000   0.67095000000000   Si   (8h)
 -0.17034000000000   0.79399000000000   0.67095000000000   Si   (8h)
```

In_4Se_3: A4B3_oP28_58_4g_3g - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'In4Se3'
_chemical_formula_sum 'In4 Se3'

loop_
_publ_author_name
 'J. H. C. Hogg'
 'H. H. Sutherland'
 'D. J. Williams'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 29
_journal_year 1973
_journal_page_first 1590
_journal_page_last 1593
_publ_Section_title
;
 The crystal structure of tetraindium triselenide
;

# Found in Structural and optical properties of In$_{4}$Se$_{3}$ thin
    ↪ films obtained by flash evaporation, 1995

_aflow_title 'In$_{4}$Se$_{3}$ Structure'
_aflow_proto 'A4B3_oP28_58_4g_3g'
_aflow_params 'a,b/a,c/a,x_{1},y_{1},x_{2},y_{2},x_{3},y_{3},x_{4},y_{4
    ↪ },x_{5},y_{5},x_{6},y_{6},x_{7},y_{7}'
_aflow_params_values '15.297,0.804602209584,0.266784336798,0.7111,0.3393
    ↪ ,0.8157,0.5236,-0.0325,0.6442,0.4238,0.3974,-0.0967,0.8493,
    ↪ 0.7688,0.1386,0.4239,0.156'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP28'

_symmetry_space_group_name_H-M "P 21/n 21/n 2/m"
_symmetry_Int_Tables_number 58

_cell_length_a      15.29700
_cell_length_b      12.30800
_cell_length_c      4.08100
_cell_angle_alpha   90.00000
_cell_angle_beta    90.00000
_cell_angle_gamma   90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x+1/2,y+1/2,-z+1/2
4 -x,-y,z
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x+1/2,-y+1/2,z+1/2
8 x,y,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
In1 In 4 g 0.71110 0.33930 0.00000 1.00000
In2 In 4 g 0.81570 0.52360 0.00000 1.00000
In3 In 4 g -0.03250 0.64420 0.00000 1.00000
In4 In 4 g 0.42380 0.39740 0.00000 1.00000
Se1 Se 4 g -0.09670 0.84930 0.00000 1.00000
Se2 Se 4 g 0.76880 0.13860 0.00000 1.00000
Se3 Se 4 g 0.42390 0.15600 0.00000 1.00000
```

In_4Se_3: A4B3_oP28_58_4g_3g - POSCAR

```
A4B3_oP28_58_4g_3g & a,b/a,c/a,x1,y1,x2,y2,x3,y3,x4,y4,x5,y5,x6,y6,x7,y7
    ↪ --params=15.297,0.804602209584,0.266784336798,0.7111,0.3393,
    ↪ 0.8157,0.5236,-0.0325,0.6442,0.4238,0.3974,-0.0967,0.8493,
    ↪ 0.7688,0.1386,0.4239,0.156 & Pnnm D_{2h}^{12} #58 (g^7) & oP28
    ↪ & None & In4Se3 & In4Se3 & J. H. C. Hogg and H. H. Sutherland
    ↪ and D. J. Williams, Acta Crystallogr. Sect. B Struct. Sci. 29,
    ↪ 1590-1593 (1973)
   1.00000000000000
  15.29700000000000    0.00000000000000    0.00000000000000
   0.00000000000000   12.30800000000000    0.00000000000000
   0.00000000000000    0.00000000000000    4.08100000000000
     In      Se
     16      12
Direct
   0.71110000000000    0.33930000000000    0.00000000000000   In   (4g)
  -0.71110000000000   -0.33930000000000    0.00000000000000   In   (4g)
  -0.21110000000000    0.83930000000000    0.50000000000000   In   (4g)
   1.21110000000000    0.16070000000000    0.50000000000000   In   (4g)
   0.81570000000000    0.52360000000000    0.00000000000000   In   (4g)
  -0.81570000000000   -0.52360000000000    0.00000000000000   In   (4g)
  -0.31570000000000    1.02360000000000    0.50000000000000   In   (4g)
   1.31570000000000   -0.02360000000000    0.50000000000000   In   (4g)
  -0.03250000000000    0.64420000000000    0.00000000000000   In   (4g)
   0.03250000000000   -0.64420000000000    0.00000000000000   In   (4g)
   0.53250000000000    1.14420000000000    0.50000000000000   In   (4g)
   0.46750000000000   -0.14420000000000    0.50000000000000   In   (4g)
   0.42380000000000    0.39740000000000    0.00000000000000   In   (4g)
  -0.42380000000000   -0.39740000000000    0.00000000000000   In   (4g)
   0.07620000000000    0.89740000000000    0.50000000000000   In   (4g)
   0.92380000000000    0.10260000000000    0.50000000000000   In   (4g)
  -0.09670000000000    0.84930000000000    0.00000000000000   Se   (4g)
   0.09670000000000   -0.84930000000000    0.00000000000000   Se   (4g)
   0.59670000000000    1.34930000000000    0.50000000000000   Se   (4g)
   0.40330000000000   -0.34930000000000    0.50000000000000   Se   (4g)
   0.76880000000000    0.13860000000000    0.00000000000000   Se   (4g)
  -0.76880000000000   -0.13860000000000    0.00000000000000   Se   (4g)
  -0.26880000000000    0.63860000000000    0.50000000000000   Se   (4g)
   1.26880000000000    0.36140000000000    0.50000000000000   Se   (4g)
   0.42390000000000    0.15600000000000    0.00000000000000   Se   (4g)
  -0.42390000000000   -0.15600000000000    0.00000000000000   Se   (4g)
   0.07610000000000    0.65600000000000    0.50000000000000   Se   (4g)
   0.92390000000000    0.34400000000000    0.50000000000000   Se   (4g)
```

Adamite [$Zn_2(AsO_4)(OH)$, $H2_7$]: ABC5D2_oP36_58_g_g_3gh_eg - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Adamite'
_chemical_formula_sum 'As H O5 Zn2'

loop_
```

```
_publ_author_name
 'R. J. Hill'
_journal_name_full_name
;
 American Mineralogist
;
_journal_volume 61
_journal_year 1976
_journal_page_first 979
_journal_page_last 986
_publ_Section_title
;
 The crystal structure and infrared properties of adamite
;

_aflow_title 'Adamite [Zn$_{2}$(AsO$_{4}$)(OH), $H2_{7}$] Structure'
_aflow_proto 'ABC5D2_oP36_58_g_g_3gh_eg'
_aflow_params 'a,b/a,c/a,z_{1},x_{2},y_{2},x_{3},y_{3},x_{4},y_{4},x_{5
      ↪ },y_{5},x_{6},y_{6},x_{7},y_{7},x_{8},y_{8},z_{8}'
_aflow_params_values '8.306,1.02624608717,0.727546352035,0.24737,0.24952
      ↪ ,0.74394,0.2,0.13,0.424,0.6447,0.1079,0.1268,0.104,0.6063,
      ↪ 0.13482,0.36423,0.2685,0.3615,0.2778'
_aflow_Strukturbericht '$H2_{7}$'
_aflow_Pearson 'oP36'

_symmetry_space_group_name_H-M "P 21/n 21/n 2/m"
_symmetry_Int_Tables_number 58

_cell_length_a    8.30600
_cell_length_b    8.52400
_cell_length_c    6.04300
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x+1/2,y+1/2,-z+1/2
4 -x,-y,z
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x+1/2,-y+1/2,z+1/2
8 x,y,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Zn1 Zn 4 e 0.00000 0.00000 0.24737 1.00000
As1 As 4 g 0.24952 0.74394 0.00000 1.00000
H1 H 4 g 0.20000 0.13000 0.00000 1.00000
O1 O 4 g 0.42400 0.64470 0.00000 1.00000
O2 O 4 g 0.10790 0.12680 0.00000 1.00000
O3 O 4 g 0.10400 0.60630 0.00000 1.00000
Zn2 Zn 4 g 0.13482 0.36423 0.00000 1.00000
O4 O 8 h 0.26850 0.36150 0.27780 1.00000
```

Adamite [$Zn_2(AsO_4)(OH)$, $H2_7$]: ABC5D2_oP36_58_g_g_3gh_eg - POSCAR

```
ABC5D2_oP36_58_g_g_3gh_eg & a,b/a,c/a,z1,x2,y2,x3,y3,x4,y4,x5,y5,x6,y6,
      ↪ x7,y7,x8,y8,z8 --params=8.306,1.02624608717,0.727546352035,
      ↪ 0.24737,0.24952,0.74394,0.2,0.13,0.424,0.6447,0.1079,0.1268,
      ↪ 0.104,0.6063,0.13482,0.36423,0.2685,0.3615,0.2778 & Pnnm D_{2h
      ↪ }^{12} #58 (eg^6h) & oP36 & $H2_{7}$ & AsHO5Zn2 & Adamite & R.
      ↪ J. Hill, Am. Mineral. 61, 979-986 (1976)
   1.00000000000000
   8.30600000000000    0.00000000000000    0.00000000000000
   0.00000000000000    8.52400000000000    0.00000000000000
   0.00000000000000    0.00000000000000    6.04300000000000
     As   H   O   Zn
      4   4  20    8
Direct
   0.24952000000000    0.74394000000000    0.00000000000000  As   (4g)
  -0.24952000000000   -0.74394000000000    0.00000000000000  As   (4g)
   0.25048000000000    1.24394000000000    0.50000000000000  As   (4g)
   0.74952000000000   -0.24394000000000    0.50000000000000  As   (4g)
   0.20000000000000    0.13000000000000    0.00000000000000  H    (4g)
  -0.20000000000000   -0.13000000000000    0.00000000000000  H    (4g)
   0.30000000000000    0.63000000000000    0.50000000000000  H    (4g)
   0.70000000000000    0.37000000000000    0.50000000000000  H    (4g)
   0.42400000000000    0.64470000000000    0.00000000000000  O    (4g)
  -0.42400000000000   -0.64470000000000    0.00000000000000  O    (4g)
   0.07600000000000    1.14470000000000    0.50000000000000  O    (4g)
   0.92400000000000   -0.14470000000000    0.50000000000000  O    (4g)
   0.10790000000000    0.12680000000000    0.00000000000000  O    (4g)
  -0.10790000000000   -0.12680000000000    0.00000000000000  O    (4g)
   0.39210000000000    0.62680000000000    0.50000000000000  O    (4g)
   0.60790000000000    0.37320000000000    0.50000000000000  O    (4g)
   0.10400000000000    0.60630000000000    0.00000000000000  O    (4g)
  -0.10400000000000   -0.60630000000000    0.00000000000000  O    (4g)
   0.39600000000000    1.10630000000000    0.50000000000000  O    (4g)
   0.60400000000000   -0.10630000000000    0.50000000000000  O    (4g)
   0.26850000000000    0.36150000000000    0.27780000000000  O    (8h)
  -0.26850000000000   -0.36150000000000    0.27780000000000  O    (8h)
   0.23150000000000    0.86150000000000    0.22220000000000  O    (8h)
   0.76850000000000    0.13850000000000    0.22220000000000  O    (8h)
  -0.26850000000000   -0.36150000000000   -0.27780000000000  O    (8h)
   0.26850000000000    0.36150000000000   -0.27780000000000  O    (8h)
   0.76850000000000    0.13850000000000    0.77780000000000  O    (8h)
   0.23150000000000    0.86150000000000    0.77780000000000  O    (8h)
   0.00000000000000    0.00000000000000    0.24737000000000  Zn   (4e)
   0.50000000000000    0.50000000000000    0.25263000000000  Zn   (4e)
   0.00000000000000    0.00000000000000   -0.24737000000000  Zn   (4e)
   0.50000000000000    0.50000000000000    0.74737000000000  Zn   (4e)
   0.13482000000000    0.36423000000000    0.00000000000000  Zn   (4g)
  -0.13482000000000   -0.36423000000000    0.00000000000000  Zn   (4g)
   0.36518000000000    0.86423000000000    0.50000000000000  Zn   (4g)
   0.63482000000000    0.13577000000000    0.50000000000000  Zn   (4g)
```

InS: AB_oP8_58_g_g - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'InS'
_chemical_formula_sum 'In S'

loop_
_publ_author_name
 'K. Schubert'
 'E. D\"{o}rre'
 'E. G\"{u}nzel'
_journal_name_full_name
;
 Naturwissenschaften
;
_journal_volume 41
_journal_year 1954
_journal_page_first 448
_journal_page_last 448
_publ_Section_title
;
 Kristallchemische Ergebnisse an Phasen aus B-Elementen
;

# Found in Pearson's Handbook of Crystallographic Data for Intermetallic
      ↪ Phases, 1991

_aflow_title 'InS Structure'
_aflow_proto 'AB_oP8_58_g_g'
_aflow_params 'a,b/a,c/a,x_{1},y_{1},x_{2},y_{2}'
_aflow_params_values '4.443,2.39522844925,0.886788206167,0.125,0.121,-
      ↪ 0.005,0.355'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP8'

_symmetry_space_group_name_H-M "P 21/n 21/n 2/m"
_symmetry_Int_Tables_number 58

_cell_length_a    4.44300
_cell_length_b    10.64200
_cell_length_c    3.94000
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x+1/2,y+1/2,-z+1/2
4 -x,-y,z
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x+1/2,-y+1/2,z+1/2
8 x,y,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
In1 In 4 g 0.12500 0.12100 0.00000 1.00000
S1 S 4 g -0.00500 0.35500 0.00000 1.00000
```

InS: AB_oP8_58_g_g - POSCAR

```
AB_oP8_58_g_g & a,b/a,c/a,x1,y1,x2,y2 --params=4.443,2.39522844925,
      ↪ 0.886788206167,0.125,0.121,-0.005,0.355 & Pnnm D_{2h}^{12} #58
      ↪ (g^2) & oP8 & None & InS & InS & K. Schubert and E. D\"{o}rre
      ↪ and E. G\"{u}nzel, Naturwissenschaften 41, 448(1954)
   1.00000000000000
   4.44300000000000    0.00000000000000    0.00000000000000
   0.00000000000000   10.64200000000000    0.00000000000000
   0.00000000000000    0.00000000000000    3.94000000000000
     In   S
      4   4
Direct
   0.12500000000000    0.12100000000000    0.00000000000000  In   (4g)
  -0.12500000000000   -0.12100000000000    0.00000000000000  In   (4g)
   0.37500000000000    0.62100000000000    0.50000000000000  In   (4g)
   0.62500000000000    0.37900000000000    0.50000000000000  In   (4g)
  -0.00500000000000    0.35500000000000    0.00000000000000  S    (4g)
   0.00500000000000   -0.35500000000000    0.00000000000000  S    (4g)
   0.50500000000000    0.85500000000000    0.50000000000000  S    (4g)
   0.49500000000000    0.14500000000000    0.50000000000000  S    (4g)
```

RuB_2: A2B_oP6_59_f_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'B2Ru'
_chemical_formula_sum 'B2 Ru'

loop_
_publ_author_name
 'B. Aronsson'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 17
_journal_year 1963
_journal_page_first 2036
_journal_page_last 2050
_publ_Section_title
;
 The Crystal Structure of RuB$_{2}$, OsB$_{2}$, and IrB$_{1.35}$ and
      ↪ Some General Comments on the Crystal Chemistry of Borides in
      ↪ the Composition Range MeB - MeB$_{3}$
;

_aflow_title 'RuB$_{2}$ Structure'
_aflow_proto 'A2B_oP6_59_f_a'
_aflow_params 'a,b/a,c/a,z_{1},x_{2},z_{2}'
_aflow_params_values '4.645,0.616792249731,0.870828848224,0.1508,0.059,
      ↪ 0.639'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP6'

_symmetry_space_group_name_H-M "P 21/m 21/m 2/n (origin choice 2)"
_symmetry_Int_Tables_number 59

_cell_length_a    4.64500
_cell_length_b    2.86500
_cell_length_c    4.04500
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y,-z
3 -x,y+1/2,-z
4 -x+1/2,-y+1/2,z
5 -x,-y,-z
6 -x+1/2,y,z
7 x,-y+1/2,z
8 x+1/2,y+1/2,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ru1 Ru 2 a 0.25000 0.25000 0.15080 1.00000
B1 B 4 f 0.05900 0.25000 0.63900 1.00000
```

RuB_2: A2B_oP6_59_f_a - POSCAR

```
A2B_oP6_59_f_a & a,b/a,c/a,z1,x2,z2 --params=4.645,0.616792249731,
      ↪ 0.870828848224,0.1508,0.059,0.639 & Pmmn D_{2h}^{13} #59 (af) &
      ↪ oP6 & None & B2Ru & B2Ru & B. Aronsson, Acta Chem. Scand. 17,
      ↪ 2036-2050 (1963)
   1.00000000000000
   4.64500000000000    0.00000000000000    0.00000000000000
   0.00000000000000    2.86500000000000    0.00000000000000
   0.00000000000000    0.00000000000000    4.04500000000000
     B     Ru
     4      2
Direct
   0.05900000000000    0.25000000000000    0.63900000000000     B    (4f)
   0.44100000000000    0.25000000000000    0.63900000000000     B    (4f)
  -0.05900000000000    0.75000000000000   -0.63900000000000     B    (4f)
   0.55900000000000    0.75000000000000   -0.63900000000000     B    (4f)
   0.25000000000000    0.25000000000000    0.15080000000000    Ru    (2a)
   0.75000000000000    0.75000000000000   -0.15080000000000    Ru    (2a)
```

NH_4NO_3 IV ($G0_{11}$): A4B2C3_oP18_59_ef_ab_af - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'H4N2O3'
_chemical_formula_sum 'H4 N2 O3'

loop_
_publ_author_name
 'C. S. Choi'
 'J. E. Mapes'
 'E. Prince'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 28
_journal_year 1972
_journal_page_first 1357
_journal_page_last 1361
_publ_Section_title
;
 The structure of ammonium nitrate (IV)
;

_aflow_title 'NH$_{4}$NO$_{3}$ IV ($G0_{11}$) Structure'
_aflow_proto 'A4B2C3_oP18_59_ef_ab_af'
_aflow_params 'a,b/a,c/a,z_{1},z_{2},z_{3},y_{4},z_{4},x_{5},z_{5},x_{6
      ↪ },z_{6}'
_aflow_params_values '5.745,0.946562228024,0.860226283725,0.5067,0.7629,
      ↪ 0.0836,0.8989,-0.0324,0.6045,0.8102,0.43442,0.3832'
_aflow_Strukturbericht '$G0_{11}$'
_aflow_Pearson 'oP18'

_symmetry_space_group_name_H-M "P 21/m 21/m 2/n (origin choice 2)"
_symmetry_Int_Tables_number 59

_cell_length_a    5.74500
_cell_length_b    5.43800
_cell_length_c    4.94200
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y,-z
3 -x,y+1/2,-z
4 -x+1/2,-y+1/2,z
5 -x,-y,-z
6 -x+1/2,y,z
7 x,-y+1/2,z
8 x+1/2,y+1/2,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
N1 N 2 a 0.25000 0.25000 0.50670 1.00000
O1 O 2 a 0.25000 0.25000 0.76290 1.00000
N2 N 2 b 0.25000 0.75000 0.08360 1.00000
H1 H 4 e 0.25000 0.89890 -0.03240 1.00000
H2 H 4 f 0.60450 0.25000 0.81020 1.00000
O2 O 4 f 0.43442 0.25000 0.38320 1.00000
```

NH_4NO_3 IV ($G0_{11}$): A4B2C3_oP18_59_ef_ab_af - POSCAR

```
A4B2C3_oP18_59_ef_ab_af & a,b/a,c/a,z1,z2,z3,y4,z4,x5,z5,x6,z6 --params=
      ↪ 5.745,0.946562228024,0.860226283725,0.5067,0.7629,0.0836,0.8989
      ↪ ,-0.0324,0.6045,0.8102,0.43442,0.3832 & Pmmn D_{2h}^{13} #59 (a
      ↪ ^2bef^2) & oP18 & $G0_{11}$ & H4N2O3 & H4N2O3 & C. S. Choi and
      ↪ J. E. Mapes and E. Prince, Acta Crystallogr. Sect. B Struct.
      ↪ Sci. 28, 1357-1361 (1972)
   1.00000000000000
   5.74500000000000    0.00000000000000    0.00000000000000
   0.00000000000000    5.43800000000000    0.00000000000000
   0.00000000000000    0.00000000000000    4.94200000000000
     H     N     O
     8     4     6
Direct
   0.25000000000000    0.89890000000000   -0.03240000000000     H    (4e)
   0.25000000000000   -0.39890000000000   -0.03240000000000     H    (4e)
   0.75000000000000    1.39890000000000    0.03240000000000     H    (4e)
   0.75000000000000   -0.89890000000000    0.03240000000000     H    (4e)
   0.60450000000000    0.25000000000000    0.81020000000000     H    (4f)
  -0.10450000000000    0.25000000000000    0.81020000000000     H    (4f)
  -0.60450000000000    0.75000000000000   -0.81020000000000     H    (4f)
   1.10450000000000    0.75000000000000   -0.81020000000000     H    (4f)
   0.25000000000000    0.25000000000000    0.50670000000000     N    (2a)
   0.75000000000000    0.75000000000000   -0.50670000000000     N    (2a)
   0.25000000000000    0.75000000000000    0.08360000000000     N    (2b)
   0.75000000000000    0.25000000000000   -0.08360000000000     N    (2b)
   0.25000000000000    0.25000000000000    0.76290000000000     O    (2a)
   0.75000000000000    0.75000000000000   -0.76290000000000     O    (2a)
   0.43442000000000    0.25000000000000    0.38320000000000     O    (4f)
   0.06558000000000    0.25000000000000    0.38320000000000     O    (4f)
  -0.43442000000000    0.75000000000000   -0.38320000000000     O    (4f)
   0.93442000000000    0.75000000000000   -0.38320000000000     O    (4f)
```

Shcherbinaite (V_2O_5) (*revised*): A5B2_oP14_59_a2f_f - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Shcherbinaite'
_chemical_formula_sum 'O5 V2'

loop_
_publ_author_name
 'R. Enjalbert'
 'J. Galy'
```

```
_journal_name_full_name
;
 Acta Crystallographica Section C: Structural Chemistry
;
_journal_volume 42
_journal_year 1986
_journal_page_first 1467
_journal_page_last 1469
_publ_Section_title
;
 A Refinement of the Structure of V$_{2}$O$_{5}$
;

_aflow_title 'Shcherbinaite (V$_{2}$O$_{5}$) ({\em{revised}}) Structure'
_aflow_proto 'A5B2_oP14_59_a2f_f'
_aflow_params 'a,b/a,c/a,z_{1},x_{2},z_{2},x_{3},z_{3},x_{4},z_{4}'
_aflow_params_values '11.512,0.309589993051,0.379430159833,0.001,0.1043,
      ↪ 0.531,-0.0689,0.003,0.10118,0.891'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP14'

_symmetry_space_group_name_H-M "P 21/m 21/m 2/n (origin choice 2)"
_symmetry_Int_Tables_number 59

_cell_length_a    11.51200
_cell_length_b    3.56400
_cell_length_c    4.36800
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y,-z
3 -x,y+1/2,-z
4 -x+1/2,-y+1/2,z
5 -x,-y,-z
6 -x+1/2,y,z
7 x,-y+1/2,z
8 x+1/2,y+1/2,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 2 a 0.25000 0.25000 0.00100 1.00000
O2 O 4 f 0.10430 0.25000 0.53100 1.00000
O3 O 4 f -0.06890 0.25000 0.00300 1.00000
V1 V 4 f 0.10118 0.25000 0.89100 1.00000
```

Shcherbinaite (V_2O_5) (*revised*): A5B2_oP14_59_a2f_f - POSCAR

```
A5B2_oP14_59_a2f_f & a,b/a,c/a,z1,x2,z2,x3,z3,x4,z4 --params=11.512,
      ↪ 0.309589993051,0.379430159833,0.001,0.1043,0.531,-0.0689,0.003,
      ↪ 0.10118,0.891 & Pmmn D_{2h}^{13} #59 (af^3) & oP14 & None &
      ↪ O5V2 & Shcherbinaite & R. Enjalbert and J. Galy, Acta
      ↪ Crystallogr. C 42, 1467-1469 (1986)
   1.00000000000000
  11.51200000000000    0.00000000000000    0.00000000000000
   0.00000000000000    3.56400000000000    0.00000000000000
   0.00000000000000    0.00000000000000    4.36800000000000
     O     V
    10     4
Direct
   0.25000000000000    0.25000000000000    0.00100000000000  O  (2a)
   0.75000000000000    0.75000000000000   -0.00100000000000  O  (2a)
   0.10430000000000    0.25000000000000    0.53100000000000  O  (4f)
   0.39570000000000    0.25000000000000    0.53100000000000  O  (4f)
  -0.10430000000000    0.75000000000000   -0.53100000000000  O  (4f)
   0.60430000000000    0.75000000000000   -0.53100000000000  O  (4f)
  -0.06890000000000    0.25000000000000    0.00300000000000  O  (4f)
   0.56890000000000    0.25000000000000    0.00300000000000  O  (4f)
   0.06890000000000    0.75000000000000   -0.00300000000000  O  (4f)
   0.43110000000000    0.75000000000000   -0.00300000000000  O  (4f)
   0.10118000000000    0.25000000000000    0.89100000000000  V  (4f)
   0.39882000000000    0.25000000000000    0.89100000000000  V  (4f)
  -0.10118000000000    0.75000000000000   -0.89100000000000  V  (4f)
   0.60118000000000    0.75000000000000   -0.89100000000000  V  (4f)
```

CaB_2O_4 I ($E3_2$): A2BC4_oP28_60_d_c_2d - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'B2CaO4'
_chemical_formula_sum 'B2 Ca O4'

loop_
_publ_author_name
 'M. Marezio'
 'H. A. Plettinger'
 'W. H. Zachariasen'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 16
_journal_year 1963
_journal_page_first 390
_journal_page_last 392
_publ_Section_title
;
 Refinement of the calcium metaborate structure
;

# Found in The crystal structure of the high-pressure phase CaB$_{2}$O$_
      ↪ {4}$(IV), and polymorphism in CaB$_{2}$O$_{4}$, 1969

_aflow_title 'CaB$_{2}$O$_{4}$ I ($E3_{2}$) Structure'
_aflow_proto 'A2BC4_oP28_60_d_c_2d'
_aflow_params 'a,b/a,c/a,y_{1},x_{2},y_{2},z_{2},x_{3},y_{3},z_{3},x_{4
      ↪ },y_{4},z_{4}'
_aflow_params_values '11.604,0.369269217511,0.535504998276,0.2726,0.1924
      ↪ ,0.8296,0.1258,0.0862,0.7268,0.0917,0.2078,0.1485,0.1478'
_aflow_Strukturbericht '$E3_{2}$'
_aflow_Pearson 'oP28'

_symmetry_space_group_name_H-M "P 21/b 2/c 21/n"
_symmetry_Int_Tables_number 60

_cell_length_a    11.60400
_cell_length_b    4.28500
_cell_length_c    6.21400
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z
3 -x,y,-z+1/2
4 -x+1/2,-y+1/2,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z
7 x,-y,z+1/2
8 x+1/2,y+1/2,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ca1 Ca 4 c 0.00000 0.27260 0.25000 1.00000
B1 B 8 d 0.19240 0.82960 0.12580 1.00000
O1 O 8 d 0.08620 0.72680 0.09170 1.00000
O2 O 8 d 0.20780 0.14850 0.14780 1.00000
```

CaB_2O_4 I ($E3_2$): A2BC4_oP28_60_d_c_2d - POSCAR

```
A2BC4_oP28_60_d_c_2d & a,b/a,c/a,y1,x2,y2,z2,x3,y3,z3,x4,y4,z4 --params=
      ↪ 11.604,0.369269217511,0.535504998276,0.2726,0.1924,0.8296,
      ↪ 0.1258,0.0862,0.7268,0.0917,0.2078,0.1485,0.1478 & Pbcn D_{2h
      ↪ }^{14} #60 (cd^3) & oP28 & $E3_{2}$ & B2CaO4 & B2CaO4 & M.
      ↪ Marezio and H. A. Plettinger and W. H. Zachariasen, Acta Cryst.
      ↪ 16, 390-392 (1963)
   1.00000000000000
  11.60400000000000    0.00000000000000    0.00000000000000
   0.00000000000000    4.28500000000000    0.00000000000000
   0.00000000000000    0.00000000000000    6.21400000000000
     B     Ca     O
     8     4    16
Direct
   0.19240000000000    0.82960000000000    0.12580000000000  B   (8d)
   0.30760000000000   -0.32960000000000    0.62580000000000  B   (8d)
  -0.19240000000000    0.82960000000000    0.37420000000000  B   (8d)
   0.69240000000000   -0.32960000000000   -0.12580000000000  B   (8d)
  -0.19240000000000   -0.82960000000000   -0.12580000000000  B   (8d)
   0.69240000000000    1.32960000000000    0.37420000000000  B   (8d)
   0.19240000000000   -0.82960000000000    0.62580000000000  B   (8d)
   0.30760000000000    1.32960000000000    0.12580000000000  B   (8d)
   0.00000000000000    0.27260000000000    0.25000000000000  Ca  (4c)
   0.50000000000000    0.22740000000000    0.75000000000000  Ca  (4c)
   0.00000000000000   -0.27260000000000    0.75000000000000  Ca  (4c)
   0.50000000000000    0.77260000000000    0.25000000000000  Ca  (4c)
   0.08620000000000    0.72680000000000    0.09170000000000  O   (8d)
   0.41380000000000   -0.22680000000000    0.59170000000000  O   (8d)
  -0.08620000000000    0.72680000000000    0.40830000000000  O   (8d)
   0.58620000000000   -0.22680000000000   -0.09170000000000  O   (8d)
  -0.08620000000000   -0.72680000000000   -0.09170000000000  O   (8d)
   0.58620000000000    1.22680000000000    0.40830000000000  O   (8d)
   0.08620000000000   -0.72680000000000    0.59170000000000  O   (8d)
   0.41380000000000    1.22680000000000    0.09170000000000  O   (8d)
   0.20780000000000    0.14850000000000    0.14780000000000  O   (8d)
   0.29220000000000    0.35150000000000    0.64780000000000  O   (8d)
  -0.20780000000000    0.14850000000000    0.35220000000000  O   (8d)
   0.70780000000000    0.35150000000000   -0.14780000000000  O   (8d)
  -0.20780000000000   -0.14850000000000   -0.14780000000000  O   (8d)
   0.70780000000000    0.64850000000000    0.35220000000000  O   (8d)
   0.20780000000000   -0.14850000000000    0.64780000000000  O   (8d)
   0.29220000000000    0.64850000000000    0.14780000000000  O   (8d)
```

ζ-Fe_2N: A2B_oP12_60_d_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM
```

```
_chemical_name_mineral 'Fe2N'
_chemical_formula_sum 'Fe2 N'

loop_
_publ_author_name
 'D. Rechenbach'
 'H. Jacobs'
_journal_name_full_name
;
 Journal of Alloys and Compounds
;
_journal_volume 235
_journal_year 1996
_journal_page_first 15
_journal_page_last 22
_publ_Section_title
;
 Structure determination of $\zeta$-Fe$_{2}$N by neutron and synchrotron
      ↪ powder diffraction
;

_aflow_title '$\zeta$-Fe$_{2}$N Structure'
_aflow_proto 'A2B_oP12_60_d_c'
_aflow_params 'a,b/a,c/a,y_{1},x_{2},y_{2},z_{2}'
_aflow_params_values '4.4373,1.24879994591,1.09140693665,0.364,0.249,
      ↪ 0.128,0.0827'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP12'

_symmetry_space_group_name_H-M "P 21/b 2/c 21/n"
_symmetry_Int_Tables_number 60

_cell_length_a     4.43730
_cell_length_b     5.54130
_cell_length_c     4.84290
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z
3 -x,y,-z+1/2
4 -x+1/2,-y+1/2,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z
7 x,-y,z+1/2
8 x+1/2,y+1/2,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
N1 N 4 c 0.00000 0.36400 0.25000 1.00000
Fe1 Fe 8 d 0.24900 0.12800 0.08270 1.00000
```

ζ-Fe_2N: A2B_oP12_60_d_c - POSCAR

```
A2B_oP12_60_d_c & a,b/a,c/a,y1,x2,y2,z2 --params=4.4373,1.24879994591,
      ↪ 1.09140693665,0.364,0.249,0.128,0.0827 & Pbcn D_{2h}^{14} #60 (
      ↪ cd) & oP12 & None & Fe2N & Fe2N & D. Rechenbach and H. Jacobs,
      ↪ J. Alloys Compd. 235, 15-22 (1996)
   1.00000000000000
   4.43730000000000    0.00000000000000    0.00000000000000
   0.00000000000000    5.54130000000000    0.00000000000000
   0.00000000000000    0.00000000000000    4.84290000000000
     Fe      N
      8      4
Direct
   0.24900000000000    0.12800000000000    0.08270000000000   Fe   (8d)
   0.25100000000000    0.37200000000000    0.58270000000000   Fe   (8d)
  -0.24900000000000    0.12800000000000    0.41730000000000   Fe   (8d)
   0.74900000000000    0.37200000000000   -0.08270000000000   Fe   (8d)
  -0.24900000000000   -0.12800000000000   -0.08270000000000   Fe   (8d)
   0.74900000000000    0.62800000000000    0.41730000000000   Fe   (8d)
   0.24900000000000   -0.12800000000000    0.58270000000000   Fe   (8d)
   0.25100000000000    0.62800000000000    0.08270000000000   Fe   (8d)
   0.00000000000000    0.36400000000000    0.25000000000000    N   (4c)
   0.50000000000000    0.13600000000000    0.75000000000000    N   (4c)
   0.00000000000000   -0.36400000000000    0.75000000000000    N   (4c)
   0.50000000000000    0.86400000000000    0.25000000000000    N   (4c)
```

α-PbO_2: A2B_oP12_60_d_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'O2Pb'
_chemical_formula_sum 'O2 Pb'

loop_
_publ_author_name
 'R. J. Hill'
_journal_name_full_name
;
 Materials Research Bulletin
;
_journal_volume 17
_journal_year 1982
_journal_page_first 769
_journal_page_last 784
_publ_Section_title
;
 The Crystal Structures of Lead Dioxides from the Positive Plate of the
      ↪ Lead/Acid Battery
;

# Found in Pearson's Handbook of Crystallographic Data for Intermetallic
      ↪ Phases, 1991

_aflow_title '$\alpha$-PbO$_{2}$ Structure'
_aflow_proto 'A2B_oP12_60_d_c'
_aflow_params 'a,b/a,c/a,y_{1},x_{2},y_{2},z_{2}'
_aflow_params_values '4.9898,1.19191149946,1.09535452323,0.1779,0.2685,
      ↪ 0.401,0.0248'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP12'

_symmetry_space_group_name_H-M "P 21/b 2/c 21/n"
_symmetry_Int_Tables_number 60

_cell_length_a     4.98980
_cell_length_b     5.94740
_cell_length_c     5.46560
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z
3 -x,y,-z+1/2
4 -x+1/2,-y+1/2,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z
7 x,-y,z+1/2
8 x+1/2,y+1/2,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Pb1 Pb 4 c 0.00000 0.17790 0.25000 0.49000
O1 O 8 d 0.26850 0.40100 0.02480 1.00000
```

α-PbO_2: A2B_oP12_60_d_c - POSCAR

```
A2B_oP12_60_d_c & a,b/a,c/a,y1,x2,y2,z2 --params=4.9898,1.19191149946,
      ↪ 1.09535452323,0.1779,0.2685,0.401,0.0248 & Pbcn D_{2h}^{14} #60
      ↪ (cd) & oP12 & None & O2Pb & O2Pb & R. J. Hill, Mater. Res.
      ↪ Bull. 17, 769-784 (1982)
   1.00000000000000
   4.98980000000000    0.00000000000000    0.00000000000000
   0.00000000000000    5.94740000000000    0.00000000000000
   0.00000000000000    0.00000000000000    5.46560000000000
     O     Pb
     8      4
Direct
   0.26850000000000    0.40100000000000    0.02480000000000    O   (8d)
   0.23150000000000    0.09900000000000    0.52480000000000    O   (8d)
  -0.26850000000000    0.40100000000000    0.47520000000000    O   (8d)
   0.76850000000000    0.09900000000000   -0.02480000000000    O   (8d)
  -0.26850000000000   -0.40100000000000   -0.02480000000000    O   (8d)
   0.76850000000000    0.90100000000000    0.47520000000000    O   (8d)
   0.26850000000000   -0.40100000000000    0.52480000000000    O   (8d)
   0.23150000000000    0.90100000000000    0.02480000000000    O   (8d)
   0.00000000000000    0.17790000000000    0.25000000000000   Pb   (4c)
   0.50000000000000    0.32210000000000    0.75000000000000   Pb   (4c)
   0.00000000000000   -0.17790000000000    0.75000000000000   Pb   (4c)
   0.50000000000000    0.67790000000000    0.25000000000000   Pb   (4c)
```

Cr_5O_{12}: A5B12_oP68_60_c2d_6d - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Cr5O12'
_chemical_formula_sum 'Cr5 O12'

loop_
_publ_author_name
 'K.-A. Wilhelmi'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 19
_journal_year 1965
_journal_page_first 165
_journal_page_last 176
_publ_Section_title
;
 The Crystal Structure of Cr$_{5}$O$_{12}$
```

```
;

_aflow_title 'Cr$_{5}$O$_{12}$ Structure'
_aflow_proto 'A5B12_oP68_60_c2d_6d'
_aflow_params 'a,b/a,c/a,y_{1},x_{2},y_{2},z_{2},x_{3},y_{3},z_{3},x_{4
    ↪ },y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{6},x_{7},y_{7},
    ↪ z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9}'
_aflow_params_values '12.044,0.681833277981,0.678927266689,0.3939,0.1719
    ↪ ,0.8878,0.4768,0.0857,0.7517,0.1119,0.2454,0.7314,0.0906,0.075,
    ↪ 0.7736,0.3535,0.1003,-0.0101,0.1032,0.0806,0.508,0.1356,0.0707,
    ↪ 0.27,0.3741,0.2538,0.4894,0.3625'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP68'

_symmetry_space_group_name_H-M "P 21/b 2/c 21/n"
_symmetry_Int_Tables_number 60

_cell_length_a 12.04400
_cell_length_b 8.21200
_cell_length_c 8.17700
_cell_angle_alpha 90.00000
_cell_angle_beta 90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z
3 -x,y,-z+1/2
4 -x+1/2,-y+1/2,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z
7 x,-y,z+1/2
8 x+1/2,y+1/2,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cr1 Cr 4 c 0.00000 0.39390 0.25000 1.00000
Cr2 Cr 8 d 0.17190 0.88780 0.47680 1.00000
Cr3 Cr 8 d 0.08570 0.75170 0.11190 1.00000
O1 O 8 d 0.24540 0.73140 0.09060 1.00000
O2 O 8 d 0.07500 0.77360 0.35350 1.00000
O3 O 8 d 0.10030 -0.01010 0.10320 1.00000
O4 O 8 d 0.08060 0.50800 0.13560 1.00000
O5 O 8 d 0.07070 0.27000 0.37410 1.00000
O6 O 8 d 0.25380 0.48940 0.36250 1.00000
```

Cr_5O_{12}: A5B12_oP68_60_c2d_6d - POSCAR

```
A5B12_oP68_60_c2d_6d & a,b/a,c/a,y1,x2,y2,z2,x3,y3,z3,x4,y4,z4,x5,y5,z5,
    ↪ x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9 --params=12.044,
    ↪ 0.681833277981,0.678927266689,0.3939,0.1719,0.8878,0.4768,
    ↪ 0.0857,0.7517,0.1119,0.2454,0.7314,0.0906,0.075,0.7736,0.3535,
    ↪ 0.1003,-0.0101,0.1032,0.0806,0.508,0.1356,0.0707,0.27,0.3741,
    ↪ 0.2538,0.4894,0.3625 & Pbcn D_{2h}^{14} #60 (cd^8) & oP68 &
    ↪ None & Cr5O12 & Cr5O12 & K.-A. Wilhelmi, Acta Chem. Scand. 19,
    ↪ 165-176 (1965)
   1.00000000000000
  12.04400000000000    0.00000000000000    0.00000000000000
   0.00000000000000    8.21200000000000    0.00000000000000
   0.00000000000000    0.00000000000000    8.17700000000000
     Cr      O
     20     48
Direct
   0.00000000000000    0.39390000000000    0.25000000000000   Cr   (4c)
   0.50000000000000    0.10610000000000    0.75000000000000   Cr   (4c)
   0.00000000000000   -0.39390000000000    0.75000000000000   Cr   (4c)
   0.50000000000000    0.89390000000000    0.25000000000000   Cr   (4c)
   0.17190000000000    0.88780000000000    0.47680000000000   Cr   (8d)
   0.32810000000000   -0.38780000000000    0.97680000000000   Cr   (8d)
  -0.17190000000000    0.88780000000000    0.02320000000000   Cr   (8d)
   0.67190000000000   -0.38780000000000   -0.47680000000000   Cr   (8d)
  -0.17190000000000   -0.88780000000000   -0.47680000000000   Cr   (8d)
   0.67190000000000    1.38780000000000    0.02320000000000   Cr   (8d)
   0.17190000000000   -0.88780000000000    0.97680000000000   Cr   (8d)
   0.32810000000000    1.38780000000000    0.47680000000000   Cr   (8d)
   0.08570000000000    0.75170000000000    0.11190000000000   Cr   (8d)
   0.41430000000000   -0.25170000000000    0.61190000000000   Cr   (8d)
  -0.08570000000000    0.75170000000000    0.38810000000000   Cr   (8d)
   0.58570000000000   -0.25170000000000   -0.11190000000000   Cr   (8d)
  -0.08570000000000   -0.75170000000000   -0.11190000000000   Cr   (8d)
   0.58570000000000    1.25170000000000    0.38810000000000   Cr   (8d)
   0.08570000000000   -0.75170000000000    0.61190000000000   Cr   (8d)
   0.41430000000000    1.25170000000000    0.11190000000000   Cr   (8d)
   0.24540000000000    0.73140000000000    0.09060000000000   O    (8d)
   0.25460000000000   -0.23140000000000    0.59060000000000   O    (8d)
  -0.24540000000000    0.73140000000000    0.40940000000000   O    (8d)
   0.74540000000000   -0.23140000000000   -0.09060000000000   O    (8d)
  -0.24540000000000   -0.73140000000000   -0.09060000000000   O    (8d)
   0.74540000000000    1.23140000000000    0.40940000000000   O    (8d)
   0.24540000000000   -0.73140000000000    0.59060000000000   O    (8d)
   0.25460000000000    1.23140000000000    0.09060000000000   O    (8d)
   0.07500000000000    0.77360000000000    0.35350000000000   O    (8d)
   0.42500000000000   -0.27360000000000    0.85350000000000   O    (8d)
  -0.07500000000000    0.77360000000000    0.14650000000000   O    (8d)
   0.57500000000000   -0.27360000000000   -0.35350000000000   O    (8d)
  -0.07500000000000   -0.77360000000000   -0.35350000000000   O    (8d)
   0.57500000000000    1.27360000000000    0.14650000000000   O    (8d)
   0.07500000000000   -0.77360000000000    0.85350000000000   O    (8d)
   0.42500000000000    1.27360000000000    0.35350000000000   O    (8d)
   0.10030000000000   -0.01010000000000    0.10320000000000   O    (8d)
   0.39970000000000    0.51010000000000    0.60320000000000   O    (8d)
  -0.10030000000000   -0.01010000000000    0.39680000000000   O    (8d)
   0.60030000000000    0.51010000000000   -0.10320000000000   O    (8d)
  -0.10030000000000    0.01010000000000   -0.10320000000000   O    (8d)
   0.60030000000000    0.48990000000000    0.39680000000000   O    (8d)
   0.10030000000000    0.01010000000000    0.60320000000000   O    (8d)
   0.39970000000000    0.48990000000000    0.10320000000000   O    (8d)
   0.08060000000000    0.50800000000000    0.13560000000000   O    (8d)
   0.41940000000000   -0.00800000000000    0.63560000000000   O    (8d)
  -0.08060000000000    0.50800000000000    0.36440000000000   O    (8d)
   0.58060000000000   -0.00800000000000   -0.13560000000000   O    (8d)
  -0.08060000000000   -0.50800000000000   -0.13560000000000   O    (8d)
   0.58060000000000    1.00800000000000    0.36440000000000   O    (8d)
   0.08060000000000   -0.50800000000000    0.63560000000000   O    (8d)
   0.41940000000000    1.00800000000000    0.13560000000000   O    (8d)
   0.07070000000000    0.27000000000000    0.37410000000000   O    (8d)
   0.42930000000000    0.23000000000000    0.87410000000000   O    (8d)
  -0.07070000000000    0.27000000000000    0.12590000000000   O    (8d)
   0.57070000000000    0.23000000000000   -0.37410000000000   O    (8d)
  -0.07070000000000   -0.27000000000000   -0.37410000000000   O    (8d)
   0.57070000000000    0.77000000000000    0.12590000000000   O    (8d)
   0.07070000000000   -0.27000000000000    0.87410000000000   O    (8d)
   0.42930000000000    0.77000000000000    0.37410000000000   O    (8d)
   0.25380000000000    0.48940000000000    0.36250000000000   O    (8d)
   0.24620000000000    0.01060000000000    0.86250000000000   O    (8d)
  -0.25380000000000    0.48940000000000    0.13750000000000   O    (8d)
   0.75380000000000    0.01060000000000   -0.36250000000000   O    (8d)
  -0.25380000000000   -0.48940000000000   -0.36250000000000   O    (8d)
   0.75380000000000    0.98940000000000    0.13750000000000   O    (8d)
   0.25380000000000   -0.48940000000000    0.86250000000000   O    (8d)
   0.24620000000000    0.98940000000000    0.36250000000000   O    (8d)
```

Columbite ($FeNb_2O_6$, $E5_1$): AB2C6_oP36_60_c_d_3d - CIF

```
# CIF file
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_audit_creation_method FINDSYM

_chemical_name_mineral 'Columbite'
_chemical_formula_sum 'Fe Nb2 O6'

loop_
_publ_author_name
 'P. Bordet'
 'A. {McHale}'
 'A. Santoro'
 'R. S. Roth'
_journal_name_full_name
;
 Journal of Solid State Chemistry
;
_journal_volume 64
_journal_year 1986
_journal_page_first 30
_journal_page_last 46
_publ_Section_title
;
 Powder neutron diffraction study of ZrTiO$_{4}$, Zr$_{5}$Ti$_{7}$O$_{24
    ↪ }$, and FeNb$_{2}$O$_{6}$
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'Columbite (FeNb$_{2}$O$_{6}$, $E5_{1}$) Structure'
_aflow_proto 'AB2C6_oP36_60_c_d_3d'
_aflow_params 'a,b/a,c/a,y_{1},x_{2},y_{2},z_{2},x_{3},y_{3},z_{3},x_{4
    ↪ },y_{4},z_{4},x_{5},y_{5},z_{5}'
_aflow_params_values '14.2661,0.401889794688,0.353950974688,0.3311,
    ↪ 0.3389,0.3191,0.2506,0.0963,0.1041,0.0727,0.4189,0.1163,0.099,
    ↪ 0.756,0.1236,0.0793'
_aflow_Strukturbericht '$E5_{1}$'
_aflow_Pearson 'oP36'

_symmetry_space_group_name_H-M "P 21/b 2/c 21/n"
_symmetry_Int_Tables_number 60

_cell_length_a 14.26610
_cell_length_b 5.73340
_cell_length_c 5.04950
_cell_angle_alpha 90.00000
_cell_angle_beta 90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z
3 -x,y,-z+1/2
4 -x+1/2,-y+1/2,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z
7 x,-y,z+1/2
8 x+1/2,y+1/2,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
```

```
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Fe1 Fe 4 c 0.00000 0.33110 0.25000 1.00000
Nb1 Nb 8 d 0.33890 0.31910 0.25060 1.00000
O1 O 8 d 0.09630 0.10410 0.07270 1.00000
O2 O 8 d 0.41890 0.11630 0.09900 1.00000
O3 O 8 d 0.75600 0.12360 0.07930 1.00000
```

Columbite ($FeNb_2O_6$, $E5_1$): AB2C6_oP36_60_c_d_3d - POSCAR

```
AB2C6_oP36_60_c_d_3d & a,b/a,c/a,y1,x2,y2,z2,x3,y3,z3,x4,y4,z4,x5,y5,z5
    ↪ --params=14.2661,0.401889794688,0.353950974688,0.3311,0.3389,
    ↪ 0.3191,0.2506,0.0963,0.1041,0.0727,0.4189,0.1163,0.099,0.756,
    ↪ 0.1236,0.0793 & Pbcn D_{2h}^{14} #60 (cd^4) & oP36 & $E5_{1}$ &
    ↪  FeNb2O6 & Columbite & P. Bordet et al., J. Solid State Chem.
    ↪ 64, 30-46 (1986)
   1.00000000000000
  14.26610000000000    0.00000000000000    0.00000000000000
   0.00000000000000    5.73340000000000    0.00000000000000
   0.00000000000000    0.00000000000000    5.04950000000000
    Fe    Nb    O
     4     8    24
Direct
   0.00000000000000    0.33110000000000    0.25000000000000  Fe   (4c)
   0.50000000000000    0.16890000000000    0.75000000000000  Fe   (4c)
   0.00000000000000   -0.33110000000000    0.75000000000000  Fe   (4c)
   0.50000000000000    0.83110000000000    0.25000000000000  Fe   (4c)
   0.33890000000000    0.31910000000000    0.25060000000000  Nb   (8d)
   0.16110000000000    0.18090000000000    0.75060000000000  Nb   (8d)
  -0.33890000000000    0.31910000000000    0.24940000000000  Nb   (8d)
   0.83890000000000    0.18090000000000   -0.25060000000000  Nb   (8d)
  -0.33890000000000   -0.31910000000000   -0.25060000000000  Nb   (8d)
   0.83890000000000    0.81910000000000    0.24940000000000  Nb   (8d)
   0.33890000000000   -0.31910000000000    0.75060000000000  Nb   (8d)
   0.16110000000000    0.81910000000000    0.25060000000000  Nb   (8d)
   0.09630000000000    0.10410000000000    0.07270000000000  O    (8d)
   0.40370000000000    0.39590000000000    0.57270000000000  O    (8d)
  -0.09630000000000    0.10410000000000    0.42730000000000  O    (8d)
   0.59630000000000    0.39590000000000   -0.07270000000000  O    (8d)
  -0.09630000000000   -0.10410000000000   -0.07270000000000  O    (8d)
   0.59630000000000    0.60410000000000    0.42730000000000  O    (8d)
   0.09630000000000   -0.10410000000000    0.57270000000000  O    (8d)
   0.40370000000000    0.60410000000000    0.07270000000000  O    (8d)
   0.41890000000000    0.11630000000000    0.09900000000000  O    (8d)
   0.08110000000000    0.38370000000000    0.59900000000000  O    (8d)
  -0.41890000000000    0.11630000000000    0.40100000000000  O    (8d)
   0.91890000000000    0.38370000000000   -0.09900000000000  O    (8d)
  -0.41890000000000   -0.11630000000000   -0.09900000000000  O    (8d)
   0.91890000000000    0.61630000000000    0.40100000000000  O    (8d)
   0.41890000000000   -0.11630000000000    0.59900000000000  O    (8d)
   0.08110000000000    0.61630000000000    0.09900000000000  O    (8d)
   0.75600000000000    0.12360000000000    0.07930000000000  O    (8d)
  -0.25600000000000    0.37640000000000    0.57930000000000  O    (8d)
  -0.75600000000000    0.12360000000000    0.42070000000000  O    (8d)
   1.25600000000000    0.37640000000000   -0.07930000000000  O    (8d)
  -0.75600000000000   -0.12360000000000   -0.07930000000000  O    (8d)
   1.25600000000000    0.62360000000000    0.42070000000000  O    (8d)
   0.75600000000000   -0.12360000000000    0.57930000000000  O    (8d)
  -0.25600000000000    0.62360000000000    0.07930000000000  O    (8d)
```

Ca_2RuO_4: A2B4C_oP28_61_c_2c_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Ca2O4Ru'
_chemical_formula_sum 'Ca2 O4 Ru'

loop_
_publ_author_name
 'O. Friedt'
 'M. Braden'
 'G. Andr\\'{e}'
 'P. Adelmann'
 'S. Nakatsuji'
 'Y. Maeno'
_journal_name_full_name
;
 Physical Review B
;
_journal_volume 63
_journal_year 2001
_journal_page_first 174432
_journal_page_last 174432
_publ_Section_title
;
 Structural and magnetic aspects of the metal-insulator transition in
    ↪ Ca$_{2-x}$Sr$_{x}$RuO$_{4}$
;

# Found in Unique Crystal Structure of Ca$_{2}$RuO$_{4}$ in the Current
    ↪ Stabilized Semimetallic State, 2019

_aflow_title 'Ca$_{2}$RuO$_{4}$ Structure'
_aflow_proto 'A2B4C_oP28_61_c_2c_a'
_aflow_params 'a,b/a,c/a,x_{2},y_{2},z_{2},x_{3},y_{3},z_{3},x_{4},y_{4
    ↪ },z_{4}'
_aflow_params_values '5.3945,1.03807581796,2.18098062842,0.0042,0.0559,
    ↪ 0.3524,0.1961,0.3018,0.0264,0.0673,0.4782,0.3355'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP28'

_symmetry_space_group_name_H-M "P 21/b 21/c 21/a"
_symmetry_Int_Tables_number 61

_cell_length_a    5.39450
_cell_length_b    5.59990
_cell_length_c    11.76530
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z
3 -x,y+1/2,-z+1/2
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z
7 x,-y+1/2,z+1/2
8 x+1/2,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ru1 Ru 4 a 0.00000 0.00000 0.00000 1.00000
Ca1 Ca 8 c 0.00420 0.05590 0.35240 1.00000
O1 O 8 c 0.19610 0.30180 0.02640 1.00000
O2 O 8 c 0.06730 0.47820 0.33550 1.00000
```

Ca_2RuO_4: A2B4C_oP28_61_c_2c_a - POSCAR

```
A2B4C_oP28_61_c_2c_a & a,b/a,c/a,x2,y2,z2,x3,y3,z3,x4,y4,z4 --params=
    ↪ 5.3945,1.03807581796,2.18098062842,0.0042,0.0559,0.3524,0.1961,
    ↪ 0.3018,0.0264,0.0673,0.4782,0.3355 & Pbca D_{2h}^{15} #61 (ac^3
    ↪ ) & oP28 & None & Ca2O4Ru & Ca2O4Ru & O. Friedt et al., Phys.
    ↪ Rev. B 63, 174432(2001)
   1.00000000000000
   5.39450000000000    0.00000000000000    0.00000000000000
   0.00000000000000    5.59990000000000    0.00000000000000
   0.00000000000000    0.00000000000000   11.76530000000000
    Ca    O    Ru
     8    16    4
Direct
   0.00420000000000    0.05590000000000    0.35240000000000  Ca   (8c)
   0.49580000000000   -0.05590000000000    0.85240000000000  Ca   (8c)
  -0.00420000000000    0.55590000000000    0.14760000000000  Ca   (8c)
   0.50420000000000    0.44410000000000   -0.35240000000000  Ca   (8c)
  -0.00420000000000   -0.05590000000000   -0.35240000000000  Ca   (8c)
   0.50420000000000    0.05590000000000    0.14760000000000  Ca   (8c)
   0.00420000000000    0.44410000000000    0.85240000000000  Ca   (8c)
   0.49580000000000    0.55590000000000    0.35240000000000  Ca   (8c)
   0.19610000000000    0.30180000000000    0.02640000000000  O    (8c)
   0.30390000000000   -0.30180000000000    0.52640000000000  O    (8c)
  -0.19610000000000    0.80180000000000    0.47360000000000  O    (8c)
   0.69610000000000    0.19820000000000   -0.02640000000000  O    (8c)
  -0.19610000000000   -0.30180000000000   -0.02640000000000  O    (8c)
   0.69610000000000    0.30180000000000    0.47360000000000  O    (8c)
   0.19610000000000    0.19820000000000    0.52640000000000  O    (8c)
   0.30390000000000    0.80180000000000    0.02640000000000  O    (8c)
   0.06730000000000    0.47820000000000    0.33550000000000  O    (8c)
   0.43270000000000   -0.47820000000000    0.83550000000000  O    (8c)
  -0.06730000000000    0.97820000000000    0.16450000000000  O    (8c)
   0.56730000000000    0.02180000000000   -0.33550000000000  O    (8c)
  -0.06730000000000   -0.47820000000000   -0.33550000000000  O    (8c)
   0.56730000000000    0.47820000000000    0.16450000000000  O    (8c)
   0.06730000000000    0.02180000000000    0.83550000000000  O    (8c)
   0.43270000000000    0.97820000000000    0.33550000000000  O    (8c)
   0.00000000000000    0.00000000000000    0.00000000000000  Ru   (4a)
   0.50000000000000    0.00000000000000    0.50000000000000  Ru   (4a)
   0.00000000000000    0.50000000000000    0.50000000000000  Ru   (4a)
   0.50000000000000    0.50000000000000    0.00000000000000  Ru   (4a)
```

Tellurite (β-TeO_2, $C52$): A2B_oP24_61_2c_c - CIF

```
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_audit_creation_method FINDSYM

_chemical_name_mineral 'Tellurite'
_chemical_formula_sum 'O2 Te'

loop_
_publ_author_name
 'H. Beyer'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 124
_journal_year 1967
_journal_page_first 228
_journal_page_last 237
_publ_Section_title
;
 Verfeinerung der Kristallstruktur von Tellurit, dem rhombischen TeO$_{2
    ↪ }$
;
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```
# Found in {\em Ab initio} study of the vibrational properties of
      ↪ crystalline TeO$_{2}$: The $\alpha$, $\beta$, and $\gamma$
      ↪ phases, 2006

_aflow_title 'Tellurite ($\beta$-TeO$_{2}$, $C52$) Structure'
_aflow_proto 'A2B_oP24_61_2c_c'
_aflow_params 'a,b/a,c/a,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3},y_{3
      ↪ },z_{3}'
_aflow_params_values '15.035,0.363418689724,0.372929830396,0.028,0.634,
      ↪ 0.171,0.168,0.221,0.081,0.1181,0.0252,0.3378'
_aflow_Strukturbericht '$C52$'
_aflow_Pearson 'oP24'

_symmetry_space_group_name_H-M "P 21/b 21/c 21/a"
_symmetry_Int_Tables_number 61

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_cell_length_b     5.46400
_cell_length_c     5.60700
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
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2 x+1/2,-y+1/2,-z
3 -x,y+1/2,-z+1/2
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z
7 x,-y+1/2,z+1/2
8 x+1/2,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 8 c 0.02800 0.63400 0.17100 1.00000
O2 O 8 c 0.16800 0.22100 0.08100 1.00000
Te1 Te 8 c 0.11810 0.02520 0.33780 1.00000
```

Tellurite (β-TeO_2, $C52$): A2B_oP24_61_2c_c - POSCAR

```
A2B_oP24_61_2c_c & a,b/a,c/a,x1,y1,z1,x2,y2,z2,x3,y3,z3 --params=15.035,
      ↪ 0.363418689724,0.372929830396,0.028,0.634,0.171,0.168,0.221,
      ↪ 0.081,0.1181,0.0252,0.3378 & Pbca D_{2h}^{15} #61 (c^3) & oP24
      ↪ & $C52$ & O2Te & Tellurite & H. Beyer, Zeitschrift f{"u}r
      ↪ Kristallographie - Crystalline Materials 124, 228-237 (1967)
   1.00000000000000
  15.03500000000000    0.00000000000000    0.00000000000000
   0.00000000000000    5.46400000000000    0.00000000000000
   0.00000000000000    0.00000000000000    5.60700000000000
     O     Te
    16      8
Direct
   0.02800000000000    0.63400000000000    0.17100000000000     O     (8c)
   0.47200000000000   -0.63400000000000    0.67100000000000     O     (8c)
  -0.02800000000000    1.13400000000000    0.32900000000000     O     (8c)
   0.52800000000000   -0.13400000000000   -0.17100000000000     O     (8c)
  -0.02800000000000   -0.63400000000000   -0.17100000000000     O     (8c)
   0.52800000000000    0.63400000000000    0.32900000000000     O     (8c)
   0.02800000000000   -0.13400000000000    0.67100000000000     O     (8c)
   0.47200000000000    1.13400000000000    0.17100000000000     O     (8c)
   0.16800000000000    0.22100000000000    0.08100000000000     O     (8c)
   0.33200000000000   -0.22100000000000    0.58100000000000     O     (8c)
  -0.16800000000000    0.72100000000000    0.41900000000000     O     (8c)
   0.66800000000000    0.27900000000000   -0.08100000000000     O     (8c)
  -0.16800000000000   -0.22100000000000   -0.08100000000000     O     (8c)
   0.66800000000000    0.22100000000000    0.41900000000000     O     (8c)
   0.16800000000000    0.27900000000000    0.58100000000000     O     (8c)
   0.33200000000000    0.72100000000000    0.08100000000000     O     (8c)
   0.11810000000000    0.02520000000000    0.33780000000000    Te     (8c)
   0.38190000000000   -0.02520000000000    0.83780000000000    Te     (8c)
  -0.11810000000000    0.52520000000000    0.16220000000000    Te     (8c)
   0.61810000000000    0.47480000000000   -0.33780000000000    Te     (8c)
  -0.11810000000000   -0.02520000000000   -0.33780000000000    Te     (8c)
   0.61810000000000    0.02520000000000    0.16220000000000    Te     (8c)
   0.11810000000000    0.47480000000000    0.83780000000000    Te     (8c)
   0.38190000000000    0.52520000000000    0.33780000000000    Te     (8c)
```

$(TiCl_4{\cdot}POCl_3)_2$: A7BCD_oP80_61_7c_c_c_c - CIF

```
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_audit_creation_method FINDSYM

_chemical_name_mineral 'Cl7OPTi'
_chemical_formula_sum 'Cl7 O P Ti'

loop_
_publ_author_name
 'C.-I. Br\"{a}nd\\'{e}n'
 'I. Lindqvist'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 14
_journal_year 1960
_journal_page_first 726
_journal_page_last 732
_publ_Section_title
;
 The Crystal Structure of (TiCl$_{4}$$\cdot$POCl$_{3}$)$_{2}$
;

_aflow_title '(TiCl$_{4}$$\cdot$POCl$_{3}$)$_{2}$ Structure'
_aflow_proto 'A7BCD_oP80_61_7c_c_c_c'
_aflow_params 'a,b/a,c/a,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3},y_{3
      ↪ },z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{6},
      ↪ x_{7},y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{10},y_
      ↪ {10},z_{10}'
_aflow_params_values '12.42,1.02737520129,1.08776167472,0.1062,0.5507,
      ↪ 0.5424,0.2364,0.3205,0.5231,0.0098,0.2107,0.4306,0.0092,0.3285,
      ↪ 0.6506,0.2852,0.5483,0.3103,0.2769,0.3103,0.2522,0.1207,0.4706,
      ↪ 0.1548,0.1075,0.4124,0.3564,0.1875,0.4353,0.2819,0.0629,0.3616,
      ↪ 0.4986'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP80'

_symmetry_space_group_name_H-M "P 21/b 21/c 21/a"
_symmetry_Int_Tables_number 61

_cell_length_a     12.42000
_cell_length_b     12.76000
_cell_length_c     13.51000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z
3 -x,y+1/2,-z+1/2
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z
7 x,-y+1/2,z+1/2
8 x+1/2,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cl1 Cl 8 c 0.10620 0.55070 0.54240 1.00000
Cl2 Cl 8 c 0.23640 0.32050 0.52310 1.00000
Cl3 Cl 8 c 0.00980 0.21070 0.43060 1.00000
Cl4 Cl 8 c 0.00920 0.32850 0.65060 1.00000
Cl5 Cl 8 c 0.28520 0.54830 0.31030 1.00000
Cl6 Cl 8 c 0.27690 0.31030 0.25220 1.00000
Cl7 Cl 8 c 0.12070 0.47060 0.15480 1.00000
O1 O 8 c 0.10750 0.41240 0.35640 1.00000
P1 P 8 c 0.18750 0.43530 0.28190 1.00000
Ti1 Ti 8 c 0.06290 0.36160 0.49860 1.00000
```

$(TiCl_4{\cdot}POCl_3)_2$: A7BCD_oP80_61_7c_c_c_c - POSCAR

```
A7BCD_oP80_61_7c_c_c_c & a,b/a,c/a,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,z4,
      ↪ x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10 --
      ↪ params=12.42,1.02737520129,1.08776167472,0.1062,0.5507,0.5424,
      ↪ 0.2364,0.3205,0.5231,0.0098,0.2107,0.4306,0.0092,0.3285,0.6506,
      ↪ 0.2852,0.5483,0.3103,0.2769,0.3103,0.2522,0.1207,0.4706,0.1548,
      ↪ 0.1075,0.4124,0.3564,0.1875,0.4353,0.2819,0.0629,0.3616,0.4986
      ↪ & Pbca D_{2h}^{15} #61 (c^10) & oP80 & None & Cl7OPTi & Cl7OPTi
      ↪ & C.-I. Br\"{a}nd\'{e}n and I. Lindqvist, Acta Chem. Scand. 14
      ↪ , 726-732 (1960)
   1.00000000000000
  12.42000000000000    0.00000000000000    0.00000000000000
   0.00000000000000   12.76000000000000    0.00000000000000
   0.00000000000000    0.00000000000000   13.51000000000000
     Cl      O      P      Ti
     56      8      8      8
Direct
   0.10620000000000    0.55070000000000    0.54240000000000     Cl     (8c)
   0.39380000000000   -0.55070000000000    1.04240000000000     Cl     (8c)
  -0.10620000000000    1.05070000000000   -0.04240000000000     Cl     (8c)
   0.60620000000000   -0.05070000000000   -0.54240000000000     Cl     (8c)
  -0.10620000000000   -0.55070000000000   -0.54240000000000     Cl     (8c)
   0.60620000000000    0.55070000000000   -0.04240000000000     Cl     (8c)
   0.10620000000000   -0.05070000000000    1.04240000000000     Cl     (8c)
   0.39380000000000    1.05070000000000    0.54240000000000     Cl     (8c)
   0.23640000000000    0.32050000000000    0.52310000000000     Cl     (8c)
   0.26360000000000   -0.32050000000000    1.02310000000000     Cl     (8c)
  -0.23640000000000    0.82050000000000   -0.02310000000000     Cl     (8c)
   0.73640000000000    0.17950000000000   -0.52310000000000     Cl     (8c)
  -0.23640000000000   -0.32050000000000   -0.52310000000000     Cl     (8c)
   0.73640000000000    0.32050000000000   -0.02310000000000     Cl     (8c)
   0.23640000000000    0.17950000000000    1.02310000000000     Cl     (8c)
   0.26360000000000    0.82050000000000    0.52310000000000     Cl     (8c)
   0.00980000000000    0.21070000000000    0.43060000000000     Cl     (8c)
   0.49020000000000   -0.21070000000000    0.93060000000000     Cl     (8c)
  -0.00980000000000    0.71070000000000    0.06940000000000     Cl     (8c)
   0.50980000000000    0.28930000000000   -0.43060000000000     Cl     (8c)
  -0.00980000000000   -0.21070000000000   -0.43060000000000     Cl     (8c)
   0.50980000000000    0.21070000000000    0.06940000000000     Cl     (8c)
   0.00980000000000    0.28930000000000    0.93060000000000     Cl     (8c)
```

```
   0.49020000000000   0.71070000000000   0.43060000000000   Cl   (8c)
   0.00920000000000   0.32850000000000   0.65060000000000   Cl   (8c)
   0.49080000000000  -0.32850000000000   1.15060000000000   Cl   (8c)
  -0.00920000000000   0.82850000000000  -0.15060000000000   Cl   (8c)
   0.50920000000000   0.17150000000000  -0.65060000000000   Cl   (8c)
  -0.00920000000000  -0.32850000000000  -0.65060000000000   Cl   (8c)
   0.50920000000000   0.32850000000000  -0.15060000000000   Cl   (8c)
   0.00920000000000   0.17150000000000   1.15060000000000   Cl   (8c)
   0.49080000000000   0.82850000000000   0.65060000000000   Cl   (8c)
   0.28520000000000   0.54830000000000   0.31030000000000   Cl   (8c)
   0.21480000000000  -0.54830000000000   0.81030000000000   Cl   (8c)
  -0.28520000000000   1.04830000000000   0.18970000000000   Cl   (8c)
   0.78520000000000  -0.04830000000000  -0.31030000000000   Cl   (8c)
  -0.28520000000000  -0.54830000000000  -0.31030000000000   Cl   (8c)
   0.78520000000000   0.54830000000000   0.18970000000000   Cl   (8c)
   0.28520000000000  -0.04830000000000   0.81030000000000   Cl   (8c)
   0.21480000000000   1.04830000000000   0.31030000000000   Cl   (8c)
   0.27690000000000   0.31030000000000   0.25220000000000   Cl   (8c)
   0.22310000000000  -0.31030000000000   0.75220000000000   Cl   (8c)
  -0.27690000000000   0.81030000000000   0.24780000000000   Cl   (8c)
   0.77690000000000   0.18970000000000  -0.25220000000000   Cl   (8c)
  -0.27690000000000  -0.31030000000000  -0.25220000000000   Cl   (8c)
   0.77690000000000   0.31030000000000   0.24780000000000   Cl   (8c)
   0.27690000000000   0.18970000000000   0.75220000000000   Cl   (8c)
   0.22310000000000   0.81030000000000   0.25220000000000   Cl   (8c)
   0.12070000000000   0.47060000000000   0.15480000000000   Cl   (8c)
   0.37930000000000  -0.47060000000000   0.65480000000000   Cl   (8c)
  -0.12070000000000   0.97060000000000   0.34520000000000   Cl   (8c)
   0.62070000000000   0.02940000000000  -0.15480000000000   Cl   (8c)
  -0.12070000000000  -0.47060000000000  -0.15480000000000   Cl   (8c)
   0.62070000000000   0.47060000000000   0.34520000000000   Cl   (8c)
   0.12070000000000   0.02940000000000   0.65480000000000   Cl   (8c)
   0.37930000000000   0.97060000000000   0.15480000000000   Cl   (8c)
   0.10750000000000   0.41240000000000   0.35640000000000   O    (8c)
   0.39250000000000  -0.41240000000000   0.85640000000000   O    (8c)
  -0.10750000000000   0.91240000000000   0.14360000000000   O    (8c)
   0.60750000000000   0.08760000000000  -0.35640000000000   O    (8c)
  -0.10750000000000  -0.41240000000000  -0.35640000000000   O    (8c)
   0.60750000000000   0.41240000000000   0.14360000000000   O    (8c)
   0.10750000000000   0.08760000000000   0.85640000000000   O    (8c)
   0.39250000000000   0.91240000000000   0.35640000000000   O    (8c)
   0.18750000000000   0.43530000000000   0.28190000000000   P    (8c)
   0.31250000000000  -0.43530000000000   0.78190000000000   P    (8c)
  -0.18750000000000   0.93530000000000   0.21810000000000   P    (8c)
   0.68750000000000   0.06470000000000  -0.28190000000000   P    (8c)
  -0.18750000000000  -0.43530000000000  -0.28190000000000   P    (8c)
   0.68750000000000   0.43530000000000   0.21810000000000   P    (8c)
   0.18750000000000   0.06470000000000   0.78190000000000   P    (8c)
   0.31250000000000   0.93530000000000   0.28190000000000   P    (8c)
   0.06290000000000   0.36160000000000   0.49860000000000   Ti   (8c)
   0.43710000000000  -0.36160000000000   0.99860000000000   Ti   (8c)
  -0.06290000000000   0.86160000000000   0.00140000000000   Ti   (8c)
   0.56290000000000   0.13840000000000  -0.49860000000000   Ti   (8c)
  -0.06290000000000  -0.36160000000000  -0.49860000000000   Ti   (8c)
   0.56290000000000   0.36160000000000   0.00140000000000   Ti   (8c)
   0.06290000000000   0.13840000000000   0.99860000000000   Ti   (8c)
   0.43710000000000   0.86160000000000   0.49860000000000   Ti   (8c)
```

Hambergite [$Be_2BO_3(OH)$, $G7_2$]: AB2CD4_oP64_61_c_2c_c_4c - CIF

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_chemical_formula_sum 'B Be2 H O4'

loop_
_publ_author_name
 'G. D. Gatta'
 'G. J. {McIntyre}'
 'G. Bromiley'
 'A. Guastoni'
 'F. Nestola'
_journal_name_full_name
;
 American Mineralogist
;
_journal_volume 97
_journal_year 2012
_journal_page_first 1891
_journal_page_last 1897
_publ_Section_title
;
A single-crystal neutron diffraction study of hambergite, Be$_{2}$BO$_{
    ↪ 3}$(OH,F)
;

_aflow_title 'Hambergite [Be$_{2}$BO$_{3}$(OH), $G7_{2}$] Structure'
_aflow_proto 'AB2CD4_oP64_61_c_2c_c_4c'
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    ↪ },z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{6},
    ↪ x_{7},y_{7},z_{7},x_{8},y_{8},z_{8}'
_aflow_params_values '9.762,1.24984634296,0.453800450727,0.10617,0.60704
    ↪ ,0.77298,0.00261,0.68871,0.26018,0.23724,0.56757,0.27717,0.3138
    ↪ ,0.7228,0.4574,0.0376,0.68766,0.61914,0.1012,0.60302,0.08204,
    ↪ 0.18691,0.5345,0.61701,0.33976,0.67302,0.296'
_aflow_Strukturbericht '$G7_{2}$'
_aflow_Pearson 'oP64'

_symmetry_space_group_name_H-M "P 21/b 21/c 21/a"
_symmetry_Int_Tables_number 61

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_cell_length_b    12.20100
_cell_length_c    4.43000
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

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_space_group_symop_operation_xyz
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2 x+1/2,-y+1/2,-z
3 -x,y+1/2,-z+1/2
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z
7 x,-y+1/2,z+1/2
8 x+1/2,y,-z+1/2

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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
B1 B 8 c 0.10617 0.60704 0.77298 1.00000
Be1 Be 8 c 0.00261 0.68871 0.26018 1.00000
Be2 Be 8 c 0.23724 0.56757 0.27717 1.00000
H1 H 8 c 0.31380 0.72280 0.45740 1.00000
O1 O 8 c 0.03760 0.68766 0.61914 1.00000
O2 O 8 c 0.10120 0.60302 0.08204 1.00000
O3 O 8 c 0.18691 0.53450 0.61701 1.00000
O4 O 8 c 0.33976 0.67302 0.29600 1.00000
```

Hambergite [$Be_2BO_3(OH)$, $G7_2$]: AB2CD4_oP64_61_c_2c_c_4c - POSCAR

```
AB2CD4_oP64_61_c_2c_c_4c & a,b/a,c/a,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,z4
    ↪ ,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8 --params=9.762,
    ↪ 1.24984634296,0.453800450727,0.10617,0.60704,0.77298,0.00261,
    ↪ 0.68871,0.26018,0.23724,0.56757,0.27717,0.3138,0.7228,0.4574,
    ↪ 0.0376,0.68766,0.61914,0.1012,0.60302,0.08204,0.18691,0.5345,
    ↪ 0.61701,0.33976,0.67302,0.296 & Pbca D_{2h}^{15} #61 (c^8) &
    ↪ oP64 & $G7_{2}$ & BBe2HO4 & Hambergite & G. D. Gatta et al.,
    ↪ Am. Mineral. 97, 1891-1897 (2012)
  1.00000000000000
  9.76200000000000   0.00000000000000   0.00000000000000
  0.00000000000000  12.20100000000000   0.00000000000000
  0.00000000000000   0.00000000000000   4.43000000000000
    B    Be    H    O
    8    16    8    32
Direct
   0.10617000000000   0.60704000000000   0.77298000000000   B    (8c)
   0.39383000000000  -0.60704000000000   1.27298000000000   B    (8c)
  -0.10617000000000   1.10704000000000  -0.27298000000000   B    (8c)
   0.60617000000000  -0.10704000000000  -0.77298000000000   B    (8c)
  -0.10617000000000  -0.60704000000000  -0.77298000000000   B    (8c)
   0.60617000000000   0.60704000000000  -0.27298000000000   B    (8c)
   0.10617000000000  -0.10704000000000   1.27298000000000   B    (8c)
   0.39383000000000   1.10704000000000   0.77298000000000   B    (8c)
   0.00261000000000   0.68871000000000   0.26018000000000   Be   (8c)
   0.49739000000000  -0.68871000000000   0.76018000000000   Be   (8c)
  -0.00261000000000   1.18871000000000   0.23982000000000   Be   (8c)
   0.50261000000000  -0.18871000000000  -0.26018000000000   Be   (8c)
  -0.00261000000000  -0.68871000000000  -0.26018000000000   Be   (8c)
   0.50261000000000   0.68871000000000   0.23982000000000   Be   (8c)
   0.00261000000000  -0.18871000000000   0.76018000000000   Be   (8c)
   0.49739000000000   1.18871000000000   0.26018000000000   Be   (8c)
   0.23724000000000   0.56757000000000   0.27717000000000   Be   (8c)
   0.26276000000000  -0.56757000000000   0.77717000000000   Be   (8c)
  -0.23724000000000   1.06757000000000   0.22283000000000   Be   (8c)
   0.73724000000000  -0.06757000000000  -0.27717000000000   Be   (8c)
  -0.23724000000000  -0.56757000000000  -0.27717000000000   Be   (8c)
   0.73724000000000   0.56757000000000   0.22283000000000   Be   (8c)
   0.23724000000000  -0.06757000000000   0.77717000000000   Be   (8c)
   0.26276000000000   1.06757000000000   0.27717000000000   Be   (8c)
   0.31380000000000   0.72280000000000   0.45740000000000   H    (8c)
   0.18620000000000  -0.72280000000000   0.95740000000000   H    (8c)
  -0.31380000000000   1.22280000000000   0.04260000000000   H    (8c)
   0.81380000000000  -0.22280000000000  -0.45740000000000   H    (8c)
  -0.31380000000000  -0.72280000000000  -0.45740000000000   H    (8c)
   0.81380000000000   0.72280000000000   0.04260000000000   H    (8c)
   0.31380000000000  -0.22280000000000   0.95740000000000   H    (8c)
   0.18620000000000   1.22280000000000   0.45740000000000   H    (8c)
   0.03760000000000   0.68766000000000   0.61914000000000   O    (8c)
   0.46240000000000  -0.68766000000000   1.11914000000000   O    (8c)
  -0.03760000000000   1.18766000000000  -0.11914000000000   O    (8c)
   0.53760000000000  -0.18766000000000  -0.61914000000000   O    (8c)
  -0.03760000000000  -0.68766000000000  -0.61914000000000   O    (8c)
   0.53760000000000   0.68766000000000  -0.11914000000000   O    (8c)
   0.03760000000000  -0.18766000000000   1.11914000000000   O    (8c)
   0.46240000000000   1.18766000000000   0.61914000000000   O    (8c)
   0.10120000000000   0.60302000000000   0.08204000000000   O    (8c)
   0.39880000000000  -0.60302000000000   0.58204000000000   O    (8c)
  -0.10120000000000   1.10302000000000   0.41796000000000   O    (8c)
   0.60120000000000  -0.10302000000000  -0.08204000000000   O    (8c)
  -0.10120000000000  -0.60302000000000  -0.08204000000000   O    (8c)
   0.60120000000000   0.60302000000000   0.41796000000000   O    (8c)
   0.10120000000000  -0.10302000000000   0.58204000000000   O    (8c)
   0.39880000000000   1.10302000000000   0.08204000000000   O    (8c)
   0.18691000000000   0.53450000000000   0.61701000000000   O    (8c)
   0.31309000000000  -0.53450000000000   1.11701000000000   O    (8c)
  -0.18691000000000   1.03450000000000  -0.11701000000000   O    (8c)
   0.68691000000000  -0.03450000000000  -0.61701000000000   O    (8c)
  -0.18691000000000  -0.53450000000000  -0.61701000000000   O    (8c)
```

```
   0.68691000000000   0.53450000000000  −0.11701000000000   O   (8c)
   0.18691000000000  −0.03450000000000   1.11701000000000   O   (8c)
   0.31309000000000   1.03450000000000   0.61701000000000   O   (8c)
   0.33976000000000   0.67302000000000   0.29600000000000   O   (8c)
   0.16024000000000  −0.67302000000000   0.79600000000000   O   (8c)
  −0.33976000000000   1.17302000000000   0.20400000000000   O   (8c)
   0.83976000000000  −0.17302000000000  −0.29600000000000   O   (8c)
  −0.33976000000000  −0.67302000000000  −0.29600000000000   O   (8c)
   0.83976000000000   0.67302000000000   0.20400000000000   O   (8c)
   0.33976000000000  −0.17302000000000   0.79600000000000   O   (8c)
   0.16024000000000   1.17302000000000   0.29600000000000   O   (8c)
```

Enstatite ($MgSiO_3$, $S4_3$): AB3C_oP80_61_2c_6c_2c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Enstatite'
_chemical_formula_sum 'Mg O3 Si'

loop_
_publ_author_name
 'B. E. Warren'
 'D. I. Modell'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 75
_journal_year 1930
_journal_page_first 1
_journal_page_last 14
_publ_Section_title
;
 The Structure of Enstatite MgSiO$_{3}$
;

_aflow_title 'Enstatite (MgSiO$_{3}$, $S4_{3}$) Structure'
_aflow_proto 'AB3C_oP80_61_2c_6c_2c'
_aflow_params 'a,b/a,c/a,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3},y_{3
      ↪ },z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{6},
      ↪ x_{7},y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{10},y_
      ↪ {10},z_{10}'
_aflow_params_values '18.2,0.486813186813,0.285714285714,0.13,0.33,0.37,
      ↪ 0.13,0.96,0.37,0.06,0.14,0.2,0.06,0.5,0.2,0.05,0.75,0.05,0.19,
      ↪ 0.35,0.06,0.19,0.01,0.05,0.2,0.75,0.3,0.03,0.65,0.29,0.22,0.85,
      ↪ 0.04'
_aflow_Strukturbericht '$S4_{3}$'
_aflow_Pearson 'oP80'

_symmetry_space_group_name_H-M "P 21/b 21/c 21/a"
_symmetry_Int_Tables_number 61

_cell_length_a     18.20000
_cell_length_b     8.86000
_cell_length_c     5.20000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z
3 -x,y+1/2,-z+1/2
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z
7 x,-y+1/2,z+1/2
8 x+1/2,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Mg1 Mg 8 c 0.13000 0.33000 0.37000 1.00000
Mg2 Mg 8 c 0.13000 0.96000 0.37000 1.00000
O1 O 8 c 0.06000 0.14000 0.20000 1.00000
O2 O 8 c 0.06000 0.50000 0.20000 1.00000
O3 O 8 c 0.05000 0.75000 0.05000 1.00000
O4 O 8 c 0.19000 0.35000 0.06000 1.00000
O5 O 8 c 0.19000 0.01000 0.05000 1.00000
O6 O 8 c 0.20000 0.75000 0.30000 1.00000
Si1 Si 8 c 0.03000 0.65000 0.29000 1.00000
Si2 Si 8 c 0.22000 0.85000 0.04000 1.00000
```

Enstatite ($MgSiO_3$, $S4_3$): AB3C_oP80_61_2c_6c_2c - POSCAR

```
AB3C_oP80_61_2c_6c_2c & a,b/a,c/a,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,z4,x5
      ↪ ,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10 --params
      ↪ =18.2,0.486813186813,0.285714285714,0.13,0.33,0.37,0.13,0.96,
      ↪ 0.37,0.06,0.14,0.2,0.06,0.5,0.2,0.05,0.75,0.05,0.19,0.35,0.06,
      ↪ 0.19,0.01,0.05,0.2,0.75,0.3,0.03,0.65,0.29,0.22,0.85,0.04 &
      ↪ Pbca D_{2h}^{15} #61 (c^10) & oP80 & $S4_{3}$ & MgO3Si &
      ↪ Enstatite & B. E. Warren and D. I. Modell, Zeitschrift f{"u}r
      ↪ Kristallographie - Crystalline Materials 75, 1-14 (1930)
   1.00000000000000
  18.20000000000000     0.00000000000000     0.00000000000000
   0.00000000000000     8.86000000000000     0.00000000000000
   0.00000000000000     0.00000000000000     5.20000000000000
     Mg    O    Si
     16   48    16
Direct
   0.13000000000000   0.33000000000000   0.37000000000000   Mg   (8c)
   0.37000000000000  -0.33000000000000   0.87000000000000   Mg   (8c)
  -0.13000000000000   0.83000000000000   0.13000000000000   Mg   (8c)
   0.63000000000000   0.17000000000000  -0.37000000000000   Mg   (8c)
  -0.13000000000000  -0.33000000000000  -0.37000000000000   Mg   (8c)
   0.63000000000000   0.33000000000000   0.13000000000000   Mg   (8c)
   0.13000000000000   0.17000000000000   0.87000000000000   Mg   (8c)
   0.37000000000000   0.83000000000000   0.37000000000000   Mg   (8c)
   0.13000000000000   0.96000000000000   0.37000000000000   Mg   (8c)
   0.37000000000000  -0.96000000000000   0.87000000000000   Mg   (8c)
  -0.13000000000000   1.46000000000000   0.13000000000000   Mg   (8c)
   0.63000000000000  -0.46000000000000  -0.37000000000000   Mg   (8c)
  -0.13000000000000  -0.96000000000000  -0.37000000000000   Mg   (8c)
   0.63000000000000   0.96000000000000   0.13000000000000   Mg   (8c)
   0.13000000000000  -0.46000000000000   0.87000000000000   Mg   (8c)
   0.37000000000000   1.46000000000000   0.37000000000000   Mg   (8c)
   0.06000000000000   0.14000000000000   0.20000000000000   O   (8c)
   0.44000000000000  -0.14000000000000   0.70000000000000   O   (8c)
  -0.06000000000000   0.64000000000000   0.30000000000000   O   (8c)
   0.56000000000000   0.36000000000000  -0.20000000000000   O   (8c)
  -0.06000000000000  -0.14000000000000  -0.20000000000000   O   (8c)
   0.56000000000000   0.14000000000000   0.30000000000000   O   (8c)
   0.06000000000000   0.36000000000000   0.70000000000000   O   (8c)
   0.44000000000000   0.64000000000000   0.20000000000000   O   (8c)
   0.06000000000000   0.50000000000000   0.20000000000000   O   (8c)
   0.44000000000000  -0.50000000000000   0.70000000000000   O   (8c)
  -0.06000000000000   1.00000000000000   0.30000000000000   O   (8c)
   0.56000000000000   0.00000000000000  -0.20000000000000   O   (8c)
  -0.06000000000000  -0.50000000000000  -0.20000000000000   O   (8c)
   0.56000000000000   0.50000000000000   0.30000000000000   O   (8c)
   0.06000000000000   0.00000000000000   0.70000000000000   O   (8c)
   0.44000000000000   1.00000000000000   0.20000000000000   O   (8c)
   0.05000000000000   0.75000000000000   0.05000000000000   O   (8c)
   0.45000000000000  -0.75000000000000   0.55000000000000   O   (8c)
  -0.05000000000000   1.25000000000000   0.45000000000000   O   (8c)
   0.55000000000000  -0.25000000000000  -0.05000000000000   O   (8c)
  -0.05000000000000  -0.75000000000000  -0.05000000000000   O   (8c)
   0.55000000000000   0.75000000000000   0.45000000000000   O   (8c)
   0.05000000000000  -0.25000000000000   0.55000000000000   O   (8c)
   0.45000000000000   1.25000000000000   0.05000000000000   O   (8c)
   0.19000000000000   0.35000000000000   0.06000000000000   O   (8c)
   0.31000000000000  -0.35000000000000   0.56000000000000   O   (8c)
  -0.19000000000000   0.85000000000000   0.44000000000000   O   (8c)
   0.69000000000000   0.15000000000000  -0.06000000000000   O   (8c)
  -0.19000000000000  -0.35000000000000  -0.06000000000000   O   (8c)
   0.69000000000000   0.35000000000000   0.44000000000000   O   (8c)
   0.19000000000000   0.15000000000000   0.56000000000000   O   (8c)
   0.31000000000000   0.85000000000000   0.06000000000000   O   (8c)
   0.19000000000000   0.01000000000000   0.05000000000000   O   (8c)
   0.31000000000000  -0.01000000000000   0.55000000000000   O   (8c)
  -0.19000000000000   0.51000000000000   0.45000000000000   O   (8c)
   0.69000000000000   0.49000000000000  -0.05000000000000   O   (8c)
  -0.19000000000000  -0.01000000000000  -0.05000000000000   O   (8c)
   0.69000000000000   0.01000000000000   0.45000000000000   O   (8c)
   0.19000000000000   0.49000000000000   0.55000000000000   O   (8c)
   0.31000000000000   0.51000000000000   0.05000000000000   O   (8c)
   0.20000000000000   0.75000000000000   0.30000000000000   O   (8c)
   0.30000000000000  -0.75000000000000   0.80000000000000   O   (8c)
  -0.20000000000000   1.25000000000000   0.20000000000000   O   (8c)
   0.70000000000000  -0.25000000000000  -0.30000000000000   O   (8c)
  -0.20000000000000  -0.75000000000000  -0.30000000000000   O   (8c)
   0.70000000000000   0.75000000000000   0.20000000000000   O   (8c)
   0.20000000000000  -0.25000000000000   0.80000000000000   O   (8c)
   0.30000000000000   1.25000000000000   0.30000000000000   O   (8c)
   0.03000000000000   0.65000000000000   0.29000000000000   Si   (8c)
   0.47000000000000  -0.65000000000000   0.79000000000000   Si   (8c)
  -0.03000000000000   1.15000000000000   0.21000000000000   Si   (8c)
   0.53000000000000  -0.15000000000000  -0.29000000000000   Si   (8c)
  -0.03000000000000  -0.65000000000000  -0.29000000000000   Si   (8c)
   0.53000000000000   0.65000000000000   0.21000000000000   Si   (8c)
   0.03000000000000  -0.15000000000000   0.79000000000000   Si   (8c)
   0.47000000000000   1.15000000000000   0.29000000000000   Si   (8c)
   0.22000000000000   0.85000000000000   0.04000000000000   Si   (8c)
   0.28000000000000  -0.85000000000000   0.54000000000000   Si   (8c)
  -0.22000000000000   1.35000000000000   0.46000000000000   Si   (8c)
   0.72000000000000  -0.35000000000000  -0.04000000000000   Si   (8c)
  -0.22000000000000  -0.85000000000000  -0.04000000000000   Si   (8c)
   0.72000000000000   0.85000000000000   0.46000000000000   Si   (8c)
   0.22000000000000  -0.35000000000000   0.54000000000000   Si   (8c)
   0.28000000000000   1.35000000000000   0.04000000000000   Si   (8c)
```

COCl: ABC_oP24_61_c_c_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'CClO'
_chemical_formula_sum 'C Cl O'

loop_
_publ_author_name
 'P. Groth'
 'O. Hassel'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 16
_journal_year 1962
```

```
_journal_page_first 2311
_journal_page_last 2317
_publ_Section_title
;
 Crystal Structures of Oxalyl Bromide and Oxalyl Chloride
;

_aflow_title 'COCl Structure'
_aflow_proto 'ABC_oP24_61_c_c_c'
_aflow_params 'a,b/a,c/a,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3},y_{3
      ↪ },z_{3}'
_aflow_params_values '6.44,0.944099378882,1.85248447205,-0.006,0.062,
      ↪ 0.053,-0.034,-0.083,0.167,0.118,0.23,0.05'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP24'

_symmetry_space_group_name_H-M "P 21/b 21/c 21/a"
_symmetry_Int_Tables_number 61

_cell_length_a     6.44000
_cell_length_b     6.08000
_cell_length_c     11.93000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z
3 -x,y+1/2,-z+1/2
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z
7 x,-y+1/2,z+1/2
8 x+1/2,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
C1 C 8 c -0.00600 0.06200 0.05300 1.00000
Cl1 Cl 8 c -0.03400 -0.08300 0.16700 1.00000
O1 O 8 c 0.11800 0.23000 0.05000 1.00000
```

COCl: ABC_oP24_61_c_c_c - POSCAR

```
ABC_oP24_61_c_c_c & a,b/a,c/a,x1,y1,z1,x2,y2,z2,x3,y3,z3 --params=6.44,
      ↪ 0.944099378882,1.85248447205,-0.006,0.062,0.053,-0.034,-0.083,
      ↪ 0.167,0.118,0.23,0.05 & Pbca D_{2h}^{15} #61 (c^3) & oP24 &
      ↪ None & ClIO & ClIO & P. Groth and O. Hassel, Acta Chem. Scand.
      ↪ 16, 2311-2317 (1962)
   1.00000000000000
   6.44000000000000    0.00000000000000    0.00000000000000
   0.00000000000000    6.08000000000000    0.00000000000000
   0.00000000000000    0.00000000000000   11.93000000000000
     C     Cl     O
     8      8      8
Direct
  -0.00600000000000    0.06200000000000    0.05300000000000    C    (8c)
   0.50600000000000   -0.06200000000000    0.55300000000000    C    (8c)
   0.00600000000000    0.56200000000000    0.44700000000000    C    (8c)
   0.49400000000000    0.43800000000000   -0.05300000000000    C    (8c)
   0.00600000000000   -0.06200000000000   -0.05300000000000    C    (8c)
   0.49400000000000    0.06200000000000    0.44700000000000    C    (8c)
  -0.00600000000000    0.43800000000000    0.55300000000000    C    (8c)
   0.50600000000000    0.56200000000000    0.05300000000000    C    (8c)
  -0.03400000000000   -0.08300000000000    0.16700000000000   Cl    (8c)
   0.53400000000000    0.08300000000000    0.66700000000000   Cl    (8c)
   0.03400000000000    0.41700000000000    0.33300000000000   Cl    (8c)
   0.46600000000000    0.58300000000000   -0.16700000000000   Cl    (8c)
   0.03400000000000    0.08300000000000   -0.16700000000000   Cl    (8c)
   0.46600000000000   -0.08300000000000    0.33300000000000   Cl    (8c)
  -0.03400000000000    0.58300000000000    0.66700000000000   Cl    (8c)
   0.53400000000000    0.41700000000000    0.16700000000000   Cl    (8c)
   0.11800000000000    0.23000000000000    0.05000000000000    O    (8c)
   0.38200000000000   -0.23000000000000    0.55000000000000    O    (8c)
  -0.11800000000000    0.73000000000000    0.45000000000000    O    (8c)
   0.61800000000000    0.27000000000000   -0.05000000000000    O    (8c)
  -0.11800000000000   -0.23000000000000   -0.05000000000000    O    (8c)
   0.61800000000000    0.23000000000000    0.45000000000000    O    (8c)
   0.11800000000000    0.27000000000000    0.55000000000000    O    (8c)
   0.38200000000000    0.73000000000000    0.05000000000000    O    (8c)
```

Topaz ($Al_2SiO_4F_2$, $S0_5$): A2B2C4D_oP36_62_d_d_2cd_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Topaz'
_chemical_formula_sum 'Al2 F2 O4 Si'

loop_
_publ_author_name
 'K. Komatsu'
 'T. Kuribayashi'
 'Y. Kudoh'
_journal_name_full_name
;
 Journal of Mineralogical and Petrological Sciences
;
_journal_volume 98
_journal_year 2003
_journal_page_first 167
_journal_page_last 180
_publ_Section_title
;
 Effect of temperature and pressure on the crystal structure of topaz,
      ↪ Al$_{2}$SiO$_{4}$(OH,F)$_{2}$
;

_aflow_title 'Topaz (Al$_{2}$SiO$_{4}$F$_{2}$, $S0_{5}$) Structure'
_aflow_proto 'A2B2C4D_oP36_62_d_d_2cd_c'
_aflow_params 'a,b/a,c/a,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3},x_{4},y_{4
      ↪ },z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{6}'
_aflow_params_values '8.7935,0.954762040143,0.528674589185,0.468,0.7957,
      ↪ 0.2436,0.4566,0.55945,0.1022,0.63102,0.5824,0.40358,0.75265,
      ↪ 0.55741,0.0981,0.4893,0.5923,0.711'
_aflow_Strukturbericht '$S0_{5}$'
_aflow_Pearson 'oP36'

_symmetry_space_group_name_H-M "P 21/n 21/m 21/a"
_symmetry_Int_Tables_number 62

_cell_length_a     8.79350
_cell_length_b     8.39570
_cell_length_c     4.64890
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x,y+1/2,-z
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x,-y+1/2,z
8 x+1/2,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 4 c 0.46800 0.25000 0.79570 1.00000
O2 O 4 c 0.24360 0.25000 0.45660 1.00000
Si1 Si 4 c 0.55945 0.25000 0.10220 1.00000
Al1 Al 8 d 0.63102 0.58240 0.40358 1.00000
F1 F 8 d 0.75265 0.55741 0.09810 1.00000
O3 O 8 d 0.48930 0.59230 0.71100 1.00000
```

Topaz ($Al_2SiO_4F_2$, $S0_5$): A2B2C4D_oP36_62_d_d_2cd_c - POSCAR

```
A2B2C4D_oP36_62_d_d_2cd_c & a,b/a,c/a,x1,z1,x2,z2,x3,z3,x4,y4,z4,x5,y5,
      ↪ z5,x6,y6,z6 --params=8.7935,0.954762040143,0.528674589185,0.468
      ↪ ,0.7957,0.2436,0.4566,0.55945,0.1022,0.63102,0.5824,0.40358,
      ↪ 0.75265,0.55741,0.0981,0.4893,0.5923,0.711 & Pnma D_{2h}^{16} #
      ↪ 62 (c^3d^3) & oP36 & $S0_{5}$ & Al2F2O4Si & Topaz & K. Komatsu
      ↪ and T. Kuribayashi and Y. Kudoh, J. Miner. Petrol. Sci. 98,
      ↪ 167-180 (2003)
   1.00000000000000
   8.79350000000000    0.00000000000000    0.00000000000000
   0.00000000000000    8.39570000000000    0.00000000000000
   0.00000000000000    0.00000000000000    4.64890000000000
     Al      F      O     Si
      8      8     16      4
Direct
   0.63102000000000    0.58240000000000    0.40358000000000   Al    (8d)
  -0.13102000000000   -0.58240000000000    0.90358000000000   Al    (8d)
  -0.63102000000000    1.08240000000000   -0.40358000000000   Al    (8d)
   1.13102000000000   -0.08240000000000    0.09642000000000   Al    (8d)
  -0.63102000000000   -0.58240000000000   -0.40358000000000   Al    (8d)
   1.13102000000000    0.58240000000000    0.09642000000000   Al    (8d)
   0.63102000000000   -0.08240000000000    0.40358000000000   Al    (8d)
  -0.13102000000000    1.08240000000000    0.90358000000000   Al    (8d)
   0.75265000000000    0.55741000000000    0.09810000000000    F    (8d)
  -0.25265000000000   -0.55741000000000    0.59810000000000    F    (8d)
  -0.75265000000000    1.05741000000000   -0.09810000000000    F    (8d)
   1.25265000000000   -0.05741000000000    0.40190000000000    F    (8d)
  -0.75265000000000   -0.55741000000000   -0.09810000000000    F    (8d)
   1.25265000000000    0.55741000000000    0.40190000000000    F    (8d)
   0.75265000000000   -0.05741000000000    0.09810000000000    F    (8d)
  -0.25265000000000    1.05741000000000    0.59810000000000    F    (8d)
   0.46800000000000    0.25000000000000    0.79570000000000    O    (4c)
   0.03200000000000    0.75000000000000    1.29570000000000    O    (4c)
  -0.46800000000000    0.75000000000000   -0.79570000000000    O    (4c)
   0.96800000000000    0.25000000000000   -0.29570000000000    O    (4c)
   0.24360000000000    0.25000000000000    0.45660000000000    O    (4c)
   0.25640000000000    0.75000000000000    0.95660000000000    O    (4c)
  -0.24360000000000    0.75000000000000   -0.45660000000000    O    (4c)
   0.74360000000000    0.25000000000000    0.04340000000000    O    (4c)
   0.48930000000000    0.59230000000000    0.71100000000000    O    (8d)
   0.01070000000000   -0.59230000000000    1.21100000000000    O    (8d)
  -0.48930000000000    1.09230000000000   -0.71100000000000    O    (8d)
   0.98930000000000   -0.09230000000000   -0.21100000000000    O    (8d)
```

```
 −0.48930000000000  −0.59230000000000  −0.71100000000000   O   (8d)
  0.98930000000000   0.59230000000000  −0.21100000000000   O   (8d)
  0.48930000000000  −0.09230000000000   0.71100000000000   O   (8d)
  0.01070000000000   1.09230000000000   1.21100000000000   O   (8d)
  0.55945000000000   0.25000000000000   0.10220000000000   Si  (4c)
 −0.05945000000000   0.75000000000000   0.60220000000000   Si  (4c)
 −0.55945000000000   0.75000000000000  −0.10220000000000   Si  (4c)
  1.05945000000000   0.25000000000000   0.39780000000000   Si  (4c)
```

Norbergite [$Mg(F,OH)_2 \cdot Mg_2SiO_4$, $S0_7$]: A2B3C4D_oP40_62_d_cd_2cd_c - CIF

```
# CIF file
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_audit_creation_method FINDSYM

_chemical_name_mineral 'Norbergite'
_chemical_formula_sum 'F2 Mg3 O4 Si'

loop_
_publ_author_name
 'G. V. Gibbs'
 'P. H. Ribbe'
_journal_name_full_name
;
 American Mineralogist
;
_journal_volume 54
_journal_year 1969
_journal_page_first 376
_journal_page_last 390
_publ_Section_title
;
 The crystal structures of the humite minerals: I. Norbergite
;

# Found in The RRUFF$^\mathroman{TM}$ Project, {Norbergite, RRUFF ID:
      ↪ R050280},

_aflow_title 'Norbergite [Mg(F,OH)$_{2}$ $\cdot$ Mg$_{2}$SiO$_{4}$, $S0_
      ↪ {7}$] Structure'
_aflow_proto 'A2B3C4D_oP40_62_d_cd_2cd_c'
_aflow_params 'a,b/a,c/a,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3},x_{4},z_{4
      ↪ },x_{5},y_{5},z_{5},x_{6},y_{6},z_{6},x_{7},y_{7},z_{7}'
_aflow_params_values '10.2718,0.851613154462,0.458575906852,-0.0923,-
      ↪ 0.0076,0.7204,0.7617,0.574,0.2793,0.7196,0.4195,-0.0318,0.0834,
      ↪ 0.7295,0.633,0.4305,-0.011,0.7907,0.1034,0.269'
_aflow_Strukturbericht '$S0_{7}$'
_aflow_Pearson 'oP40'

_symmetry_space_group_name_H-M "P 21/n 21/m 21/a"
_symmetry_Int_Tables_number 62

_cell_length_a     10.27180
_cell_length_b     8.74760
_cell_length_c     4.71040
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x,y+1/2,-z
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x,-y+1/2,z
8 x+1/2,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Mg1 Mg 4 c -0.09230 0.25000 -0.00760 1.00000
O1 O 4 c 0.72040 0.25000 0.76170 1.00000
O2 O 4 c 0.57400 0.25000 0.27930 1.00000
Si1 Si 4 c 0.71960 0.25000 0.41950 1.00000
F1 F 8 d -0.03180 0.08340 0.72950 1.00000
Mg2 Mg 8 d 0.63300 0.43050 -0.01100 1.00000
O3 O 8 d 0.79070 0.10340 0.26900 1.00000
```

Norbergite [$Mg(F,OH)_2 \cdot Mg_2SiO_4$, $S0_7$]: A2B3C4D_oP40_62_d_cd_2cd_c - POSCAR

```
A2B3C4D_oP40_62_d_cd_2cd_c & a,b/a,c/a,x1,z1,x2,z2,x3,z3,x4,z4,x5,y5,z5,
      ↪ x6,y6,z6,x7,y7,z7 --params=10.2718,0.851613154462,
      ↪ 0.458575906852,-0.0923,-0.0076,0.7204,0.7617,0.574,0.2793,
      ↪ 0.7196,0.4195,-0.0318,0.0834,0.7295,0.633,0.4305,-0.011,0.7907,
      ↪ 0.1034,0.269 & Pnma D_{2h}^{16} #62 (c^4d^3) & oP40 & $S0_{7}$
      ↪ & F2Mg3O4Si & Norbergite & G. V. Gibbs and P. H. Ribbe, Am.
      ↪ Mineral. 54, 376-390 (1969)
   1.00000000000000
  10.27180000000000     0.00000000000000     0.00000000000000
   0.00000000000000     8.74760000000000     0.00000000000000
   0.00000000000000     0.00000000000000     4.71040000000000
     F    Mg    O    Si
     8    12    16    4
Direct
 -0.03180000000000   0.08340000000000   0.72950000000000   F   (8d)
  0.53180000000000  -0.08340000000000   1.22950000000000   F   (8d)
  0.03180000000000   0.58340000000000  -0.72950000000000   F   (8d)
  0.46820000000000   0.41660000000000  -0.22950000000000   F   (8d)
  0.03180000000000  -0.08340000000000  -0.72950000000000   F   (8d)
  0.46820000000000   0.08340000000000  -0.22950000000000   F   (8d)
 -0.03180000000000   0.41660000000000   0.72950000000000   F   (8d)
  0.53180000000000   0.58340000000000   1.22950000000000   F   (8d)
 -0.09230000000000   0.25000000000000  -0.00760000000000   Mg  (4c)
  0.59230000000000   0.75000000000000   0.49240000000000   Mg  (4c)
  0.09230000000000   0.75000000000000   0.00760000000000   Mg  (4c)
  0.40770000000000   0.25000000000000   0.50760000000000   Mg  (4c)
  0.63300000000000   0.43050000000000  -0.01100000000000   Mg  (8d)
 -0.13300000000000  -0.43050000000000   0.48900000000000   Mg  (8d)
 -0.63300000000000   0.93050000000000   0.01100000000000   Mg  (8d)
  1.13300000000000   0.06950000000000   0.51100000000000   Mg  (8d)
 -0.63300000000000  -0.43050000000000   0.01100000000000   Mg  (8d)
  1.13300000000000   0.43050000000000   0.51100000000000   Mg  (8d)
  0.63300000000000   0.06950000000000  -0.01100000000000   Mg  (8d)
 -0.13300000000000   0.93050000000000   0.48900000000000   Mg  (8d)
  0.72040000000000   0.25000000000000   0.76170000000000   O   (4c)
 -0.22040000000000   0.75000000000000   1.26170000000000   O   (4c)
 -0.72040000000000   0.75000000000000  -0.76170000000000   O   (4c)
  1.22040000000000   0.25000000000000  -0.26170000000000   O   (4c)
  0.57400000000000   0.25000000000000   0.27930000000000   O   (4c)
 -0.07400000000000   0.75000000000000   0.77930000000000   O   (4c)
 -0.57400000000000   0.75000000000000  -0.27930000000000   O   (4c)
  1.07400000000000   0.25000000000000   0.22070000000000   O   (4c)
  0.79070000000000   0.10340000000000   0.26900000000000   O   (8d)
 -0.29070000000000  -0.10340000000000   0.76900000000000   O   (8d)
 -0.79070000000000   0.60340000000000  -0.26900000000000   O   (8d)
  1.29070000000000   0.39660000000000   0.23100000000000   O   (8d)
 -0.79070000000000  -0.10340000000000  -0.26900000000000   O   (8d)
  1.29070000000000   0.10340000000000   0.23100000000000   O   (8d)
  0.79070000000000   0.39660000000000   0.26900000000000   O   (8d)
 -0.29070000000000   0.60340000000000   0.76900000000000   O   (8d)
  0.71960000000000   0.25000000000000   0.41950000000000   Si  (4c)
 -0.21960000000000   0.75000000000000   0.91950000000000   Si  (4c)
 -0.71960000000000   0.75000000000000  -0.41950000000000   Si  (4c)
  1.21960000000000   0.25000000000000   0.08050000000000   Si  (4c)
```

Arcanite (K_2SO_4, $H1_6$): A2B4C_oP28_62_2c_2cd_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Arcanite'
_chemical_formula_sum 'K2 O4 S'

loop_
_publ_author_name
 'J. A. {McGinnety}'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 28
_journal_year 1972
_journal_page_first 2845
_journal_page_last 2852
_publ_Section_title
;
 Redetermination of the structures of potassium sulphate and potassium
      ↪ chromate: the effect of electrostatic crystal forces upon
      ↪ observed bond lengths
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'Arcanite (K$_{2}$SO$_{4}$, $H1_{6}$) Structure'
_aflow_proto 'A2B4C_oP28_62_2c_2cd_c'
_aflow_params 'a,b/a,c/a,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3},x_{4},z_{4
      ↪ },x_{5},z_{5},x_{6},y_{6},z_{6}'
_aflow_params_values '7.476,0.770866773676,1.34711075441,0.82623,0.21062
      ↪ ,0.51104,0.20406,0.4621,-0.0834,0.2037,0.0582,0.26702,-0.08029,
      ↪ 0.1991,0.0412,0.8522'
_aflow_Strukturbericht '$H1_{6}$'
_aflow_Pearson 'oP28'

_symmetry_space_group_name_H-M "P 21/n 21/m 21/a"
_symmetry_Int_Tables_number 62

_cell_length_a     7.47600
_cell_length_b     5.76300
_cell_length_c     10.07100
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x,y+1/2,-z
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x,-y+1/2,z
8 x+1/2,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
```

```
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
K1 K 4 c 0.82623 0.25000 0.21062 1.00000
K2 K 4 c 0.51104 0.25000 0.20406 1.00000
O1 O 4 c 0.46210 0.25000 −0.08340 1.00000
O2 O 4 c 0.20370 0.25000 0.05820 1.00000
S1 S 4 c 0.26702 0.25000 −0.08029 1.00000
O3 O 8 d 0.19910 0.04120 0.85220 1.00000
```

Arcanite (K_2SO_4, $H1_6$): A2B4C_oP28_62_2c_2cd_c - POSCAR

```
A2B4C_oP28_62_2c_2cd_c & a,b/a,c/a,x1,z1,x2,z2,x3,z3,x4,z4,x5,z5,x6,y6,
    ↪ z6 --params=7.476,0.770866773676,1.34711075441,0.82623,0.21062,
    ↪ 0.51104,0.20406,0.4621,−0.0834,0.2037,0.0582,0.26702,−0.08029,
    ↪ 0.1991,0.0412,0.8522 & Pnma D_{2h}^{16} #62 (c^5d) & oP28 &
    ↪ $H1_{6}$ & K2O4S & Arcanite & J. A. {McGinnety}, Acta
    ↪ Crystallogr. Sect. B Struct. Sci. 28, 2845−2852 (1972)
  1.00000000000000
  7.47600000000000   0.00000000000000   0.00000000000000
  0.00000000000000   5.76300000000000   0.00000000000000
  0.00000000000000   0.00000000000000  10.07100000000000
    K    O    S
    8   16    4
Direct
   0.82623000000000   0.25000000000000   0.21062000000000  K   (4c)
  −0.32623000000000   0.75000000000000   0.71062000000000  K   (4c)
  −0.82623000000000   0.75000000000000  −0.21062000000000  K   (4c)
   1.32623000000000   0.25000000000000   0.28938000000000  K   (4c)
   0.51104000000000   0.25000000000000   0.20406000000000  K   (4c)
  −0.01104000000000   0.75000000000000   0.70406000000000  K   (4c)
  −0.51104000000000   0.75000000000000  −0.20406000000000  K   (4c)
   1.01104000000000   0.25000000000000   0.29594000000000  K   (4c)
   0.46210000000000   0.25000000000000  −0.08340000000000  O   (4c)
   0.03790000000000   0.75000000000000   0.41660000000000  O   (4c)
  −0.46210000000000   0.75000000000000   0.08340000000000  O   (4c)
   0.96210000000000   0.25000000000000   0.58340000000000  O   (4c)
   0.20370000000000   0.25000000000000   0.05820000000000  O   (4c)
   0.29630000000000   0.75000000000000   0.55820000000000  O   (4c)
  −0.20370000000000   0.75000000000000  −0.05820000000000  O   (4c)
   0.70370000000000   0.25000000000000   0.44180000000000  O   (4c)
   0.19910000000000   0.04120000000000   0.85220000000000  O   (8d)
   0.30090000000000  −0.04120000000000   1.35220000000000  O   (8d)
  −0.19910000000000   0.54120000000000  −0.85220000000000  O   (8d)
   0.69910000000000   0.45880000000000  −0.35220000000000  O   (8d)
  −0.19910000000000  −0.04120000000000  −0.85220000000000  O   (8d)
   0.69910000000000   0.04120000000000  −0.35220000000000  O   (8d)
   0.19910000000000   0.45880000000000   0.85220000000000  O   (8d)
   0.30090000000000   0.54120000000000   1.35220000000000  O   (8d)
   0.26702000000000   0.25000000000000  −0.08029000000000  S   (4c)
   0.23298000000000   0.75000000000000   0.41971000000000  S   (4c)
  −0.26702000000000   0.75000000000000   0.08029000000000  S   (4c)
   0.76702000000000   0.25000000000000   0.58029000000000  S   (4c)
```

Anthophyllite ($Mg_5Fe_2Si_8O_{22}(OH)_2$, $S4_2$): A2B5C22D2E8_oP156_62_d_c2d_2c10d_2c_4d - CIF

```
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_audit_creation_method FINDSYM

_chemical_name_mineral 'Anthophyllite'
_chemical_formula_sum 'Fe2 Mg5 O22 (OH)2 Si8'

loop_
_publ_author_name
 'E. M. Walitzi'
 'F. Walter'
 'K. Ettinger'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 188
_journal_year 1989
_journal_page_first 237
_journal_page_last 244
_publ_Section_title
;
 Verfeinerung der Kristallstruktur von Anthophyllit vom Ochsenkogel/
    ↪ Gleinalpe, \"{O}sterreich
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'Anthophyllite (Mg$_{5}$Fe$_{2}$Si$_{8}$O$_{22}$(OH)$_{2}$,
    ↪ $S4_{4}$) Structure'
_aflow_proto 'A2B5C22D2E8_oP156_62_d_c2d_2c10d_2c_4d'
_aflow_params 'a,b/a,c/a,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3},x_{4},z_{4
    ↪ },x_{5},z_{5},x_{6},y_{6},z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},
    ↪ z_{8},x_{9},y_{9},z_{9},x_{10},y_{10},z_{10},x_{11},y_{11},z_{
    ↪ 11},x_{12},y_{12},z_{12},x_{13},y_{13},z_{13},x_{14},y_{14},z_{
    ↪ 14},x_{15},y_{15},z_{15},x_{16},y_{16},z_{16},x_{17},y_{17},z_{
    ↪ 17},x_{18},y_{18},z_{18},x_{19},y_{19},z_{19},x_{20},y_{20},z_{
    ↪ 20},x_{21},y_{21},z_{21},x_{22},y_{22},z_{22}'
_aflow_params_values '18.544,0.972066436583,0.284836065574,0.1255,0.893,
    ↪ 0.7975,0.4595,0.9536,0.778,0.1831,0.5573,0.0697,0.227,0.1239,−
    ↪ 0.0094,0.3897,0.1251,0.1634,0.393,0.125,0.0731,0.8196,0.1825,
    ↪ 0.1635,0.0599,0.0682,0.1636,0.7275,0.1857,0.0773,0.563,0.0631,
    ↪ 0.0772,0.2209,0.1868,−0.0012,0.0743,0.0668,−0.0068,0.7092,
    ↪ 0.1973,0.8822,0.3321,0.0504,0.8882,0.0565,0.2003,0.869,0.8288,
    ↪ 0.0486,0.8596,0.5504,0.2305,0.8345,0.567,0.0186,0.8336,0.2743,
    ↪ 0.2271,−0.0798,0.0637,0.0245,−0.0817,0.7754'
_aflow_Strukturbericht '$S4_{4}$'
_aflow_Pearson 'oP156'

_symmetry_space_group_name_H-M "P 21/n 21/m 21/a"
_symmetry_Int_Tables_number 62

_cell_length_a     18.54400
_cell_length_b     18.02600
_cell_length_c     5.28200
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
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2 x+1/2,−y+1/2,−z+1/2
3 −x,y+1/2,−z
4 −x+1/2,−y,z+1/2
5 −x,−y,−z
6 −x+1/2,y+1/2,z+1/2
7 x,−y+1/2,z
8 x+1/2,y,−z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Mg1 Mg 4 c 0.12550 0.25000 0.89300 1.00000
O1 O 4 c 0.79750 0.25000 0.45950 1.00000
O2 O 4 c 0.95360 0.25000 0.77800 1.00000
OH1 OH 4 c 0.18310 0.25000 0.55730 1.00000
OH2 OH 4 c 0.06970 0.25000 0.22700 1.00000
Fe1 Fe 8 d 0.12390 −0.00940 0.38970 1.00000
Mg2 Mg 8 d 0.12510 0.16340 0.39300 1.00000
Mg3 Mg 8 d 0.12500 0.07310 0.81960 1.00000
O3 O 8 d 0.18250 0.16350 0.05990 1.00000
O4 O 8 d 0.06820 0.16360 0.72750 1.00000
O5 O 8 d 0.18570 0.07730 0.56300 1.00000
O6 O 8 d 0.06310 0.07720 0.22090 1.00000
O7 O 8 d 0.18680 −0.00120 0.07430 1.00000
O8 O 8 d 0.06680 −0.00680 0.70920 1.00000
O9 O 8 d 0.19730 0.88220 0.33210 1.00000
O10 O 8 d 0.05040 0.88820 0.05650 1.00000
O11 O 8 d 0.20030 0.86900 0.82880 1.00000
O12 O 8 d 0.04860 0.85960 0.55040 1.00000
Si1 Si 8 d 0.23050 0.83450 0.56700 1.00000
Si2 Si 8 d 0.01860 0.83360 0.27430 1.00000
Si3 Si 8 d 0.22710 −0.07980 0.06370 1.00000
Si4 Si 8 d 0.02450 −0.08170 0.77540 1.00000
```

Anthophyllite ($Mg_5Fe_2Si_8O_{22}(OH)_2$, $S4_2$): A2B5C22D2E8_oP156_62_d_c2d_2c10d_2c_4d - POSCAR

```
A2B5C22D2E8_oP156_62_d_c2d_2c10d_2c_4d & a,b/a,c/a,x1,z1,x2,z2,x3,z3,x4,
    ↪ z4,x5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10,x11,
    ↪ y11,z11,x12,y12,z12,x13,y13,z13,x14,y14,z14,x15,y15,z15,x16,y16
    ↪ ,z16,x17,y17,z17,x18,y18,z18,x19,y19,z19,x20,y20,z20,x21,y21,
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    ↪ 0.1255,0.893,0.7975,0.4595,0.9536,0.778,0.1831,0.5573,0.0697,
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    ↪ 0.8196,0.1825,0.1635,0.0599,0.0682,0.1636,0.7275,0.1857,0.0773,
    ↪ 0.563,0.0631,0.0772,0.2209,0.1868,−0.0012,0.0743,0.0668,−0.0068
    ↪ ,0.7092,0.1973,0.8822,0.3321,0.0504,0.8882,0.0565,0.2003,0.869,
    ↪ 0.8288,0.0486,0.8596,0.5504,0.2305,0.8345,0.567,0.0186,0.8336,
    ↪ 0.2743,0.2271,−0.0798,0.0637,0.0245,−0.0817,0.7754 & Pnma D_{2h
    ↪ }^{16} #62 (c^5d^17) & oP156 & $S4_{4}$ & Fe2Mg5O22(OH)2Si8 &
    ↪ Anthophyllite & E. M. Walitzi and F. Walter and K. Ettinger,
    ↪ Zeitschrift f{"u}r Kristallographie - Crystalline Materials 188
    ↪ , 237−244 (1989)
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  0.00000000000000   0.00000000000000   5.28200000000000
    Fe    Mg     O    OH    Si
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Direct
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   0.37610000000000   0.00940000000000   0.88970000000000  Fe   (8d)
  −0.12390000000000   0.49060000000000  −0.38970000000000  Fe   (8d)
   0.62390000000000   0.50940000000000   0.11030000000000  Fe   (8d)
  −0.12390000000000   0.00940000000000  −0.38970000000000  Fe   (8d)
   0.62390000000000  −0.00940000000000   0.11030000000000  Fe   (8d)
   0.12390000000000   0.50940000000000   0.38970000000000  Fe   (8d)
   0.37610000000000   0.49060000000000   0.88970000000000  Fe   (8d)
   0.12550000000000   0.25000000000000   0.89300000000000  Mg   (4c)
   0.37450000000000   0.75000000000000   1.39300000000000  Mg   (4c)
  −0.12550000000000   0.75000000000000  −0.89300000000000  Mg   (4c)
   0.62550000000000   0.25000000000000  −0.39300000000000  Mg   (4c)
   0.12510000000000   0.16340000000000   0.39300000000000  Mg   (8d)
   0.37490000000000  −0.16340000000000   0.89300000000000  Mg   (8d)
  −0.12510000000000   0.66340000000000  −0.39300000000000  Mg   (8d)
   0.62510000000000   0.33660000000000   0.10700000000000  Mg   (8d)
  −0.12510000000000  −0.16340000000000  −0.39300000000000  Mg   (8d)
   0.62510000000000   0.16340000000000   0.10700000000000  Mg   (8d)
   0.12510000000000   0.33660000000000   0.39300000000000  Mg   (8d)
   0.37490000000000   0.66340000000000   0.89300000000000  Mg   (8d)
   0.12500000000000   0.07310000000000   0.81960000000000  Mg   (8d)
   0.37500000000000  −0.07310000000000   1.31960000000000  Mg   (8d)
  −0.12500000000000   0.57310000000000  −0.81960000000000  Mg   (8d)
   0.62500000000000   0.42690000000000  −0.31960000000000  Mg   (8d)
```

−0.12500000000000	−0.07310000000000	−0.81960000000000	Mg	(8d)
0.62500000000000	0.07310000000000	−0.31960000000000	Mg	(8d)
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0.37500000000000	0.57310000000000	1.31960000000000	Mg	(8d)
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−0.29750000000000	0.75000000000000	0.95950000000000	O	(4c)
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−0.45360000000000	0.75000000000000	1.27800000000000	O	(4c)
−0.95360000000000	0.75000000000000	−0.77800000000000	O	(4c)
1.45360000000000	0.25000000000000	−0.27800000000000	O	(4c)
0.18250000000000	0.16350000000000	0.05990000000000	O	(8d)
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0.68250000000000	0.33650000000000	0.44010000000000	O	(8d)
−0.18250000000000	−0.16350000000000	−0.05990000000000	O	(8d)
0.68250000000000	0.16350000000000	0.44010000000000	O	(8d)
0.18250000000000	0.33650000000000	0.05990000000000	O	(8d)
0.31750000000000	0.66350000000000	0.55990000000000	O	(8d)
0.06820000000000	0.16360000000000	0.72750000000000	O	(8d)
0.43180000000000	−0.16360000000000	1.22750000000000	O	(8d)
−0.06820000000000	0.66360000000000	−0.72750000000000	O	(8d)
0.56820000000000	0.33640000000000	−0.22750000000000	O	(8d)
−0.06820000000000	−0.16360000000000	−0.72750000000000	O	(8d)
0.56820000000000	0.16360000000000	−0.22750000000000	O	(8d)
0.06820000000000	0.33640000000000	0.72750000000000	O	(8d)
0.43180000000000	0.66360000000000	1.22750000000000	O	(8d)
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0.68570000000000	0.42270000000000	−0.06300000000000	O	(8d)
−0.18570000000000	−0.07730000000000	−0.56300000000000	O	(8d)
0.68570000000000	0.07730000000000	−0.06300000000000	O	(8d)
0.18570000000000	0.42270000000000	0.56300000000000	O	(8d)
0.31430000000000	0.57730000000000	1.06300000000000	O	(8d)
0.06310000000000	0.07720000000000	0.22090000000000	O	(8d)
0.43690000000000	−0.07720000000000	0.72090000000000	O	(8d)
−0.06310000000000	0.57720000000000	−0.22090000000000	O	(8d)
0.56310000000000	0.42280000000000	0.27910000000000	O	(8d)
−0.06310000000000	−0.07720000000000	−0.22090000000000	O	(8d)
0.56310000000000	0.07720000000000	0.27910000000000	O	(8d)
0.06310000000000	0.42280000000000	0.22090000000000	O	(8d)
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0.18680000000000	−0.00120000000000	0.07430000000000	O	(8d)
0.31320000000000	0.00120000000000	0.57430000000000	O	(8d)
−0.18680000000000	0.49880000000000	−0.07430000000000	O	(8d)
0.68680000000000	0.50120000000000	0.42570000000000	O	(8d)
−0.18680000000000	0.00120000000000	−0.07430000000000	O	(8d)
0.68680000000000	−0.00120000000000	0.42570000000000	O	(8d)
0.18680000000000	0.50120000000000	0.07430000000000	O	(8d)
0.31320000000000	0.49880000000000	0.57430000000000	O	(8d)
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0.43320000000000	0.00680000000000	1.20920000000000	O	(8d)
−0.06680000000000	0.49320000000000	−0.70920000000000	O	(8d)
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−0.06680000000000	0.00680000000000	−0.70920000000000	O	(8d)
0.56680000000000	−0.00680000000000	−0.20920000000000	O	(8d)
0.06680000000000	0.50680000000000	0.70920000000000	O	(8d)
0.43320000000000	0.49320000000000	1.20920000000000	O	(8d)
0.19730000000000	0.88220000000000	0.33210000000000	O	(8d)
0.30270000000000	−0.88220000000000	0.83210000000000	O	(8d)
−0.19730000000000	1.38220000000000	−0.33210000000000	O	(8d)
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−0.19730000000000	−0.88220000000000	−0.33210000000000	O	(8d)
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0.19730000000000	−0.38220000000000	0.33210000000000	O	(8d)
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0.44960000000000	1.38820000000000	0.55650000000000	O	(8d)
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0.29970000000000	−0.86900000000000	1.32880000000000	O	(8d)
−0.20030000000000	1.36900000000000	−0.82880000000000	O	(8d)
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−0.20030000000000	−0.86900000000000	−0.82880000000000	O	(8d)
0.70030000000000	0.86900000000000	−0.32880000000000	O	(8d)
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0.45140000000000	1.35960000000000	1.05040000000000	O	(8d)
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0.68310000000000	0.25000000000000	−0.05730000000000	OH	(4c)
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0.56970000000000	0.25000000000000	0.27300000000000	OH	(4c)
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0.26950000000000	−0.83450000000000	1.06700000000000	Si	(8d)
−0.23050000000000	1.33450000000000	−0.56700000000000	Si	(8d)
0.73050000000000	−0.33450000000000	−0.06700000000000	Si	(8d)
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0.01860000000000	0.83360000000000	0.27430000000000	Si	(8d)
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0.51860000000000	0.83360000000000	0.22570000000000	Si	(8d)
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0.48140000000000	1.33360000000000	0.77430000000000	Si	(8d)
0.22710000000000	−0.07980000000000	0.06370000000000	Si	(8d)
0.27290000000000	0.07980000000000	0.56370000000000	Si	(8d)
−0.22710000000000	0.42020000000000	−0.06370000000000	Si	(8d)
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−0.22710000000000	0.07980000000000	−0.06370000000000	Si	(8d)
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−0.02450000000000	0.08170000000000	−0.77540000000000	Si	(8d)
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Sillimanite (Al_2SiO_5, $S0_3$): A2B5C_oP32_62_bc_3cd_c - CIF

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 'R. M. Hazen'
 'L. W. Finger'
 'C. T. Prewitt'
 'R. T. Downs'
_journal_name_full_name
;
 Physics and Chemistry of Minerals
;
_journal_volume 25
_journal_year 1997
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 Compressibility and crystal structure of sillimanite, Al$_{2}$SiO$_{5}$
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;

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_symmetry_Int_Tables_number 62

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_space_group_symop_operation_xyz
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3 -x,y+1/2,-z
4 -x+1/2,-y,z+1/2
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8 x+1/2,y,-z+1/2

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Al2 Al 4 c 0.15480 0.25000 0.85830 1.00000
O1 O 4 c -0.09110 0.25000 0.36020 1.00000
O2 O 4 c 0.06610 0.25000 0.64340 1.00000
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Si1 Si 4 c 0.84040 0.25000 0.15320 1.00000
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Sillimanite (Al_2SiO_5, $S0_3$): A2B5C_oP32_62_bc_3cd_c - POSCAR

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    ↪ & Al2O5Si & Sillimanite & H. Yang et al., Phys. Chem. Miner. 25
    ↪ , 39-47 (1997)
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   0.00000000000000   5.77510000000000   0.00000000000000
   0.00000000000000   0.00000000000000   7.48570000000000
     Al    O    Si
      8   20     4
Direct
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   0.50000000000000   0.00000000000000   0.00000000000000   Al   (4b)
   0.00000000000000   0.50000000000000   0.50000000000000   Al   (4b)
   0.50000000000000   0.50000000000000   0.00000000000000   Al   (4b)
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   0.34520000000000   0.75000000000000   1.35830000000000   Al   (4c)
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$K_2S_3O_6$ ($K5_1$): A2B6C3_oP44_62_2c_2c2d_3c - CIF

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 'J. T. Szyma\\'{n}ski'
_journal_name_full_name
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 Acta Crystallographica Section B: Structural Science
;
_journal_volume 35
_journal_year 1979
_journal_page_first 1967
_journal_page_last 1970
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A redetermination of the crystal structure of potassium trithionate,
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_aflow_params 'a,b/a,c/a,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3},x_{4},z_{4
    ↪ },x_{5},z_{5},x_{6},z_{6},x_{7},z_{7},x_{8},y_{8},z_{8},x_{9},
    ↪ y_{9},z_{9}'
_aflow_params_values '9.794,0.58627731264,1.3977945681,0.1324,0.4084,
    ↪ 0.1792,0.736,0.8837,0.1626,0.3229,-0.0759,0.0298,0.1757,0.0907,
    ↪ 0.0289,0.3015,0.0293,-0.0807,0.5411,0.779,0.6492,0.541,-0.076'
_aflow_Strukturbericht '$K5_{1}$'
_aflow_Pearson 'oP44'

_symmetry_space_group_name_H-M "P 21/n 21/m 21/a"
_symmetry_Int_Tables_number 62

_cell_length_a     9.79400
_cell_length_b     5.74200
_cell_length_c     13.69000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x,y+1/2,-z
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x,-y+1/2,z
8 x+1/2,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
K1 K 4 c 0.13240 0.25000 0.40840 1.00000
K2 K 4 c 0.17920 0.25000 0.73600 1.00000
O1 O 4 c 0.88370 0.25000 0.16260 1.00000
O2 O 4 c 0.32290 0.25000 -0.07590 1.00000
S1 S 4 c 0.02980 0.25000 0.17570 1.00000
S2 S 4 c 0.09070 0.25000 0.02890 1.00000
S3 S 4 c 0.30150 0.25000 0.02930 1.00000
O3 O 8 d -0.08070 0.54110 0.77900 1.00000
O4 O 8 d 0.64920 0.54100 -0.07600 1.00000
```

$K_2S_3O_6$ ($K5_1$): A2B6C3_oP44_62_2c_2c2d_3c - POSCAR

```
A2B6C3_oP44_62_2c_2c2d_3c & a,b/a,c/a,x1,z1,x2,z2,x3,z3,x4,z4,x5,z5,x6,
    ↪ z6,x7,z7,x8,y8,z8,x9,y9,z9 --params=9.794,0.58627731264,
    ↪ 1.3977945681,0.1324,0.4084,0.1792,0.736,0.8837,0.1626,0.3229,-
    ↪ 0.0759,0.0298,0.1757,0.0907,0.0289,0.3015,0.0293,-0.0807,0.5411
    ↪ ,0.779,0.6492,0.541,-0.076 & Pnma D_{2h}^{16} #62 (c^7d^2) &
    ↪ oP44 & $K5_{1}$ & K2O6S3 & K2O6S3 & J. M. Stewart and J. T.
    ↪ Szyma\'{n}ski, Acta Crystallogr. Sect. B Struct. Sci. 35,
    ↪ 1967-1970 (1979)
   1.00000000000000
   9.79400000000000   0.00000000000000   0.00000000000000
   0.00000000000000   5.74200000000000   0.00000000000000
   0.00000000000000   0.00000000000000  13.69000000000000
     K    O    S
     8   24   12
Direct
   0.13240000000000   0.25000000000000   0.40840000000000   K   (4c)
   0.36760000000000   0.75000000000000   0.90840000000000   K   (4c)
  -0.13240000000000   0.75000000000000  -0.40840000000000   K   (4c)
   0.63240000000000   0.25000000000000   0.09160000000000   K   (4c)
   0.17920000000000   0.25000000000000   0.73600000000000   K   (4c)
   0.32080000000000   0.75000000000000   1.23600000000000   K   (4c)
  -0.17920000000000   0.75000000000000  -0.73600000000000   K   (4c)
   0.67920000000000   0.25000000000000  -0.23600000000000   K   (4c)
   0.88370000000000   0.25000000000000   0.16260000000000   O   (4c)
  -0.38370000000000   0.75000000000000   0.66260000000000   O   (4c)
  -0.88370000000000   0.75000000000000  -0.16260000000000   O   (4c)
   1.38370000000000   0.25000000000000   0.33740000000000   O   (4c)
   0.32290000000000   0.25000000000000  -0.07590000000000   O   (4c)
   0.17710000000000   0.75000000000000   0.42410000000000   O   (4c)
  -0.32290000000000   0.75000000000000   0.07590000000000   O   (4c)
   0.82290000000000   0.25000000000000   0.57590000000000   O   (4c)
  -0.08070000000000   0.54110000000000   0.77900000000000   O   (8d)
   0.58070000000000  -0.54110000000000   1.27900000000000   O   (8d)
   0.08070000000000   1.04110000000000  -0.77900000000000   O   (8d)
   0.41930000000000  -0.04110000000000  -0.27900000000000   O   (8d)
   0.08070000000000  -0.54110000000000  -0.77900000000000   O   (8d)
   0.41930000000000   0.54110000000000  -0.27900000000000   O   (8d)
  -0.08070000000000  -0.04110000000000   0.77900000000000   O   (8d)
   0.58070000000000   1.04110000000000   1.27900000000000   O   (8d)
   0.64920000000000   0.54100000000000  -0.07600000000000   O   (8d)
  -0.14920000000000  -0.54100000000000   0.42400000000000   O   (8d)
  -0.64920000000000   1.04100000000000   0.07600000000000   O   (8d)
   1.14920000000000  -0.04100000000000   0.57600000000000   O   (8d)
  -0.64920000000000  -0.54100000000000   0.07600000000000   O   (8d)
   1.14920000000000   0.54100000000000   0.57600000000000   O   (8d)
   0.64920000000000  -0.04100000000000  -0.07600000000000   O   (8d)
  -0.14920000000000   1.04100000000000   0.42400000000000   O   (8d)
   0.02980000000000   0.25000000000000   0.17570000000000   S   (4c)
   0.47020000000000   0.75000000000000   0.67570000000000   S   (4c)
  -0.02980000000000   0.75000000000000  -0.17570000000000   S   (4c)
   0.52980000000000   0.25000000000000   0.32430000000000   S   (4c)
   0.09070000000000   0.25000000000000   0.02890000000000   S   (4c)
   0.40930000000000   0.75000000000000   0.52890000000000   S   (4c)
  -0.09070000000000   0.75000000000000  -0.02890000000000   S   (4c)
   0.59070000000000   0.25000000000000   0.47110000000000   S   (4c)
   0.30150000000000   0.25000000000000   0.02930000000000   S   (4c)
   0.19850000000000   0.75000000000000   0.52930000000000   S   (4c)
  -0.30150000000000   0.75000000000000  -0.02930000000000   S   (4c)
   0.80150000000000   0.25000000000000   0.47070000000000   S   (4c)
```

Danburite ($CaB_2Si_2O_8$, $S6_3$): A2BC8D2_oP52_62_d_c_2c3d_d - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Danburite'
_chemical_formula_sum 'B2 Ca O8 Si2'

loop_
_publ_author_name
 'K. Sugiyama'
 'Y. Tak\\'{e}uchi'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 173
_journal_year 1985
```

```
_journal_page_first 293
_journal_page_last 304
_publ_Section_title
;
 Unusual thermal expansion of a B–O bond in the structure of danburite
      ↪ CaB$_{2}$Si$_{2}$O$_{8}$
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'Danburite (CaB$_{2}$Si$_{2}$O$_{8}$, $S6_{3}$) Structure'
_aflow_proto 'A2BC8D2_oP52_62_d_c_2c3d_d'
_aflow_params 'a,b/a,c/a,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3},x_{4},y_{4
      ↪ },z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{6},x_{7},y_{7},z_{7},
      ↪ x_{8},y_{8},z_{8}'
_aflow_params_values '8.037,0.960781386089,1.09120318527,0.61445,0.38657
      ↪ ,0.4863,0.664,0.8162,0.4282,0.7411,0.4206,0.4192,0.80709,-
      ↪ 0.00324,0.06797,0.87368,-0.04233,0.36496,0.60035,0.0782,0.31351
      ↪ ,-0.05333,-0.05574,0.1925'
_aflow_Strukturbericht '$S6_{3}$'
_aflow_Pearson 'oP52'

_symmetry_space_group_name_H-M "P 21/n 21/m 21/a"
_symmetry_Int_Tables_number 62

_cell_length_a      8.03700
_cell_length_b      7.72180
_cell_length_c      8.77000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x,y+1/2,-z
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x,-y+1/2,z
8 x+1/2,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ca1 Ca 4 c 0.61445 0.25000 0.38657 1.00000
O1 O 4 c 0.48630 0.25000 0.66400 1.00000
O2 O 4 c 0.81620 0.25000 0.42820 1.00000
B1 B 8 d 0.74110 0.42060 0.41920 1.00000
O3 O 8 d 0.80709 -0.00324 0.06797 1.00000
O4 O 8 d 0.87368 -0.04233 0.36496 1.00000
O5 O 8 d 0.60035 0.07820 0.31351 1.00000
Si1 Si 8 d -0.05333 -0.05574 0.19250 1.00000
```

Danburite ($CaB_2Si_2O_8$, $S6_3$): A2BC8D2_oP52_62_d_c_2c3d_d - POSCAR

```
A2BC8D2_oP52_62_d_c_2c3d_d & a,b/a,c/a,x1,z1,x2,z2,x3,z3,x4,y4,z4,x5,y5,
      ↪ z5,x6,y6,z6,x7,y7,z7,x8,y8,z8 --params=8.037,0.960781386089,
      ↪ 1.09120318527,0.61445,0.38657,0.4863,0.664,0.8162,0.4282,0.7411
      ↪ ,0.4206,0.4192,0.80709,-0.00324,0.06797,0.87368,-0.04233,
      ↪ 0.36496,0.60035,0.0782,0.31351,-0.05333,-0.05574,0.1925 & Pnma
      ↪ D_{2h}^{16} #62 (c^3d^5) & oP52 & $S6_{3}$ & B2CaO8Si2 &
      ↪ Danburite & K. Sugiyama and Y. Tak\'{e}uchi, Zeitschrift f{"u}r
      ↪  Kristallographie - Crystalline Materials 173, 293-304 (1985)
   1.00000000000000
   8.03700000000000   0.00000000000000   0.00000000000000
   0.00000000000000   7.72180000000000   0.00000000000000
   0.00000000000000   0.00000000000000   8.77000000000000
     B    Ca    O    Si
     8     4    32    8
Direct
   0.74110000000000   0.42060000000000   0.41920000000000   B   (8d)
  -0.24110000000000  -0.42060000000000   0.91920000000000   B   (8d)
  -0.74110000000000   0.92060000000000  -0.41920000000000   B   (8d)
   1.24110000000000   0.07940000000000   0.08080000000000   B   (8d)
  -0.74110000000000  -0.42060000000000  -0.41920000000000   B   (8d)
   1.24110000000000   0.42060000000000   0.08080000000000   B   (8d)
   0.74110000000000   0.07940000000000   0.41920000000000   B   (8d)
  -0.24110000000000   0.92060000000000   0.91920000000000   B   (8d)
   0.61445000000000   0.25000000000000   0.38657000000000   Ca  (4c)
  -0.11445000000000   0.75000000000000   0.88657000000000   Ca  (4c)
  -0.61445000000000   0.75000000000000  -0.38657000000000   Ca  (4c)
   1.11445000000000   0.25000000000000   0.11343000000000   Ca  (4c)
   0.48630000000000   0.25000000000000   0.66400000000000   O   (4c)
   0.01370000000000   0.75000000000000   1.16400000000000   O   (4c)
  -0.48630000000000   0.75000000000000  -0.66400000000000   O   (4c)
   0.98630000000000   0.25000000000000  -0.16400000000000   O   (4c)
   0.81620000000000   0.25000000000000   0.42820000000000   O   (4c)
  -0.31620000000000   0.75000000000000   0.92820000000000   O   (4c)
  -0.81620000000000   0.75000000000000  -0.42820000000000   O   (4c)
   1.31620000000000   0.25000000000000   0.07180000000000   O   (4c)
   0.80709000000000  -0.00324000000000   0.06797000000000   O   (8d)
  -0.30709000000000   0.00324000000000   0.56797000000000   O   (8d)
  -0.80709000000000   0.49676000000000  -0.06797000000000   O   (8d)
   1.30709000000000   0.50324000000000   0.43203000000000   O   (8d)
  -0.80709000000000   0.00324000000000  -0.06797000000000   O   (8d)
   1.30709000000000  -0.00324000000000   0.43203000000000   O   (8d)
   0.80709000000000   0.50324000000000   0.06797000000000   O   (8d)
  -0.30709000000000   0.49676000000000   0.56797000000000   O   (8d)
   0.87368000000000  -0.04233000000000   0.36496000000000   O   (8d)
  -0.37368000000000   0.04233000000000   0.86496000000000   O   (8d)
  -0.87368000000000   0.45767000000000  -0.36496000000000   O   (8d)
   1.37368000000000   0.54233000000000   0.13504000000000   O   (8d)
  -0.87368000000000   0.04233000000000  -0.36496000000000   O   (8d)
   1.37368000000000  -0.04233000000000   0.13504000000000   O   (8d)
   0.87368000000000   0.54233000000000   0.36496000000000   O   (8d)
  -0.37368000000000   0.45767000000000   0.86496000000000   O   (8d)
   0.60035000000000   0.07820000000000   0.31351000000000   O   (8d)
  -0.10035000000000  -0.07820000000000   0.81351000000000   O   (8d)
  -0.60035000000000   0.57820000000000  -0.31351000000000   O   (8d)
   1.10035000000000   0.42180000000000   0.18649000000000   O   (8d)
  -0.60035000000000  -0.07820000000000  -0.31351000000000   O   (8d)
   1.10035000000000   0.07820000000000   0.18649000000000   O   (8d)
   0.60035000000000   0.42180000000000   0.31351000000000   O   (8d)
  -0.10035000000000   0.57820000000000   0.81351000000000   O   (8d)
  -0.05333000000000  -0.05574000000000   0.19250000000000   Si  (8d)
   0.55333000000000   0.05574000000000   0.69250000000000   Si  (8d)
   0.05333000000000   0.44426000000000  -0.19250000000000   Si  (8d)
   0.44667000000000   0.55574000000000   0.30750000000000   Si  (8d)
   0.05333000000000   0.05574000000000  -0.19250000000000   Si  (8d)
   0.44667000000000  -0.05574000000000   0.30750000000000   Si  (8d)
  -0.05333000000000   0.55574000000000   0.19250000000000   Si  (8d)
   0.55333000000000   0.44426000000000   0.69250000000000   Si  (8d)
```

$C53$ ($SrBr_2$) (*obsolete*): A2B_oP12_62_2c_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Br2Sr'
_chemical_formula_sum 'Br2 Sr'

loop_
_publ_author_name
 'M. A. Kamermans'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 101
_journal_year 1939
_journal_page_first 406
_journal_page_last 411
_publ_Section_title
;
 The Crystal Structure of SrBr$_{2}$
;

# Found in Strukturbericht Band VII 1939, 1943

_aflow_title '$C53$ (SrBr$_{2}$) ({\em{obsolete}}) Structure'
_aflow_proto 'A2B_oP12_62_2c_c'
_aflow_params 'a,b/a,c/a,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3}'
_aflow_params_values '11.42,0.376532399299,0.805604203152,0.103,0.119,
      ↪ 0.614,0.842,0.811,0.108'
_aflow_Strukturbericht '$C53$'
_aflow_Pearson 'oP12'

_symmetry_space_group_name_H-M "P 21/n 21/m 21/a"
_symmetry_Int_Tables_number 62

_cell_length_a      11.42000
_cell_length_b      4.30000
_cell_length_c      9.20000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x,y+1/2,-z
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x,-y+1/2,z
8 x+1/2,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Br1 Br 4 c 0.10300 0.25000 0.11900 1.00000
Br2 Br 4 c 0.61400 0.25000 0.84200 1.00000
Sr1 Sr 4 c 0.81100 0.25000 0.10800 1.00000
```

$C53$ ($SrBr_2$) (*obsolete*): A2B_oP12_62_2c_c - POSCAR

```
A2B_oP12_62_2c_c & a,b/a,c/a,x1,z1,x2,z2,x3,z3 --params=11.42,
      ↪ 0.376532399299,0.805604203152,0.103,0.119,0.614,0.842,0.811,
      ↪ 0.108 & Pnma D_{2h}^{16} #62 (c^3) & oP12 & $C53$ & Br2Sr &
      ↪ Br2Sr & M. A. Kamermans, Zeitschrift f{"u}r Kristallographie -
      ↪ Crystalline Materials 101, 406-411 (1939)
   1.00000000000000
```

```
   11.42000000000000     0.00000000000000     0.00000000000000
    0.00000000000000     4.30000000000000     0.00000000000000
    0.00000000000000     0.00000000000000     9.20000000000000
      Br      Sr
       8       4
Direct
    0.10300000000000     0.25000000000000     0.11900000000000    Br    (4c)
    0.39700000000000     0.75000000000000     0.61900000000000    Br    (4c)
   -0.10300000000000     0.75000000000000    -0.11900000000000    Br    (4c)
    0.60300000000000     0.25000000000000     0.38100000000000    Br    (4c)
    0.61400000000000     0.25000000000000     0.84200000000000    Br    (4c)
   -0.11400000000000     0.75000000000000     1.34200000000000    Br    (4c)
   -0.61400000000000     0.75000000000000    -0.84200000000000    Br    (4c)
    1.11400000000000     0.25000000000000    -0.34200000000000    Br    (4c)
    0.81100000000000     0.25000000000000     0.10800000000000    Sr    (4c)
   -0.31100000000000     0.75000000000000     0.60800000000000    Sr    (4c)
   -0.81100000000000     0.75000000000000    -0.10800000000000    Sr    (4c)
    1.31100000000000     0.25000000000000     0.39200000000000    Sr    (4c)
```

Cs_2Sb: A2B_oP24_62_4c_2c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Cs2Sb'
_chemical_formula_sum 'Cs2 Sb'

loop_
_publ_author_name
 'C. Hirschle'
 'C. R\"{o}hr'
_journal_name_full_name
;
 Zeitschrift fur Anorganische und Allgemeine Chemie
;
_journal_volume 626
_journal_year 2000
_journal_page_first 1992
_journal_page_last 1998
_publ_Section_title
;
 Darstellung und Kristallstruktur der bekannten Zintl-Phasen Cs$_{3}
      ↪ $Sb$_{7}$ und Cs$_{4}$Sb$_{2}$
;

# Found in PAULING FILE in: Inorganic Solid Phases (online database), {
      ↪ Cs$_{4}$Sb$_{2}$ (Cs$_{2}$Sb) Crystal Structure}, 2016

_aflow_title 'Cs$_{2}$Sb Structure'
_aflow_proto 'A2B_oP24_62_4c_2c'
_aflow_params 'a,b/a,c/a,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3},x_{4},z_{4
      ↪ },x_{5},z_{5},x_{6},z_{6}'
_aflow_params_values '15.985,0.395308101345,0.687832342821,0.00065,
      ↪ 0.6743,0.20442,0.40591,0.25609,0.73979,0.43929,0.41364,0.16182,
      ↪ 0.06288,0.34361,0.09158'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP24'

_symmetry_space_group_name_H-M "P 21/n 21/m 21/a"
_symmetry_Int_Tables_number 62

_cell_length_a     15.98500
_cell_length_b     6.31900
_cell_length_c     10.99500
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x,y+1/2,-z
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x,-y+1/2,z
8 x+1/2,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cs1 Cs 4 c 0.00065 0.25000 0.67430 1.00000
Cs2 Cs 4 c 0.20442 0.25000 0.40591 1.00000
Cs3 Cs 4 c 0.25609 0.25000 0.73979 1.00000
Cs4 Cs 4 c 0.43929 0.25000 0.41364 1.00000
Sb1 Sb 4 c 0.16182 0.25000 0.06288 1.00000
Sb2 Sb 4 c 0.34361 0.25000 0.09158 1.00000
```

Cs_2Sb: A2B_oP24_62_4c_2c - POSCAR

```
A2B_oP24_62_4c_2c & a,b/a,c/a,x1,z1,x2,z2,x3,z3,x4,z4,x5,z5,x6,z6 --
      ↪ params=15.985,0.395308101345,0.687832342821,0.00065,0.6743,
      ↪ 0.20442,0.40591,0.25609,0.73979,0.43929,0.41364,0.16182,0.06288
      ↪ ,0.34361,0.09158 & Pnma D_{2h}^{16} #62 (c^6) & oP24 & None &
      ↪ Cs2Sb & Cs2Sb & C. Hirschle and C. R\"{o}hr, Z. Anorg. Allg.
      ↪ Chem. 626, 1992-1998 (2000)
    1.00000000000000
   15.98500000000000     0.00000000000000     0.00000000000000
    0.00000000000000     6.31900000000000     0.00000000000000
    0.00000000000000     0.00000000000000    10.99500000000000
      Cs      Sb
      16       8
Direct
    0.00065000000000     0.25000000000000     0.67430000000000    Cs    (4c)
    0.49935000000000     0.75000000000000     1.17430000000000    Cs    (4c)
   -0.00065000000000     0.75000000000000    -0.67430000000000    Cs    (4c)
    0.50065000000000     0.25000000000000    -0.17430000000000    Cs    (4c)
    0.20442000000000     0.25000000000000     0.40591000000000    Cs    (4c)
    0.29558000000000     0.75000000000000     0.90591000000000    Cs    (4c)
   -0.20442000000000     0.75000000000000    -0.40591000000000    Cs    (4c)
    0.70442000000000     0.25000000000000     0.09409000000000    Cs    (4c)
    0.25609000000000     0.25000000000000     0.73979000000000    Cs    (4c)
    0.24391000000000     0.75000000000000     1.23979000000000    Cs    (4c)
   -0.25609000000000     0.75000000000000    -0.73979000000000    Cs    (4c)
    0.75609000000000     0.25000000000000    -0.23979000000000    Cs    (4c)
    0.43929000000000     0.25000000000000     0.41364000000000    Cs    (4c)
    0.06071000000000     0.75000000000000     0.91364000000000    Cs    (4c)
   -0.43929000000000     0.75000000000000    -0.41364000000000    Cs    (4c)
    0.93929000000000     0.25000000000000     0.08636000000000    Cs    (4c)
    0.16182000000000     0.25000000000000     0.06288000000000    Sb    (4c)
    0.33818000000000     0.75000000000000     0.56288000000000    Sb    (4c)
   -0.16182000000000     0.75000000000000    -0.06288000000000    Sb    (4c)
    0.66182000000000     0.25000000000000     0.43712000000000    Sb    (4c)
    0.34361000000000     0.25000000000000     0.09158000000000    Sb    (4c)
    0.15639000000000     0.75000000000000     0.59158000000000    Sb    (4c)
   -0.34361000000000     0.75000000000000    -0.09158000000000    Sb    (4c)
    0.84361000000000     0.25000000000000     0.40842000000000    Sb    (4c)
```

$RhCl_2(NH_3)_5Cl$ ($J1_8$): A3B15C5D_oP96_62_cd_3c6d_3cd_c - CIF

```
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_audit_creation_method FINDSYM

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_chemical_formula_sum 'Cl3 H15 N5 Rh'

loop_
_publ_author_name
 'R. S. Evans'
 'E. A. Hopcus'
 'J. Bordner'
 'A. F. Schreiner'
_journal_name_full_name
;
 Journal of Crystal and Molecular Structure
;
_journal_volume 3
_journal_year 1973
_journal_page_first 235
_journal_page_last 245
_publ_Section_title
;
 Molecular and crystal structures of halopentaamminerhodium-(III)
      ↪ complexes, [Rh(NH$_{3}$)$_{5}$Cl]Cl$_{2}$ and [Rh(NH$_{3}$)$_{
      ↪ 5}$Br]Br$_{2}$
;

# Found in Comparisons of $\pi$-bonding and hydrogen bonding in
      ↪ isomorphous compounds: [$M$(NH$_{3}$)$_{5}$Cl]Cl$_{2}$ ($M$=Cr,
      ↪ Co, Rh, Ir, Ru, Os), 1986

_aflow_title 'RhCl$_{2}$(NH$_{3}$)$_{5}$Cl ($J1_{8}$) Structure'
_aflow_proto 'A3B15C5D_oP96_62_cd_3c6d_3cd_c'
_aflow_params 'a,b/a,c/a,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3},x_{4},z_{4
      ↪ },x_{5},z_{5},x_{6},z_{6},x_{7},z_{7},x_{8},z_{8},x_{9},y_{9},
      ↪ z_{9},x_{10},y_{10},z_{10},x_{11},y_{11},z_{11},x_{12},y_{12},
      ↪ z_{12},x_{13},y_{13},z_{13},x_{14},y_{14},z_{14},x_{15},y_{15},
      ↪ z_{15},x_{16},y_{16},z_{16}'
_aflow_params_values '13.36,0.782934131737,0.504565868263,0.4705,-0.0482
      ↪ ,0.771,-0.003,0.527,0.527,0.778,0.337,0.7207,0.3777,0.7041,-
      ↪ 0.0519,0.5046,0.4129,0.6038,0.1805,0.1484,-0.0018,0.8394,0.396,
      ↪ 0.509,-0.06,0.357,0.52,0.739,0.452,0.528,0.798,0.297,0.696,
      ↪ 0.125,0.522,0.807,0.567,0.274,0.809,0.541,0.3975,0.5534,0.8224'
_aflow_Strukturbericht '$J1_{8}$'
_aflow_Pearson 'oP96'

_symmetry_space_group_name_H-M "P 21/n 21/m 21/a"
_symmetry_Int_Tables_number 62

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_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

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_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x,y+1/2,-z
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x,-y+1/2,z
8 x+1/2,y,-z+1/2

loop_
_atom_site_label
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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cl1  Cl  4  c  0.47050  0.25000  -0.04820  1.00000
H1  H  4  c  0.77100  0.25000  -0.00300  1.00000
H2  H  4  c  0.52700  0.25000  0.52700  1.00000
H3  H  4  c  0.77800  0.25000  0.33700  1.00000
N1  N  4  c  0.72070  0.25000  0.37770  1.00000
N2  N  4  c  0.70410  0.25000  -0.05190  1.00000
N3  N  4  c  0.50460  0.25000  0.41290  1.00000
Rh1  Rh  4  c  0.60380  0.25000  0.18050  1.00000
Cl2  Cl  8  d  0.14840  -0.00180  0.83940  1.00000
H4  H  8  d  0.39600  0.50900  -0.06000  1.00000
H5  H  8  d  0.35700  0.52000  0.73900  1.00000
H6  H  8  d  0.45200  0.52800  0.79800  1.00000
H7  H  8  d  0.29700  0.69600  0.12500  1.00000
H8  H  8  d  0.52200  0.80700  0.56700  1.00000
H9  H  8  d  0.27400  0.80900  0.54100  1.00000
N4  N  8  d  0.39750  0.55340  0.82240  1.00000
```

$RhCl_2(NH_3)_5Cl$ ($J1_8$): A3B15C5D_oP96_62_cd_3c6d_3cd_c - POSCAR

```
A3B15C5D_oP96_62_cd_3c6d_3cd_c & a,b/a,c/a,x1,z1,x2,z2,x3,z3,x4,z4,x5,z5
    ↪ ,x6,z6,x7,z7,x8,z8,x9,y9,z9,x10,y10,z10,x11,y11,z11,x12,y12,z12
    ↪ ,x13,y13,z13,x14,y14,z14,x15,y15,z15,x16,y16,z16 --params=13.36
    ↪ ,0.782934131737,0.504565868263,0.4705,-0.0482,0.771,-0.003,
    ↪ 0.527,0.527,0.778,0.337,0.7207,0.3777,0.7041,-0.0519,0.5046,
    ↪ 0.4129,0.6038,0.1805,0.1484,-0.0018,0.8394,0.396,0.509,-0.06,
    ↪ 0.357,0.52,0.739,0.452,0.528,0.798,0.297,0.696,0.125,0.522,
    ↪ 0.807,0.567,0.274,0.809,0.541,0.3975,0.5534,0.8224 & Pnma D_{2h
    ↪ }^{16} #62 (c^8d^8) & oP96 & $J1_{8}$ & Cl3H15N5Rh & Cl3H15N5Rh
    ↪ & R. S. Evans et al., J. Cryst. Mol. Struct. 3, 235-245 (1973)
   1.00000000000000
  13.36000000000000    0.00000000000000    0.00000000000000
   0.00000000000000   10.46000000000000    0.00000000000000
   0.00000000000000    0.00000000000000    6.74100000000000
     Cl     H     N    Rh
     12    60    20     4
Direct
   0.47050000000000   0.25000000000000  -0.04820000000000   Cl   (4c)
   0.02950000000000   0.75000000000000   0.45180000000000   Cl   (4c)
  -0.47050000000000   0.75000000000000   0.04820000000000   Cl   (4c)
   0.97050000000000   0.25000000000000   0.54820000000000   Cl   (4c)
   0.14840000000000  -0.00180000000000   0.83940000000000   Cl   (8d)
   0.35160000000000   0.00180000000000   1.33940000000000   Cl   (8d)
  -0.14840000000000   0.49820000000000  -0.83940000000000   Cl   (8d)
   0.64840000000000   0.50180000000000  -0.33940000000000   Cl   (8d)
  -0.14840000000000   0.00180000000000  -0.83940000000000   Cl   (8d)
   0.64840000000000  -0.00180000000000  -0.33940000000000   Cl   (8d)
   0.14840000000000   0.50180000000000   0.83940000000000   Cl   (8d)
   0.35160000000000   0.49820000000000   1.33940000000000   Cl   (8d)
   0.77100000000000   0.25000000000000  -0.00300000000000    H   (4c)
  -0.27100000000000   0.75000000000000   0.49700000000000    H   (4c)
  -0.77100000000000   0.75000000000000   0.00300000000000    H   (4c)
   1.27100000000000   0.25000000000000   0.50300000000000    H   (4c)
   0.52700000000000   0.25000000000000   0.52700000000000    H   (4c)
  -0.02700000000000   0.75000000000000   1.02700000000000    H   (4c)
  -0.52700000000000   0.75000000000000  -0.52700000000000    H   (4c)
   1.02700000000000   0.25000000000000  -0.02700000000000    H   (4c)
   0.77800000000000   0.25000000000000   0.33700000000000    H   (4c)
  -0.27800000000000   0.75000000000000   0.83700000000000    H   (4c)
  -0.77800000000000   0.75000000000000  -0.33700000000000    H   (4c)
   1.27800000000000   0.25000000000000   0.16300000000000    H   (4c)
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   0.10400000000000  -0.50900000000000   0.44000000000000    H   (8d)
  -0.39600000000000   1.00900000000000   0.06000000000000    H   (8d)
   0.89600000000000  -0.00900000000000   0.56000000000000    H   (8d)
  -0.39600000000000  -0.50900000000000   0.06000000000000    H   (8d)
   0.89600000000000   0.50900000000000   0.56000000000000    H   (8d)
   0.39600000000000  -0.00900000000000  -0.06000000000000    H   (8d)
   0.10400000000000   1.00900000000000   0.44000000000000    H   (8d)
   0.35700000000000   0.52000000000000   0.73900000000000    H   (8d)
   0.14300000000000  -0.52000000000000   1.23900000000000    H   (8d)
  -0.35700000000000   1.02000000000000  -0.73900000000000    H   (8d)
   0.85700000000000  -0.02000000000000  -0.23900000000000    H   (8d)
  -0.35700000000000  -0.52000000000000  -0.73900000000000    H   (8d)
   0.85700000000000   0.52000000000000  -0.23900000000000    H   (8d)
   0.35700000000000  -0.02000000000000   0.73900000000000    H   (8d)
   0.14300000000000   1.02000000000000   1.23900000000000    H   (8d)
   0.45200000000000   0.52800000000000   0.79800000000000    H   (8d)
   0.04800000000000  -0.52800000000000   1.29800000000000    H   (8d)
  -0.45200000000000   1.02800000000000  -0.79800000000000    H   (8d)
   0.95200000000000  -0.02800000000000  -0.29800000000000    H   (8d)
  -0.45200000000000  -0.52800000000000  -0.79800000000000    H   (8d)
   0.95200000000000   0.52800000000000  -0.29800000000000    H   (8d)
   0.45200000000000  -0.02800000000000   0.79800000000000    H   (8d)
   0.04800000000000   1.02800000000000   1.29800000000000    H   (8d)
   0.29700000000000   0.69600000000000   0.12500000000000    H   (8d)
   0.20300000000000  -0.69600000000000   0.62500000000000    H   (8d)
  -0.29700000000000   1.19600000000000  -0.12500000000000    H   (8d)
   0.79700000000000  -0.19600000000000   0.37500000000000    H   (8d)
  -0.29700000000000  -0.69600000000000  -0.12500000000000    H   (8d)
   0.79700000000000   0.69600000000000   0.37500000000000    H   (8d)
   0.29700000000000  -0.19600000000000   0.12500000000000    H   (8d)
   0.20300000000000   1.19600000000000   0.62500000000000    H   (8d)
   0.52200000000000   0.80700000000000   0.56700000000000    H   (8d)
  -0.02200000000000  -0.80700000000000   1.06700000000000    H   (8d)
  -0.52200000000000   1.30700000000000  -0.56700000000000    H   (8d)
   1.02200000000000  -0.30700000000000  -0.06700000000000    H   (8d)
  -0.52200000000000  -0.80700000000000  -0.56700000000000    H   (8d)
   1.02200000000000   0.80700000000000  -0.06700000000000    H   (8d)
   0.52200000000000  -0.30700000000000   0.56700000000000    H   (8d)
  -0.02200000000000   1.30700000000000   1.06700000000000    H   (8d)
   0.27400000000000   0.80900000000000   0.54100000000000    H   (8d)
   0.22600000000000  -0.80900000000000   1.04100000000000    H   (8d)
  -0.27400000000000   1.30900000000000  -0.54100000000000    H   (8d)
   0.77400000000000  -0.30900000000000  -0.04100000000000    H   (8d)
  -0.27400000000000  -0.80900000000000  -0.54100000000000    H   (8d)
   0.77400000000000   0.80900000000000  -0.04100000000000    H   (8d)
   0.27400000000000  -0.30900000000000   0.54100000000000    H   (8d)
   0.22600000000000   1.30900000000000   1.04100000000000    H   (8d)
   0.72070000000000   0.25000000000000   0.37770000000000    N   (4c)
  -0.22070000000000   0.75000000000000   0.87770000000000    N   (4c)
  -0.72070000000000   0.75000000000000  -0.37770000000000    N   (4c)
   1.22070000000000   0.25000000000000   0.12230000000000    N   (4c)
   0.70410000000000   0.25000000000000  -0.05190000000000    N   (4c)
  -0.20410000000000   0.75000000000000   0.44810000000000    N   (4c)
  -0.70410000000000   0.75000000000000   0.05190000000000    N   (4c)
   1.20410000000000   0.25000000000000   0.55190000000000    N   (4c)
   0.50460000000000   0.25000000000000   0.41290000000000    N   (4c)
  -0.00460000000000   0.75000000000000   0.91290000000000    N   (4c)
  -0.50460000000000   0.75000000000000  -0.41290000000000    N   (4c)
   1.00460000000000   0.25000000000000   0.08710000000000    N   (4c)
   0.39750000000000   0.55340000000000   0.82240000000000    N   (8d)
   0.10250000000000  -0.55340000000000   1.32240000000000    N   (8d)
  -0.39750000000000   1.05340000000000  -0.82240000000000    N   (8d)
   0.89750000000000  -0.05340000000000  -0.32240000000000    N   (8d)
  -0.39750000000000  -0.55340000000000  -0.82240000000000    N   (8d)
   0.89750000000000   0.55340000000000  -0.32240000000000    N   (8d)
   0.39750000000000  -0.05340000000000   0.82240000000000    N   (8d)
   0.10250000000000   1.05340000000000   1.32240000000000    N   (8d)
   0.60380000000000   0.25000000000000   0.18050000000000   Rh   (4c)
  -0.10380000000000   0.75000000000000   0.68050000000000   Rh   (4c)
  -0.60380000000000   0.75000000000000  -0.18050000000000   Rh   (4c)
   1.10380000000000   0.25000000000000   0.31950000000000   Rh   (4c)
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NH_4I_3 ($D0_{16}$): A3B_oP16_62_3c_c - CIF

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 'A. J. T. Finney'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 26
_journal_year 1970
_journal_page_first 904
_journal_page_last 906
_publ_Section_title
;
 Refinement of the structure of ammonium triiodide, NH$_{4}$I$_{3}$
;

# Found in The Structure and Stability of Simple Tri-Iodides, {Ph.D.
    ↪ Thesis, University of Tasmania}, 1973

_aflow_title 'NH$_{4}$I$_{3}$ ($D0_{16}$) Structure'
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    ↪ }'
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    ↪ 0.15287,0.6188,-0.04894,0.42157,0.76482,0.16484,0.02957'
_aflow_Strukturbericht '$D0_{16}$'
_aflow_Pearson 'oP16'

_symmetry_space_group_name_H-M "P 21/n 21/m 21/a"
_symmetry_Int_Tables_number 62

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_cell_angle_beta    90.00000
_cell_angle_gamma   90.00000

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_space_group_symop_operation_xyz
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2 x+1/2,-y+1/2,-z+1/2
3 -x,y+1/2,-z
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x,-y+1/2,z
8 x+1/2,y,-z+1/2

loop_
_atom_site_label
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_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
I1  I  4  c  0.84315  0.25000  0.15287  1.00000
```

```
I2  I  4  c  0.61880  0.25000  −0.04894  1.00000
I3  I  4  c  0.42157  0.25000  0.76482  1.00000
NH41 NH4  4  c  0.16484  0.25000  0.02957  1.00000
```

NH_4I_3 ($D0_{16}$): A3B_oP16_62_3c_c - POSCAR

```
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    ↪ 0.613735095665,0.89305850818,0.84315,0.15287,0.6188,−0.04894,
    ↪ 0.42157,0.76482,0.16484,0.02957 & Pnma D_{2h}^{16} #62 (c^4) &
    ↪ oP16 & $D0_{16}$ & I3(NH4) & I3(NH4) & G. H. Cheesman and A. J.
    ↪  T. Finney, Acta Crystallogr. Sect. B Struct. Sci. 26, 904−906
    ↪ (1970)
  1.00000000000000
 10.81900000000000   0.00000000000000   0.00000000000000
  0.00000000000000   6.64000000000000   0.00000000000000
  0.00000000000000   0.00000000000000   9.66200000000000
    I    NH4
   12      4
Direct
  0.84315000000000   0.25000000000000   0.15287000000000    I    (4c)
 −0.34315000000000   0.75000000000000   0.65287000000000    I    (4c)
 −0.84315000000000   0.75000000000000  −0.15287000000000    I    (4c)
  1.34315000000000   0.25000000000000   0.34713000000000    I    (4c)
  0.61880000000000   0.25000000000000  −0.04894000000000    I    (4c)
 −0.11880000000000   0.75000000000000   0.45106000000000    I    (4c)
 −0.61880000000000   0.75000000000000   0.04894000000000    I    (4c)
  1.11880000000000   0.25000000000000   0.54894000000000    I    (4c)
  0.42157000000000   0.25000000000000   0.76482000000000    I    (4c)
  0.07843000000000   0.75000000000000   1.26482000000000    I    (4c)
 −0.42157000000000   0.75000000000000  −0.76482000000000    I    (4c)
  0.92157000000000   0.25000000000000  −0.26482000000000    I    (4c)
  0.16484000000000   0.25000000000000   0.02957000000000  NH4    (4c)
  0.33516000000000   0.75000000000000   0.52957000000000  NH4    (4c)
 −0.16484000000000   0.75000000000000  −0.02957000000000  NH4    (4c)
  0.66484000000000   0.25000000000000   0.47043000000000  NH4    (4c)
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Original β-WO_3 (*obsolete*): A3B_oP32_62_ab4c_2c - CIF

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_audit_creation_method FINDSYM

_chemical_name_mineral 'O3W'
_chemical_formula_sum 'O3 W'

loop_
_publ_author_name
 'P. M. Woodward'
 'A. W. Sleight'
 'T. Vogt'
_journal_name_full_name
;
 Journal of Solid State Chemistry
;
_journal_volume 131
_journal_year 1997
_journal_page_first 9
_journal_page_last 17
_publ_Section_title
;
 Ferroelectric Tungsten Trioxide
;

_aflow_title 'Original $\beta$-WO$_{3}$ ({\em obsolete}) Structure'
_aflow_proto 'A3B_oP32_62_ab4c_2c'
_aflow_params 'a,b/a,c/a,x_{3},z_{3},x_{4},z_{4},x_{5},z_{5},x_{6},z_{6
    ↪ },x_{7},z_{7},x_{8},z_{8}'
_aflow_params_values '7.57,0.969749009247,1.02430647292,0.231,0.473,
    ↪ 0.222,0.029,0.496,0.238,0.485,0.724,0.471,0.469,0.47,−0.032'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP32'

_symmetry_space_group_name_H-M "P 21/n 21/m 21/a"
_symmetry_Int_Tables_number 62

_cell_length_a      7.57000
_cell_length_b      7.34100
_cell_length_c      7.75400
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,−y+1/2,−z+1/2
3 −x,y+1/2,−z
4 −x+1/2,−y,z+1/2
5 −x,−y,−z
6 −x+1/2,y+1/2,z+1/2
7 x,−y+1/2,z
8 x+1/2,y,−z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 4 a 0.00000 0.00000 0.00000 1.00000
O2 O 4 b 0.00000 0.00000 0.50000 1.00000
O3 O 4 c 0.23100 0.25000 0.47300 1.00000
O4 O 4 c 0.22200 0.25000 0.02900 1.00000
O5 O 4 c 0.49600 0.25000 0.23800 1.00000
O6 O 4 c 0.48500 0.25000 0.72400 1.00000
W1 W 4 c 0.47100 0.25000 0.46900 1.00000
W2 W 4 c 0.47000 0.25000 −0.03200 1.00000
```

Original β-WO_3 (*obsolete*): A3B_oP32_62_ab4c_2c - POSCAR

```
A3B_oP32_62_ab4c_2c & a,b/a,c/a,x3,z3,x4,z4,x5,z5,x6,z6,x7,z7,x8,z8 --
    ↪ params=7.57,0.969749009247,1.02430647292,0.231,0.473,0.222,
    ↪ 0.029,0.496,0.238,0.485,0.724,0.471,0.469,0.47,−0.032 & Pnma D_
    ↪ {2h}^{16} #62 (abc^6) & oP32 & None & O3W & O3W & P. M.
    ↪ Woodward and A. W. Sleight and T. Vogt, J. Solid State Chem.
    ↪ 131, 9−17 (1997)
  1.00000000000000
  7.57000000000000   0.00000000000000   0.00000000000000
  0.00000000000000   7.34100000000000   0.00000000000000
  0.00000000000000   0.00000000000000   7.75400000000000
    O    W
   24    8
Direct
  0.00000000000000   0.00000000000000   0.00000000000000    O    (4a)
  0.50000000000000   0.00000000000000   0.50000000000000    O    (4a)
  0.00000000000000   0.50000000000000   0.00000000000000    O    (4a)
  0.50000000000000   0.50000000000000   0.50000000000000    O    (4a)
  0.00000000000000   0.00000000000000   0.50000000000000    O    (4b)
  0.50000000000000   0.00000000000000   0.00000000000000    O    (4b)
  0.00000000000000   0.50000000000000   0.50000000000000    O    (4b)
  0.50000000000000   0.50000000000000   0.00000000000000    O    (4b)
  0.23100000000000   0.25000000000000   0.47300000000000    O    (4c)
  0.26900000000000   0.75000000000000   0.97300000000000    O    (4c)
 −0.23100000000000   0.75000000000000  −0.47300000000000    O    (4c)
  0.73100000000000   0.25000000000000   0.02700000000000    O    (4c)
  0.22200000000000   0.25000000000000   0.02900000000000    O    (4c)
  0.27800000000000   0.75000000000000   0.52900000000000    O    (4c)
 −0.22200000000000   0.75000000000000  −0.02900000000000    O    (4c)
  0.72200000000000   0.25000000000000   0.47100000000000    O    (4c)
  0.49600000000000   0.25000000000000   0.23800000000000    O    (4c)
  0.00400000000000   0.75000000000000   0.73800000000000    O    (4c)
 −0.49600000000000   0.75000000000000  −0.23800000000000    O    (4c)
  0.99600000000000   0.25000000000000   0.26200000000000    O    (4c)
  0.48500000000000   0.25000000000000   0.72400000000000    O    (4c)
  0.01500000000000   0.75000000000000   1.22400000000000    O    (4c)
 −0.48500000000000   0.75000000000000  −0.72400000000000    O    (4c)
  0.98500000000000   0.25000000000000  −0.22400000000000    O    (4c)
  0.47100000000000   0.25000000000000   0.46900000000000    W    (4c)
  0.02900000000000   0.75000000000000   0.96900000000000    W    (4c)
 −0.47100000000000   0.75000000000000  −0.46900000000000    W    (4c)
  0.97100000000000   0.25000000000000   0.03100000000000    W    (4c)
  0.47000000000000   0.25000000000000  −0.03200000000000    W    (4c)
  0.03000000000000   0.75000000000000   0.46800000000000    W    (4c)
 −0.47000000000000   0.75000000000000   0.03200000000000    W    (4c)
  0.97000000000000   0.25000000000000   0.53200000000000    W    (4c)
```

P_4Se_3: A4B3_oP112_62_8c4d_4c4d - CIF

```
# CIF file
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_audit_creation_method FINDSYM

_chemical_name_mineral 'P4Se3'
_chemical_formula_sum 'P4 Se3'

loop_
_publ_author_name
 'E. Keulen'
 'A. Vos'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 12
_journal_year 1959
_journal_page_first 323
_journal_page_last 329
_publ_Section_title
;
 The Crystal Structure of P$_{4}$Se$_{3}$
;

_aflow_title 'P$_{4}$Se$_{3}$ Structure'
_aflow_proto 'A4B3_oP112_62_8c4d_4c4d'
_aflow_params 'a,b/a,c/a,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3},x_{4},z_{4
    ↪ },x_{5},z_{5},x_{6},z_{6},x_{7},z_{7},x_{8},z_{8},x_{9},z_{9},
    ↪ x_{10},z_{10},x_{11},z_{11},x_{12},z_{12},x_{13},y_{13},z_{13},
    ↪ x_{14},y_{14},z_{14},x_{15},y_{15},z_{15},x_{16},y_{16},z_{16},
    ↪ x_{17},y_{17},z_{17},x_{18},y_{18},z_{18},x_{19},y_{19},z_{19},
    ↪ x_{20},y_{20},z_{20}'
_aflow_params_values '11.797,0.825548868356,2.22683733152,0.086,0.627,
    ↪ 0.041,0.5,−0.034,0.282,0.238,0.331,−0.048,0.825,0.239,0.807,
    ↪ 0.192,0.104,−0.032,0.015,−0.057,0.571,0.0148,0.2567,0.107,0.87,
    ↪ 0.163,0.018,0.227,0.134,0.5875,0.304,0.634,0.6443,−0.01,0.134,
    ↪ 0.7527,−0.039,0.634,0.8663,0.1592,0.072,0.5139,0.8527,0.572,
    ↪ 0.6298,0.1723,0.072,0.7612,0.0602,0.572,−0.0645'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP112'

_symmetry_space_group_name_H-M "P 21/n 21/m 21/a"
_symmetry_Int_Tables_number 62

_cell_length_a     11.79700
_cell_length_b      9.73900
_cell_length_c     26.27000
_cell_angle_alpha  90.00000
```

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_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x,y+1/2,-z
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x,-y+1/2,z
8 x+1/2,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
P1 P 4 c 0.08600 0.25000 0.62700 1.00000
P2 P 4 c 0.04100 0.25000 0.50000 1.00000
P3 P 4 c -0.03400 0.25000 0.28200 1.00000
P4 P 4 c 0.23800 0.25000 0.33100 1.00000
P5 P 4 c -0.04800 0.25000 0.82500 1.00000
P6 P 4 c 0.23900 0.25000 0.80700 1.00000
P7 P 4 c 0.19200 0.25000 0.10400 1.00000
P8 P 4 c -0.03200 0.25000 0.01500 1.00000
Se1 Se 4 c -0.05700 0.25000 0.57100 1.00000
Se2 Se 4 c 0.01480 0.25000 0.25670 1.00000
Se3 Se 4 c 0.10700 0.25000 0.87000 1.00000
Se4 Se 4 c 0.16300 0.25000 0.01800 1.00000
P9 P 8 d 0.22700 0.13400 0.58750 1.00000
P10 P 8 d 0.30400 0.63400 0.64430 1.00000
P11 P 8 d -0.01000 0.13400 0.75270 1.00000
P12 P 8 d -0.03900 0.63400 0.86630 1.00000
Se5 Se 8 d 0.15920 0.07200 0.51390 1.00000
Se6 Se 8 d 0.85270 0.57200 0.62980 1.00000
Se7 Se 8 d 0.17230 0.07200 0.76120 1.00000
Se8 Se 8 d 0.06020 0.57200 -0.06450 1.00000
```

P_4Se_3: A4B3_oP112_62_8c4d_4c4d - POSCAR

```
A4B3_oP112_62_8c4d_4c4d & a,b/a,c/a,x1,z1,x2,z2,x3,z3,x4,z4,x5,z5,x6,z6,
    ↪ x7,z7,x8,z8,x9,z9,x10,z10,x11,z11,x12,z12,x13,y13,z13,x14,y14,
    ↪ z14,x15,y15,z15,x16,y16,z16,x17,y17,z17,x18,y18,z18,x19,y19,z19
    ↪ ,x20,y20,z20 --params=11.797,0.825548868356,2.22683733152,0.086
    ↪ ,0.627,0.041,0.5,-0.034,0.282,0.238,0.331,-0.048,0.825,0.239,
    ↪ 0.807,0.192,0.104,-0.032,0.015,-0.057,0.571,0.0148,0.2567,0.107
    ↪ ,0.87,0.163,0.018,0.227,0.134,0.5875,0.304,0.634,0.6443,-0.01,
    ↪ 0.134,0.7527,-0.039,0.634,0.8663,0.1592,0.072,0.5139,0.8527,
    ↪ 0.572,0.6298,0.1723,0.072,0.7612,0.0602,0.572,-0.0645 & Pnma D_
    ↪ {2h}^{16} #62 (c^12d^8) & oP112 & None & P4Se3 & P4Se3 & E.
    ↪ Keulen and A. Vos, Acta Cryst. 12, 323-329 (1959)
   1.00000000000000
  11.79700000000000    0.00000000000000    0.00000000000000
   0.00000000000000    9.73900000000000    0.00000000000000
   0.00000000000000    0.00000000000000   26.27000000000000
     P     Se
    64     48
Direct
   0.08600000000000    0.25000000000000    0.62700000000000    P    (4c)
   0.41400000000000    0.75000000000000    1.12700000000000    P    (4c)
  -0.08600000000000    0.75000000000000   -0.62700000000000    P    (4c)
   0.58600000000000    0.25000000000000   -0.12700000000000    P    (4c)
   0.04100000000000    0.25000000000000    0.50000000000000    P    (4c)
   0.45900000000000    0.75000000000000    1.00000000000000    P    (4c)
  -0.04100000000000    0.75000000000000   -0.50000000000000    P    (4c)
   0.54100000000000    0.25000000000000    0.00000000000000    P    (4c)
  -0.03400000000000    0.25000000000000    0.28200000000000    P    (4c)
   0.53400000000000    0.75000000000000    0.78200000000000    P    (4c)
   0.03400000000000    0.75000000000000   -0.28200000000000    P    (4c)
   0.46600000000000    0.25000000000000    0.21800000000000    P    (4c)
   0.23800000000000    0.25000000000000    0.33100000000000    P    (4c)
   0.26200000000000    0.75000000000000    0.83100000000000    P    (4c)
  -0.23800000000000    0.75000000000000   -0.33100000000000    P    (4c)
   0.73800000000000    0.25000000000000    0.16900000000000    P    (4c)
  -0.04800000000000    0.25000000000000    0.82500000000000    P    (4c)
   0.54800000000000    0.75000000000000    1.32500000000000    P    (4c)
   0.04800000000000    0.75000000000000   -0.82500000000000    P    (4c)
   0.45200000000000    0.25000000000000   -0.32500000000000    P    (4c)
   0.23900000000000    0.25000000000000    0.80700000000000    P    (4c)
   0.26100000000000    0.75000000000000    1.30700000000000    P    (4c)
  -0.23900000000000    0.75000000000000   -0.80700000000000    P    (4c)
   0.73900000000000    0.25000000000000   -0.30700000000000    P    (4c)
   0.19200000000000    0.25000000000000    0.10400000000000    P    (4c)
   0.30800000000000    0.75000000000000    0.60400000000000    P    (4c)
  -0.19200000000000    0.75000000000000   -0.10400000000000    P    (4c)
   0.69200000000000    0.25000000000000    0.39600000000000    P    (4c)
  -0.03200000000000    0.25000000000000    0.01500000000000    P    (4c)
   0.53200000000000    0.75000000000000    0.51500000000000    P    (4c)
   0.03200000000000    0.75000000000000   -0.01500000000000    P    (4c)
   0.46800000000000    0.25000000000000    0.48500000000000    P    (4c)
   0.22700000000000    0.13400000000000    0.58750000000000    P    (8d)
   0.27300000000000   -0.13400000000000    1.08750000000000    P    (8d)
  -0.22700000000000    0.63400000000000   -0.58750000000000    P    (8d)
   0.72700000000000    0.36600000000000   -0.08750000000000    P    (8d)
  -0.22700000000000   -0.13400000000000   -0.58750000000000    P    (8d)
   0.72700000000000    0.13400000000000   -0.08750000000000    P    (8d)
   0.22700000000000    0.36600000000000    0.58750000000000    P    (8d)
   0.27300000000000    0.63400000000000    1.08750000000000    P    (8d)
   0.30400000000000    0.63400000000000    0.64430000000000    P    (8d)
   0.19600000000000   -0.63400000000000    1.14430000000000    P    (8d)
  -0.30400000000000    1.13400000000000   -0.64430000000000    P    (8d)
   0.80400000000000   -0.13400000000000   -0.14430000000000    P    (8d)
  -0.30400000000000   -0.63400000000000   -0.64430000000000    P    (8d)
   0.80400000000000    0.63400000000000   -0.14430000000000    P    (8d)
   0.30400000000000   -0.13400000000000    0.64430000000000    P    (8d)
   0.19600000000000    1.13400000000000    1.14430000000000    P    (8d)
  -0.01000000000000    0.13400000000000    0.75270000000000    P    (8d)
   0.51000000000000   -0.13400000000000    1.25270000000000    P    (8d)
   0.01000000000000    0.63400000000000   -0.75270000000000    P    (8d)
   0.49000000000000    0.36600000000000   -0.25270000000000    P    (8d)
   0.01000000000000   -0.13400000000000   -0.75270000000000    P    (8d)
   0.49000000000000    0.13400000000000   -0.25270000000000    P    (8d)
  -0.01000000000000    0.36600000000000    0.75270000000000    P    (8d)
   0.51000000000000    0.63400000000000    1.25270000000000    P    (8d)
  -0.03900000000000    0.63400000000000    0.86630000000000    P    (8d)
   0.53900000000000   -0.63400000000000    1.36630000000000    P    (8d)
   0.03900000000000    1.13400000000000   -0.86630000000000    P    (8d)
   0.46100000000000   -0.13400000000000   -0.36630000000000    P    (8d)
   0.03900000000000   -0.63400000000000   -0.86630000000000    P    (8d)
   0.46100000000000    0.63400000000000   -0.36630000000000    P    (8d)
  -0.03900000000000   -0.13400000000000    0.86630000000000    P    (8d)
   0.53900000000000    1.13400000000000    1.36630000000000    P    (8d)
  -0.05700000000000    0.25000000000000    0.57100000000000   Se    (4c)
   0.55700000000000    0.75000000000000    1.07100000000000   Se    (4c)
   0.05700000000000    0.75000000000000   -0.57100000000000   Se    (4c)
   0.44300000000000    0.25000000000000   -0.07100000000000   Se    (4c)
   0.01480000000000    0.25000000000000    0.25670000000000   Se    (4c)
   0.48520000000000    0.75000000000000    0.75670000000000   Se    (4c)
  -0.01480000000000    0.75000000000000   -0.25670000000000   Se    (4c)
   0.51480000000000    0.25000000000000    0.24330000000000   Se    (4c)
   0.10700000000000    0.25000000000000    0.87000000000000   Se    (4c)
   0.39300000000000    0.75000000000000    1.37000000000000   Se    (4c)
  -0.10700000000000    0.75000000000000   -0.87000000000000   Se    (4c)
   0.60700000000000    0.25000000000000   -0.37000000000000   Se    (4c)
   0.16300000000000    0.25000000000000    0.01800000000000   Se    (4c)
   0.33700000000000    0.75000000000000    0.51800000000000   Se    (4c)
  -0.16300000000000    0.75000000000000   -0.01800000000000   Se    (4c)
   0.66300000000000    0.25000000000000    0.48200000000000   Se    (4c)
   0.15920000000000    0.07200000000000    0.51390000000000   Se    (8d)
   0.34080000000000   -0.07200000000000    1.01390000000000   Se    (8d)
  -0.15920000000000    0.57200000000000   -0.51390000000000   Se    (8d)
   0.65920000000000    0.42800000000000   -0.01390000000000   Se    (8d)
  -0.15920000000000   -0.07200000000000   -0.51390000000000   Se    (8d)
   0.65920000000000    0.07200000000000   -0.01390000000000   Se    (8d)
   0.15920000000000    0.42800000000000    0.51390000000000   Se    (8d)
   0.34080000000000    0.57200000000000    1.01390000000000   Se    (8d)
   0.85270000000000    0.57200000000000    0.62980000000000   Se    (8d)
  -0.35270000000000   -0.57200000000000    1.12980000000000   Se    (8d)
  -0.85270000000000    1.07200000000000   -0.62980000000000   Se    (8d)
   1.35270000000000   -0.07200000000000   -0.12980000000000   Se    (8d)
  -0.85270000000000   -0.57200000000000   -0.62980000000000   Se    (8d)
   1.35270000000000    0.57200000000000   -0.12980000000000   Se    (8d)
   0.85270000000000   -0.07200000000000    0.62980000000000   Se    (8d)
  -0.35270000000000    1.07200000000000    1.12980000000000   Se    (8d)
   0.17230000000000    0.07200000000000    0.76120000000000   Se    (8d)
   0.32770000000000   -0.07200000000000    1.26120000000000   Se    (8d)
  -0.17230000000000    0.57200000000000   -0.76120000000000   Se    (8d)
   0.67230000000000    0.42800000000000   -0.26120000000000   Se    (8d)
  -0.17230000000000   -0.07200000000000   -0.76120000000000   Se    (8d)
   0.67230000000000    0.07200000000000   -0.26120000000000   Se    (8d)
   0.17230000000000    0.42800000000000    0.76120000000000   Se    (8d)
   0.32770000000000    0.57200000000000    1.26120000000000   Se    (8d)
   0.06020000000000    0.57200000000000   -0.06450000000000   Se    (8d)
   0.43980000000000   -0.57200000000000    0.43550000000000   Se    (8d)
  -0.06020000000000    1.07200000000000    0.06450000000000   Se    (8d)
   0.56020000000000   -0.07200000000000    0.56450000000000   Se    (8d)
  -0.06020000000000   -0.57200000000000    0.06450000000000   Se    (8d)
   0.56020000000000    0.57200000000000    0.56450000000000   Se    (8d)
   0.06020000000000   -0.07200000000000   -0.06450000000000   Se    (8d)
   0.43980000000000    1.07200000000000    0.43550000000000   Se    (8d)
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Mo_4P_3: A4B3_oP56_62_8c_6c - CIF

```
# CIF file
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_audit_creation_method FINDSYM

_chemical_name_mineral 'Mo4P3'
_chemical_formula_sum 'Mo4 P3'

loop_
_publ_author_name
 'S. Rundqvist'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 19
_journal_year 1965
_journal_page_first 393
_journal_page_last 400
_publ_Section_title
;
 The Crystal Structure of Mo$_{4}$P$_{3}$
;

_aflow_title 'Mo$_{4}$P$_{3}$ Structure'
_aflow_proto 'A4B3_oP56_62_8c_6c'
_aflow_params 'a,b/a,c/a,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3},x_{4},z_{4
    ↪ },x_{5},z_{5},x_{6},z_{6},x_{7},z_{7},x_{8},z_{8},x_{9},z_{9},
    ↪ x_{10},z_{10},x_{11},z_{11},x_{12},z_{12},x_{13},z_{13},x_{14},
    ↪ z_{14}'
_aflow_params_values '12.428,0.254103636949,1.64467331831,0.1939,0.2262,
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```
      ↪ 0.3843 , 0.8518 , 0.1218 , 0.6891 , 0.298 , 0.0835 , 0.4007 , 0.5269 , 0.0952 ,
      ↪ 0.4665 , 0.2977 , 0.3863 , 0.0095 , 0.8272 , 0.2464 , -0.027 , 0.4824 , 0.4186 ,
      ↪ 0.3206 , 0.6371 , 0.0382 , 0.5805 , 0.0155 , 0.2745 , 0.1994 , 0.7941'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP56'

_symmetry_space_group_name_H-M "P 21/n 21/m 21/a"
_symmetry_Int_Tables_number 62

_cell_length_a      12.42800
_cell_length_b      3.15800
_cell_length_c      20.44000
_cell_angle_alpha   90.00000
_cell_angle_beta    90.00000
_cell_angle_gamma   90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x,y+1/2,-z
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x,-y+1/2,z
8 x+1/2,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Mo1 Mo 4 c 0.19390 0.25000 0.22620 1.00000
Mo2 Mo 4 c 0.38430 0.25000 0.85180 1.00000
Mo3 Mo 4 c 0.12180 0.25000 0.68910 1.00000
Mo4 Mo 4 c 0.29800 0.25000 0.08350 1.00000
Mo5 Mo 4 c 0.40070 0.25000 0.52690 1.00000
Mo6 Mo 4 c 0.09520 0.25000 0.46650 1.00000
Mo7 Mo 4 c 0.29770 0.25000 0.38630 1.00000
Mo8 Mo 4 c 0.00950 0.25000 0.82720 1.00000
P1 P 4 c 0.24640 0.25000 -0.02700 1.00000
P2 P 4 c 0.48240 0.25000 0.41860 1.00000
P3 P 4 c 0.32060 0.25000 0.63710 1.00000
P4 P 4 c 0.03820 0.25000 0.58050 1.00000
P5 P 4 c 0.01550 0.25000 0.27450 1.00000
P6 P 4 c 0.19940 0.25000 0.79410 1.00000
```

Mo_4P_3: A4B3_oP56_62_8c_6c - POSCAR

```
A4B3_oP56_62_8c_6c & a,b/a,c/a,x1,z1,x2,z2,x3,z3,x4,z4,x5,z5,x6,z6,x7,z7
      ↪ ,x8,z8,x9,z9,x10,z10,x11,z11,x12,z12,x13,z13,x14,z14 --params=
      ↪ 12.428,0.254103636949,1.64467331831,0.1939,0.2262,0.3843,0.8518
      ↪ ,0.1218,0.6891,0.298,0.0835,0.4007,0.5269,0.0952,0.4665,0.2977,
      ↪ 0.3863,0.0095,0.8272,0.2464,-0.027,0.4824,0.4186,0.3206,0.6371,
      ↪ 0.0382,0.5805,0.0155,0.2745,0.1994,0.7941 & Pnma D_{2h}^{16} #
      ↪ 62 (c^14) & oP56 & None & Mo4P3 & Mo4P3 & S. Rundqvist, Acta
      ↪ Chem. Scand. 19, 393-400 (1965)
   1.00000000000000
  12.42800000000000    0.00000000000000    0.00000000000000
   0.00000000000000    3.15800000000000    0.00000000000000
   0.00000000000000    0.00000000000000   20.44000000000000
    Mo      P
    32     24
Direct
   0.19390000000000    0.25000000000000    0.22620000000000   Mo   (4c)
   0.30610000000000    0.75000000000000    0.72620000000000   Mo   (4c)
  -0.19390000000000    0.75000000000000   -0.22620000000000   Mo   (4c)
   0.69390000000000    0.25000000000000    0.27380000000000   Mo   (4c)
   0.38430000000000    0.25000000000000    0.85180000000000   Mo   (4c)
   0.11570000000000    0.75000000000000    1.35180000000000   Mo   (4c)
  -0.38430000000000    0.75000000000000   -0.85180000000000   Mo   (4c)
   0.88430000000000    0.25000000000000   -0.35180000000000   Mo   (4c)
   0.12180000000000    0.25000000000000    0.68910000000000   Mo   (4c)
   0.37820000000000    0.75000000000000    1.18910000000000   Mo   (4c)
  -0.12180000000000    0.75000000000000   -0.68910000000000   Mo   (4c)
   0.62180000000000    0.25000000000000   -0.18910000000000   Mo   (4c)
   0.29800000000000    0.25000000000000    0.08350000000000   Mo   (4c)
   0.20200000000000    0.75000000000000    0.58350000000000   Mo   (4c)
  -0.29800000000000    0.75000000000000   -0.08350000000000   Mo   (4c)
   0.79800000000000    0.25000000000000    0.41650000000000   Mo   (4c)
   0.40070000000000    0.25000000000000    0.52690000000000   Mo   (4c)
   0.09930000000000    0.75000000000000    1.02690000000000   Mo   (4c)
  -0.40070000000000    0.75000000000000   -0.52690000000000   Mo   (4c)
   0.90070000000000    0.25000000000000   -0.02690000000000   Mo   (4c)
   0.09520000000000    0.25000000000000    0.46650000000000   Mo   (4c)
   0.40480000000000    0.75000000000000    0.96650000000000   Mo   (4c)
  -0.09520000000000    0.75000000000000   -0.46650000000000   Mo   (4c)
   0.59520000000000    0.25000000000000    0.03350000000000   Mo   (4c)
   0.29770000000000    0.25000000000000    0.38630000000000   Mo   (4c)
   0.20230000000000    0.75000000000000    0.88630000000000   Mo   (4c)
  -0.29770000000000    0.75000000000000   -0.38630000000000   Mo   (4c)
   0.79770000000000    0.25000000000000    0.11370000000000   Mo   (4c)
   0.00950000000000    0.25000000000000    0.82720000000000   Mo   (4c)
   0.49050000000000    0.75000000000000    1.32720000000000   Mo   (4c)
  -0.00950000000000    0.75000000000000   -0.82720000000000   Mo   (4c)
   0.50950000000000    0.25000000000000   -0.32720000000000   Mo   (4c)
   0.24640000000000    0.25000000000000   -0.02700000000000    P   (4c)
   0.25360000000000    0.75000000000000    0.47300000000000    P   (4c)
  -0.24640000000000    0.75000000000000    0.02700000000000    P   (4c)
   0.74640000000000    0.25000000000000    0.52700000000000    P   (4c)
   0.48240000000000    0.25000000000000    0.41860000000000    P   (4c)
   0.01760000000000    0.75000000000000    0.91860000000000    P   (4c)
  -0.48240000000000    0.75000000000000   -0.41860000000000    P   (4c)
   0.98240000000000    0.25000000000000    0.08140000000000    P   (4c)
   0.32060000000000    0.25000000000000    0.63710000000000    P   (4c)
   0.17940000000000    0.75000000000000    1.13710000000000    P   (4c)
  -0.32060000000000    0.75000000000000   -0.63710000000000    P   (4c)
   0.82060000000000    0.25000000000000   -0.13710000000000    P   (4c)
   0.03820000000000    0.25000000000000    0.58050000000000    P   (4c)
   0.46180000000000    0.75000000000000    1.08050000000000    P   (4c)
  -0.03820000000000    0.75000000000000   -0.58050000000000    P   (4c)
   0.53820000000000    0.25000000000000   -0.08050000000000    P   (4c)
   0.01550000000000    0.25000000000000    0.27450000000000    P   (4c)
   0.48450000000000    0.75000000000000    0.77450000000000    P   (4c)
  -0.01550000000000    0.75000000000000   -0.27450000000000    P   (4c)
   0.51550000000000    0.25000000000000    0.22550000000000    P   (4c)
   0.19940000000000    0.25000000000000    0.79410000000000    P   (4c)
   0.30060000000000    0.75000000000000    1.29410000000000    P   (4c)
  -0.19940000000000    0.75000000000000   -0.79410000000000    P   (4c)
   0.69940000000000    0.25000000000000   -0.29410000000000    P   (4c)
```

$K_2SnCl_4{\cdot}H_2O$ ($E3_5$): A4BC2D_oP32_62_2cd_b_2c_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Cl4(H2O)K2Sn'
_chemical_formula_sum 'Cl4 (H2O) K2 Sn'

loop_
_publ_author_name
 'H. Brasseur'
 'A. {de Rassenfosse}'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 101
_journal_year 1939
_journal_page_first 389
_journal_page_last 395
_publ_Section_title
;
 The Crystal Structure of Hydrated Potassium Chlorostannite
;

# Found in Strukturbericht Band VII 1939, 1943

_aflow_title 'K$_{2}$SnCl$_{4}\cdot H$_{2}$O ($E3_{5}$) Structure'
_aflow_proto 'A4BC2D_oP32_62_2cd_b_2c_a'
_aflow_params 'a,b/a,c/a,x_{3},z_{3},x_{4},z_{4},x_{5},z_{5},x_{6},z_{6
      ↪ },x_{7},y_{7},z_{7}'
_aflow_params_values '12.05,0.755186721992,0.684647302905,-0.05,0.21,
      ↪ 0.05,0.79,0.2,0.34,0.825,0.6,0.25,-0.04,0.12'
_aflow_Strukturbericht '$E3_{5}$'
_aflow_Pearson 'oP32'

_symmetry_space_group_name_H-M "P 21/n 21/m 21/a"
_symmetry_Int_Tables_number 62

_cell_length_a      12.05000
_cell_length_b      9.10000
_cell_length_c      8.25000
_cell_angle_alpha   90.00000
_cell_angle_beta    90.00000
_cell_angle_gamma   90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x,y+1/2,-z
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x,-y+1/2,z
8 x+1/2,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Sn1 Sn 4 a 0.00000 0.00000 0.00000 1.00000
H2O1 H2O 4 b 0.00000 0.00000 0.50000 1.00000
Cl1 Cl 4 c -0.05000 0.25000 0.21000 1.00000
Cl2 Cl 4 c 0.05000 0.25000 0.79000 1.00000
K1 K 4 c 0.20000 0.25000 0.34000 1.00000
K2 K 4 c 0.82500 0.25000 0.60000 1.00000
Cl3 Cl 8 d 0.25000 -0.04000 0.12000 1.00000
```

$K_2SnCl_4{\cdot}H_2O$ ($E3_5$): A4BC2D_oP32_62_2cd_b_2c_a - POSCAR

```
A4BC2D_oP32_62_2cd_b_2c_a & a,b/a,c/a,x3,z3,x4,z4,x5,z5,x6,z6,x7,y7,z7
      ↪ --params=12.05,0.755186721992,0.684647302905,-0.05,0.21,0.05,
      ↪ 0.79,0.2,0.34,0.825,0.6,0.25,-0.04,0.12 & Pnma D_{2h}^{16} #62
      ↪ (abc^4d) & oP32 & $E3_{5}$ & Cl4(H2O)K2Sn & Cl4(H2O)K2Sn & H.
      ↪ Brasseur and A. {de Rassenfosse}, Zeitschrift f{"u}r
      ↪ Kristallographie - Crystalline Materials 101, 389-395 (1939)
```

```
   1.00000000000000
  12.05000000000000    0.00000000000000    0.00000000000000
   0.00000000000000    9.10000000000000    0.00000000000000
   0.00000000000000    0.00000000000000    8.25000000000000
    Cl    H2O     K     Sn
    16     4     8     4
Direct
  -0.05000000000000    0.25000000000000    0.21000000000000   Cl    (4c)
   0.55000000000000    0.75000000000000    0.71000000000000   Cl    (4c)
   0.05000000000000    0.75000000000000   -0.21000000000000   Cl    (4c)
   0.45000000000000    0.25000000000000    0.29000000000000   Cl    (4c)
   0.05000000000000    0.25000000000000    0.79000000000000   Cl    (4c)
   0.45000000000000    0.75000000000000    1.29000000000000   Cl    (4c)
  -0.05000000000000    0.75000000000000   -0.79000000000000   Cl    (4c)
   0.55000000000000    0.25000000000000   -0.29000000000000   Cl    (4c)
   0.25000000000000   -0.04000000000000    0.12000000000000   Cl    (8d)
   0.25000000000000    0.04000000000000    0.62000000000000   Cl    (8d)
  -0.25000000000000    0.46000000000000   -0.12000000000000   Cl    (8d)
   0.75000000000000    0.54000000000000    0.38000000000000   Cl    (8d)
  -0.25000000000000    0.04000000000000   -0.12000000000000   Cl    (8d)
   0.75000000000000   -0.04000000000000    0.38000000000000   Cl    (8d)
   0.25000000000000    0.54000000000000    0.12000000000000   Cl    (8d)
   0.25000000000000    0.46000000000000    0.62000000000000   Cl    (8d)
   0.00000000000000    0.00000000000000    0.50000000000000   H2O   (4b)
   0.50000000000000    0.00000000000000    0.00000000000000   H2O   (4b)
   0.00000000000000    0.50000000000000    0.50000000000000   H2O   (4b)
   0.50000000000000    0.50000000000000    0.00000000000000   H2O   (4b)
   0.20000000000000    0.25000000000000    0.34000000000000   K     (4c)
   0.30000000000000    0.75000000000000    0.84000000000000   K     (4c)
  -0.20000000000000    0.75000000000000   -0.34000000000000   K     (4c)
   0.70000000000000    0.25000000000000    0.16000000000000   K     (4c)
   0.82500000000000    0.25000000000000    0.60000000000000   K     (4c)
  -0.32500000000000    0.75000000000000    1.10000000000000   K     (4c)
  -0.82500000000000    0.75000000000000   -0.60000000000000   K     (4c)
   1.32500000000000    0.25000000000000   -0.10000000000000   K     (4c)
   0.00000000000000    0.00000000000000    0.00000000000000   Sn    (4a)
   0.50000000000000    0.00000000000000    0.50000000000000   Sn    (4a)
   0.00000000000000    0.50000000000000    0.00000000000000   Sn    (4a)
   0.50000000000000    0.50000000000000    0.50000000000000   Sn    (4a)
```

$K_2SnCl_4{\cdot}H_2O$: A4BC2D_oP32_62_2cd_c_d_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Cl4(H2O)K2Sn'
_chemical_formula_sum 'Cl4 (H2O) K2 Sn'

loop_
_publ_author_name
 'B. Kamenar'
 'D. Grdeni\\'{c}'
_journal_name_full_name
;
 Journal of Inorganic and Nuclear Chemistry
;
_journal_volume 24
_journal_year 1962
_journal_page_first 1039
_journal_page_last 1045
_publ_Section_title
;
 The crystal structure of potassium chloride trichlorostannite hydrate,
      ↪ KCl, KSnCl$_{3}$, H$_{2}$O
;

_aflow_title 'K$_{2}$SnCl$_{4}$$\cdot$H$_{2}$O Structure'
_aflow_proto 'A4BC2D_oP32_62_2cd_c_d_c'
_aflow_params 'a,b/a,c/a,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3},x_{4},z_{4
      ↪ },x_{5},y_{5},z_{5},x_{6},y_{6},z_{6}'
_aflow_params_values '12.05,0.758506224066,0.683817427386,0.308,0.888,
      ↪ 0.766,0.099,0.54,0.441,0.512,0.0,0.444,-0.052,0.207,0.685,0.482
      ↪ ,0.37'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP32'

_symmetry_space_group_name_H-M "P 21/n 21/m 21/a"
_symmetry_Int_Tables_number 62

_cell_length_a    12.05000
_cell_length_b    9.14000
_cell_length_c    8.24000
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x,y+1/2,-z
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x,-y+1/2,z
8 x+1/2,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cl1 Cl 4 c 0.30800 0.25000 0.88800 1.00000
Cl2 Cl 4 c 0.76600 0.25000 0.09900 1.00000
H2O1 H2O 4 c 0.54000 0.25000 0.44100 1.00000
Sn1 Sn 4 c 0.51200 0.25000 0.00000 1.00000
Cl3 Cl 8 d 0.44400 -0.05200 0.20700 1.00000
K1 K 8 d 0.68500 0.48200 0.37000 1.00000
```

$K_2SnCl_4{\cdot}H_2O$: A4BC2D_oP32_62_2cd_c_d_c - POSCAR

```
A4BC2D_oP32_62_2cd_c_d_c & a,b/a,c/a,x1,z1,x2,z2,x3,z3,x4,z4,x5,y5,z5,x6
      ↪ ,y6,z6 --params=12.05,0.758506224066,0.683817427386,0.308,0.888
      ↪ ,0.766,0.099,0.54,0.441,0.512,0.0,0.444,-0.052,0.207,0.685,
      ↪ 0.482,0.37 & Pnma D_{2h}^{16} #62 (c^4d^2) & oP32 & None & Cl4(
      ↪ H2O)K2Sn & Cl4(H2O)K2Sn & B. Kamenar and D. Grdeni\'{c}, J.
      ↪ Inorg. Nucl. Chem. 24, 1039-1045 (1962)
   1.00000000000000
  12.05000000000000    0.00000000000000    0.00000000000000
   0.00000000000000    9.14000000000000    0.00000000000000
   0.00000000000000    0.00000000000000    8.24000000000000
    Cl    H2O     K     Sn
    16     4     8     4
Direct
   0.30800000000000    0.25000000000000    0.88800000000000   Cl    (4c)
   0.19200000000000    0.75000000000000    1.38800000000000   Cl    (4c)
  -0.30800000000000    0.75000000000000   -0.88800000000000   Cl    (4c)
   0.80800000000000    0.25000000000000   -0.38800000000000   Cl    (4c)
   0.76600000000000    0.25000000000000    0.09900000000000   Cl    (4c)
  -0.26600000000000    0.75000000000000    0.59900000000000   Cl    (4c)
  -0.76600000000000    0.75000000000000   -0.09900000000000   Cl    (4c)
   1.26600000000000    0.25000000000000    0.40100000000000   Cl    (4c)
   0.44400000000000   -0.05200000000000    0.20700000000000   Cl    (8d)
   0.05600000000000    0.05200000000000    0.70700000000000   Cl    (8d)
  -0.44400000000000    0.44800000000000   -0.20700000000000   Cl    (8d)
   0.94400000000000    0.55200000000000    0.29300000000000   Cl    (8d)
  -0.44400000000000    0.05200000000000   -0.20700000000000   Cl    (8d)
   0.94400000000000   -0.05200000000000    0.29300000000000   Cl    (8d)
   0.44400000000000    0.55200000000000    0.20700000000000   Cl    (8d)
   0.05600000000000    0.44800000000000    0.70700000000000   Cl    (8d)
   0.54000000000000    0.25000000000000    0.44100000000000   H2O   (4c)
  -0.04000000000000    0.75000000000000    0.94100000000000   H2O   (4c)
  -0.54000000000000    0.75000000000000   -0.44100000000000   H2O   (4c)
   1.04000000000000    0.25000000000000    0.05900000000000   H2O   (4c)
   0.68500000000000    0.48200000000000    0.37000000000000   K     (8d)
  -0.18500000000000   -0.48200000000000    0.87000000000000   K     (8d)
  -0.68500000000000    0.98200000000000   -0.37000000000000   K     (8d)
   1.18500000000000    0.01800000000000    0.13000000000000   K     (8d)
  -0.68500000000000   -0.48200000000000   -0.37000000000000   K     (8d)
   1.18500000000000    0.48200000000000    0.13000000000000   K     (8d)
   0.68500000000000    0.01800000000000    0.37000000000000   K     (8d)
  -0.18500000000000    0.98200000000000    0.87000000000000   K     (8d)
   0.51200000000000    0.25000000000000    0.00000000000000   Sn    (4c)
  -0.01200000000000    0.75000000000000    0.50000000000000   Sn    (4c)
  -0.51200000000000    0.75000000000000    0.00000000000000   Sn    (4c)
   1.01200000000000    0.25000000000000    0.50000000000000   Sn    (4c)
```

$VOSO_4$: A5BC_oP28_62_3cd_c_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'O5SV'
_chemical_formula_sum 'O5 S V'

loop_
_publ_author_name
 'P. Kierkegaard'
 'J. M. Longo'
 'B.-O. Marinder'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 19
_journal_year 1965
_journal_page_first 763
_journal_page_last 764
_publ_Section_title
;
 Note on the Crystal Structure of VOSO$_{4}$
;

_aflow_title 'VOSO$_{4}$ Structure'
_aflow_proto 'A5BC_oP28_62_3cd_c_c'
_aflow_params 'a,b/a,c/a,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3},x_{4},z_{4
      ↪ },x_{5},z_{5},x_{6},y_{6},z_{6}'
_aflow_params_values '7.371,0.850495183829,0.960792294126,0.7187,-0.0104
      ↪ ,0.0399,-0.0201,0.3719,0.1647,0.8761,0.8669,0.1658,0.2327,
      ↪ 0.1251,0.5733,0.2569'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP28'

_symmetry_space_group_name_H-M "P 21/n 21/m 21/a"
_symmetry_Int_Tables_number 62

_cell_length_a    7.37100
_cell_length_b    6.26900
_cell_length_c    7.08200
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000
```

```
loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x,y+1/2,-z
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x,-y+1/2,z
8 x+1/2,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 4 c 0.71870 0.25000 -0.01040 1.00000
O2 O 4 c 0.03990 0.25000 -0.02010 1.00000
O3 O 4 c 0.37190 0.25000 0.16470 1.00000
S1 S 4 c 0.87610 0.25000 0.86690 1.00000
V1 V 4 c 0.16580 0.25000 0.23270 1.00000
O4 O 8 d 0.12510 0.57330 0.25690 1.00000
```

$VOSO_4$: A5BC_oP28_62_3cd_c_c - POSCAR

```
A5BC_oP28_62_3cd_c_c & a,b/a,c/a,x1,z1,x2,z2,x3,z3,x4,z4,x5,z5,x6,y6,z6
      ↪ --params=7.371,0.850495183829,0.960792294126,0.7187,-0.0104,
      ↪ 0.0399,-0.0201,0.3719,0.1647,0.8761,0.8669,0.1658,0.2327,0.1251
      ↪ ,0.5733,0.2569 & Pnma D_{2h}^{16} #62 (c^5d) & oP28 & None &
      ↪ O5SV & O5SV & P. Kierkegaard and J. M. Longo and B.-O. Marinder
      ↪ , Acta Chem. Scand. 19, 763-764 (1965)
   1.00000000000000
   7.37100000000000    0.00000000000000    0.00000000000000
   0.00000000000000    6.26900000000000    0.00000000000000
   0.00000000000000    0.00000000000000    7.08200000000000
     O     S     V
    20     4     4
Direct
   0.71870000000000    0.25000000000000   -0.01040000000000    O    (4c)
  -0.21870000000000    0.75000000000000    0.48960000000000    O    (4c)
  -0.71870000000000    0.75000000000000    0.01040000000000    O    (4c)
   1.21870000000000    0.25000000000000    0.51040000000000    O    (4c)
   0.03990000000000    0.25000000000000   -0.02010000000000    O    (4c)
   0.46010000000000    0.75000000000000    0.47990000000000    O    (4c)
  -0.03990000000000    0.75000000000000    0.02010000000000    O    (4c)
   0.53990000000000    0.25000000000000    0.52010000000000    O    (4c)
   0.37190000000000    0.25000000000000    0.16470000000000    O    (4c)
   0.12810000000000    0.75000000000000    0.66470000000000    O    (4c)
  -0.37190000000000    0.75000000000000   -0.16470000000000    O    (4c)
   0.87190000000000    0.25000000000000    0.33530000000000    O    (4c)
   0.12510000000000    0.57330000000000    0.25690000000000    O    (8d)
   0.37490000000000   -0.57330000000000    0.75690000000000    O    (8d)
  -0.12510000000000    1.07330000000000   -0.25690000000000    O    (8d)
   0.62510000000000   -0.07330000000000    0.24310000000000    O    (8d)
  -0.12510000000000   -0.57330000000000   -0.25690000000000    O    (8d)
   0.62510000000000    0.57330000000000    0.24310000000000    O    (8d)
   0.12510000000000   -0.07330000000000    0.25690000000000    O    (8d)
   0.37490000000000    1.07330000000000    0.75690000000000    O    (8d)
   0.87610000000000    0.25000000000000    0.86690000000000    S    (4c)
  -0.37610000000000    0.75000000000000    1.36690000000000    S    (4c)
  -0.87610000000000    0.75000000000000   -0.86690000000000    S    (4c)
   1.37610000000000    0.25000000000000   -0.36690000000000    S    (4c)
   0.16580000000000    0.25000000000000    0.23270000000000    V    (4c)
   0.33420000000000    0.75000000000000    0.73270000000000    V    (4c)
  -0.16580000000000    0.75000000000000   -0.23270000000000    V    (4c)
   0.66580000000000    0.25000000000000    0.26730000000000    V    (4c)
```

Possible δ-$Y_2Si_2O_7$: A7B2C2_oP44_62_3c2d_2c_d - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'O7Si2Y2'
_chemical_formula_sum 'O7 Si2 Y2'

loop_
_publ_author_name
 'H. W. Dias'
 'F. P. Glasser'
 'R. P. Gunwardane'
 'R. A. Howie'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 191
_journal_year 1990
_journal_page_first 117
_journal_page_last 124
_publ_Section_title
;
 The crystal structure of $\delta$-yttrium pyrosilicate, $\delta$-Y$_{2}
      ↪ $Si$_{2}$O$_{7}$
;

# Found in Revision of the crystallographic data of polymorphic Y$_{2}
      ↪ $Si$_{2}$O$_{7}$ and Y$_{2}$SiO$_{5}$ compounds, 2004

_aflow_title 'Possible $\delta$-Y$_{2}$Si$_{2}$O$_{7}$ Structure'
_aflow_proto 'A7B2C2_oP44_62_3c2d_2c_d'
_aflow_params 'a,b/a,c/a,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3},x_{4},z_{4
      ↪ },x_{5},z_{5},x_{6},y_{6},z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},
      ↪ z_{8}'
_aflow_params_values '13.655,0.596045404614,0.367337971439,0.1524,0.0613
      ↪ ,0.0782,0.5588,0.8969,0.3386,0.1821,0.3711,-0.041,0.6214,0.2329
      ↪ ,0.0861,0.4875,-0.0477,0.0817,0.7994,0.3745,-0.0096,0.3395'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP44'

_symmetry_space_group_name_H-M "P 21/n 21/m 21/a"
_symmetry_Int_Tables_number 62

_cell_length_a      13.65500
_cell_length_b      8.13900
_cell_length_c      5.01600
_cell_angle_alpha   90.00000
_cell_angle_beta    90.00000
_cell_angle_gamma   90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x,y+1/2,-z
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x,-y+1/2,z
8 x+1/2,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 4 c 0.15240 0.25000 0.06130 1.00000
O2 O 4 c 0.07820 0.25000 0.55880 1.00000
O3 O 4 c 0.89690 0.25000 0.33860 1.00000
Si1 Si 4 c 0.18210 0.25000 0.37110 1.00000
Si2 Si 4 c -0.04100 0.25000 0.62140 1.00000
O4 O 8 d 0.23290 0.08610 0.48750 1.00000
O5 O 8 d -0.04770 0.08170 0.79940 1.00000
Y1 Y 8 d 0.37450 -0.00960 0.33950 1.00000
```

Possible δ-$Y_2Si_2O_7$: A7B2C2_oP44_62_3c2d_2c_d - POSCAR

```
A7B2C2_oP44_62_3c2d_2c_d & a,b/a,c/a,x1,z1,x2,z2,x3,z3,x4,z4,x5,z5,x6,y6
      ↪ ,z6,x7,y7,z7,x8,y8,z8 --params=13.655,0.596045404614,
      ↪ 0.367337971439,0.1524,0.0613,0.0782,0.5588,0.8969,0.3386,0.1821
      ↪ ,0.3711,-0.041,0.6214,0.2329,0.0861,0.4875,-0.0477,0.0817,
      ↪ 0.7994,0.3745,-0.0096,0.3395 & Pnma D_{2h}^{16} #62 (c^5d^3) &
      ↪ oP44 & None & O7Si2Y2 & O7Si2Y2 & H. W. Dias et al.,
      ↪ Zeitschrift f{"u}r Kristallographie - Crystalline Materials 191
      ↪ , 117-124 (1990)
   1.00000000000000
  13.65500000000000    0.00000000000000    0.00000000000000
   0.00000000000000    8.13900000000000    0.00000000000000
   0.00000000000000    0.00000000000000    5.01600000000000
     O     Si     Y
    28     8     8
Direct
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   0.34760000000000    0.75000000000000    0.56130000000000    O    (4c)
  -0.15240000000000    0.75000000000000   -0.06130000000000    O    (4c)
   0.65240000000000    0.25000000000000    0.43870000000000    O    (4c)
   0.07820000000000    0.25000000000000    0.55880000000000    O    (4c)
   0.42180000000000    0.75000000000000    1.05880000000000    O    (4c)
  -0.07820000000000    0.75000000000000   -0.55880000000000    O    (4c)
   0.57820000000000    0.25000000000000   -0.05880000000000    O    (4c)
   0.89690000000000    0.25000000000000    0.33860000000000    O    (4c)
  -0.39690000000000    0.75000000000000    0.83860000000000    O    (4c)
  -0.89690000000000    0.75000000000000   -0.33860000000000    O    (4c)
   1.39690000000000    0.25000000000000    0.16140000000000    O    (4c)
   0.23290000000000    0.08610000000000    0.48750000000000    O    (8d)
   0.26710000000000   -0.08610000000000    0.98750000000000    O    (8d)
  -0.23290000000000    0.58610000000000   -0.48750000000000    O    (8d)
   0.73290000000000    0.41390000000000    0.01250000000000    O    (8d)
  -0.23290000000000   -0.08610000000000   -0.48750000000000    O    (8d)
   0.73290000000000    0.08610000000000    0.01250000000000    O    (8d)
   0.23290000000000    0.41390000000000    0.48750000000000    O    (8d)
   0.26710000000000    0.58610000000000    0.98750000000000    O    (8d)
  -0.04770000000000    0.08170000000000    0.79940000000000    O    (8d)
   0.54770000000000   -0.08170000000000    1.29940000000000    O    (8d)
   0.04770000000000    0.58170000000000   -0.79940000000000    O    (8d)
   0.45230000000000    0.41830000000000   -0.29940000000000    O    (8d)
   0.04770000000000   -0.08170000000000   -0.79940000000000    O    (8d)
   0.45230000000000    0.08170000000000   -0.29940000000000    O    (8d)
  -0.04770000000000    0.41830000000000    0.79940000000000    O    (8d)
   0.54770000000000    0.58170000000000    1.29940000000000    O    (8d)
   0.18210000000000    0.25000000000000    0.37110000000000    Si   (4c)
   0.31790000000000    0.75000000000000    0.87110000000000    Si   (4c)
  -0.18210000000000    0.75000000000000   -0.37110000000000    Si   (4c)
   0.68210000000000    0.25000000000000    0.12890000000000    Si   (4c)
  -0.04100000000000    0.25000000000000    0.62140000000000    Si   (4c)
   0.54100000000000    0.75000000000000    1.12140000000000    Si   (4c)
   0.04100000000000    0.75000000000000   -0.62140000000000    Si   (4c)
   0.45900000000000    0.25000000000000   -0.12140000000000    Si   (4c)
   0.37450000000000   -0.00960000000000    0.33950000000000    Y    (8d)
   0.12550000000000    0.00960000000000    0.83950000000000    Y    (8d)
```

```
 −0.37450000000000   0.49040000000000  −0.33950000000000   Y   (8d)
  0.87450000000000   0.50960000000000   0.16050000000000   Y   (8d)
 −0.37450000000000   0.00960000000000  −0.33950000000000   Y   (8d)
  0.87450000000000  −0.00960000000000   0.16050000000000   Y   (8d)
  0.37450000000000   0.50960000000000   0.33950000000000   Y   (8d)
  0.12550000000000   0.49040000000000   0.83950000000000   Y   (8d)
```

$K_4[Mo(CN)_8]\cdot 2H_2O$ ($F2_1$): A8B4C4DE8F2_oP108_62_4c2d_2d_2cd_c_4c2d_d - CIF

```
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_audit_creation_method FINDSYM

_chemical_name_mineral 'C8H4K4MoN8O2'
_chemical_formula_sum 'C8 H4 K4 Mo N8 O2'

loop_
_publ_author_name
 'I. Typilo'
 'O. Sereda'
 'H. {Stoeckli-Evans}'
 'R. Gladyshevskii'
 'D. Semenyshyn'
_journal_name_full_name
;
 Chemistry of Metals and Alloys
;
_journal_volume 3
_journal_year 2010
_journal_page_first 49
_journal_page_last 52
_publ_Section_title
;
 Refinement of the crystal structure of potassium octacyanomolybdate(IV)
      ↪ dihydrate
;

_aflow_title 'K$_{4}$[Mo(CN)$_{8}$]$\cdot$2H$_{2}$O ($F2_{1}$) Structure
      ↪ '
_aflow_proto 'A8B4C4DE8F2_oP108_62_4c2d_2d_2cd_c_4c2d_d'
_aflow_params 'a,b/a,c/a,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3},x_{4},z_{4
      ↪ },x_{5},z_{5},x_{6},z_{6},x_{7},z_{7},x_{8},z_{8},x_{9},z_{9},
      ↪ x_{10},z_{10},x_{11},z_{11},x_{12},y_{12},z_{12},x_{13},y_{13},
      ↪ z_{13},x_{14},y_{14},z_{14},x_{15},y_{15},z_{15},x_{16},y_{16},
      ↪ z_{16},x_{17},y_{17},z_{17},x_{18},y_{18},z_{18},x_{19},y_{19},
      ↪ z_{19}'
_aflow_params_values '16.6959,0.695320408004,0.519864158266,0.0619,
      ↪ 0.3937,0.1418,0.8446,0.2168,0.3978,0.2602,0.6753,0.36535,0.025,
      ↪ 0.46252,0.52249,0.1375,0.59643,0.021,0.2866,0.2614,0.2957,0.326
      ↪ ,0.7133,0.6401,0.5221,0.0391,0.1388,0.6648,0.1731,0.0725,0.5689
      ↪ ,0.349,0.08,0.367,0.412,0.017,0.347,0.14835,0.04756,0.18788,
      ↪ 0.0123,0.5809,0.2893,0.3068,0.0216,0.05,0.3888,0.0593,0.409'
_aflow_Strukturbericht '$F2_{1}$'
_aflow_Pearson 'oP108'

_symmetry_space_group_name_H-M "P 21/n 21/m 21/a"
_symmetry_Int_Tables_number 62

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_cell_length_b     11.60900
_cell_length_c     8.67960
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
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2 x+1/2,-y+1/2,-z+1/2
3 -x,y+1/2,-z
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x,-y+1/2,z
8 x+1/2,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
C1 C 4 c 0.06190 0.25000 0.39370 1.00000
C2 C 4 c 0.14180 0.25000 0.84460 1.00000
C3 C 4 c 0.21680 0.25000 0.39780 1.00000
C4 C 4 c 0.26020 0.25000 0.67530 1.00000
K1 K 4 c 0.36535 0.25000 0.02500 1.00000
K2 K 4 c 0.46252 0.25000 0.52249 1.00000
Mo1 Mo 4 c 0.13750 0.25000 0.59643 1.00000
N1 N 4 c 0.02100 0.25000 0.28660 1.00000
N2 N 4 c 0.26140 0.25000 0.29570 1.00000
N3 N 4 c 0.32600 0.25000 0.71330 1.00000
N4 N 4 c 0.64010 0.25000 0.52210 1.00000
C5 C 8 d 0.03910 0.13880 0.66480 1.00000
C6 C 8 d 0.17310 0.07250 0.56890 1.00000
H1 H 8 d 0.34900 0.08000 0.36700 1.00000
H2 H 8 d 0.41200 0.01700 0.34700 1.00000
K3 K 8 d 0.14835 0.04756 0.18788 1.00000
N5 N 8 d 0.01230 0.58090 0.28930 1.00000
N6 N 8 d 0.30680 0.02160 0.05000 1.00000
O1 O 8 d 0.38880 0.05930 0.40900 1.00000
```

$K_4[Mo(CN)_8]\cdot 2H_2O$ ($F2_1$): A8B4C4DE8F2_oP108_62_4c2d_2d_2cd_c_4c2d_d - POSCAR

```
A8B4C4DE8F2_oP108_62_4c2d_2d_2cd_c_4c2d_d & a,b/a,c/a,x1,z1,x2,z2,x3,z3,
      ↪ x4,z4,x5,z5,x6,z6,x7,z7,x8,z8,x9,z9,x10,z10,x11,z11,x12,y12,z12
      ↪ ,x13,y13,z13,x14,y14,z14,x15,y15,z15,x16,y16,z16,x17,y17,z17,
      ↪ x18,y18,z18,x19,y19,z19 --params=16.6959,0.695320408004,
      ↪ 0.519864158266,0.0619,0.3937,0.1418,0.8446,0.2168,0.3978,0.2602
      ↪ ,0.6753,0.36535,0.025,0.46252,0.52249,0.1375,0.59643,0.021,
      ↪ 0.2866,0.2614,0.2957,0.326,0.7133,0.6401,0.5221,0.0391,0.1388,
      ↪ 0.6648,0.1731,0.0725,0.5689,0.349,0.08,0.367,0.412,0.017,0.347,
      ↪ 0.14835,0.04756,0.18788,0.0123,0.5809,0.2893,0.3068,0.0216,0.05
      ↪ ,0.3888,0.0593,0.409 & Pnma D_{2h}^{16} #62 (c^11d^8) & oP108 &
      ↪  $F2_{1}$ & C8H4K4MoN8O2 & C8H4K4MoN8O2 & I. Typilo et al.,
      ↪ Chem. Met. Alloys 3, 49-52 (2010)
  1.00000000000000
 16.69590000000000   0.00000000000000   0.00000000000000
  0.00000000000000  11.60900000000000   0.00000000000000
  0.00000000000000   0.00000000000000   8.67960000000000
     C     H     K    Mo     N     O
    32    16    16     4    32     8
Direct
  0.06190000000000   0.25000000000000   0.39370000000000   C   (4c)
  0.43810000000000   0.75000000000000   0.89370000000000   C   (4c)
 -0.06190000000000   0.75000000000000  -0.39370000000000   C   (4c)
  0.56190000000000   0.25000000000000   0.10630000000000   C   (4c)
  0.14180000000000   0.25000000000000   0.84460000000000   C   (4c)
  0.35820000000000   0.75000000000000   1.34460000000000   C   (4c)
 -0.14180000000000   0.75000000000000  -0.84460000000000   C   (4c)
  0.64180000000000   0.25000000000000  -0.34460000000000   C   (4c)
  0.21680000000000   0.25000000000000   0.39780000000000   C   (4c)
  0.28320000000000   0.75000000000000   0.89780000000000   C   (4c)
 -0.21680000000000   0.75000000000000  -0.39780000000000   C   (4c)
  0.71680000000000   0.25000000000000   0.10220000000000   C   (4c)
  0.26020000000000   0.25000000000000   0.67530000000000   C   (4c)
  0.23980000000000   0.75000000000000   1.17530000000000   C   (4c)
 -0.26020000000000   0.75000000000000  -0.67530000000000   C   (4c)
  0.76020000000000   0.25000000000000  -0.17530000000000   C   (4c)
  0.03910000000000   0.13880000000000   0.66480000000000   C   (8d)
  0.46090000000000  -0.13880000000000   1.16480000000000   C   (8d)
 -0.03910000000000   0.63880000000000  -0.66480000000000   C   (8d)
  0.53910000000000   0.36120000000000  -0.16480000000000   C   (8d)
 -0.03910000000000  -0.13880000000000  -0.66480000000000   C   (8d)
  0.53910000000000   0.13880000000000  -0.16480000000000   C   (8d)
  0.03910000000000   0.36120000000000   0.66480000000000   C   (8d)
  0.46090000000000   0.63880000000000   1.16480000000000   C   (8d)
  0.17310000000000   0.07250000000000   0.56890000000000   C   (8d)
  0.32690000000000  -0.07250000000000   1.06890000000000   C   (8d)
 -0.17310000000000   0.57250000000000  -0.56890000000000   C   (8d)
  0.67310000000000   0.42750000000000  -0.06890000000000   C   (8d)
 -0.17310000000000  -0.07250000000000  -0.56890000000000   C   (8d)
  0.67310000000000   0.07250000000000  -0.06890000000000   C   (8d)
  0.17310000000000   0.42750000000000   0.56890000000000   C   (8d)
  0.32690000000000   0.57250000000000   1.06890000000000   C   (8d)
  0.34900000000000   0.08000000000000   0.36700000000000   H   (8d)
  0.15100000000000  -0.08000000000000   0.86700000000000   H   (8d)
 -0.34900000000000   0.58000000000000  -0.36700000000000   H   (8d)
  0.84900000000000   0.42000000000000   0.13300000000000   H   (8d)
 -0.34900000000000  -0.08000000000000  -0.36700000000000   H   (8d)
  0.84900000000000   0.08000000000000   0.13300000000000   H   (8d)
  0.34900000000000   0.42000000000000   0.36700000000000   H   (8d)
  0.15100000000000   0.58000000000000   0.86700000000000   H   (8d)
  0.41200000000000   0.01700000000000   0.34700000000000   H   (8d)
  0.08800000000000  -0.01700000000000   0.84700000000000   H   (8d)
 -0.41200000000000   0.51700000000000  -0.34700000000000   H   (8d)
  0.91200000000000   0.48300000000000   0.15300000000000   H   (8d)
 -0.41200000000000  -0.01700000000000  -0.34700000000000   H   (8d)
  0.91200000000000   0.01700000000000   0.15300000000000   H   (8d)
  0.41200000000000   0.48300000000000   0.34700000000000   H   (8d)
  0.08800000000000   0.51700000000000   0.84700000000000   H   (8d)
  0.36535000000000   0.25000000000000   0.02500000000000   K   (4c)
  0.13465000000000   0.75000000000000   0.52500000000000   K   (4c)
 -0.36535000000000   0.75000000000000  -0.02500000000000   K   (4c)
  0.86535000000000   0.25000000000000   0.47500000000000   K   (4c)
  0.46252000000000   0.25000000000000   0.52249000000000   K   (4c)
  0.03748000000000   0.75000000000000   1.02249000000000   K   (4c)
 -0.46252000000000   0.75000000000000  -0.52249000000000   K   (4c)
  0.96252000000000   0.25000000000000  -0.02249000000000   K   (4c)
  0.14835000000000   0.04756000000000   0.18788000000000   K   (8d)
  0.35165000000000  -0.04756000000000   0.68788000000000   K   (8d)
 -0.14835000000000   0.54756000000000  -0.18788000000000   K   (8d)
  0.64835000000000   0.45244000000000   0.31212000000000   K   (8d)
 -0.14835000000000  -0.04756000000000  -0.18788000000000   K   (8d)
  0.64835000000000   0.04756000000000   0.31212000000000   K   (8d)
  0.14835000000000   0.45244000000000   0.18788000000000   K   (8d)
  0.35165000000000   0.54756000000000   0.68788000000000   K   (8d)
  0.13750000000000   0.25000000000000   0.59643000000000   Mo  (4c)
  0.36250000000000   0.75000000000000   1.09643000000000   Mo  (4c)
 -0.13750000000000   0.75000000000000  -0.59643000000000   Mo  (4c)
  0.63750000000000   0.25000000000000  -0.09643000000000   Mo  (4c)
  0.02100000000000   0.25000000000000   0.28660000000000   N   (4c)
  0.47900000000000   0.75000000000000   0.78660000000000   N   (4c)
 -0.02100000000000   0.75000000000000  -0.28660000000000   N   (4c)
  0.52100000000000   0.25000000000000   0.21340000000000   N   (4c)
  0.26140000000000   0.25000000000000   0.29570000000000   N   (4c)
  0.23860000000000   0.75000000000000   0.79570000000000   N   (4c)
 -0.26140000000000   0.75000000000000  -0.29570000000000   N   (4c)
  0.76140000000000   0.25000000000000   0.20430000000000   N   (4c)
  0.32600000000000   0.25000000000000   0.71330000000000   N   (4c)
  0.17400000000000   0.75000000000000   1.21330000000000   N   (4c)
 -0.32600000000000   0.75000000000000  -0.71330000000000   N   (4c)
  0.82600000000000   0.25000000000000  -0.21330000000000   N   (4c)
  0.64010000000000   0.25000000000000   0.52210000000000   N   (4c)
 -0.14010000000000   0.75000000000000   1.02210000000000   N   (4c)
 -0.64010000000000   0.75000000000000  -0.52210000000000   N   (4c)
  1.14010000000000   0.25000000000000  -0.02210000000000   N   (4c)
  0.01230000000000   0.58090000000000   0.28930000000000   N   (8d)
```

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   0.48770000000000  −0.58090000000000   0.78930000000000   N   (8d)
  −0.01230000000000   1.08090000000000  −0.28930000000000   N   (8d)
   0.51230000000000  −0.08090000000000   0.21070000000000   N   (8d)
  −0.01230000000000  −0.58090000000000  −0.28930000000000   N   (8d)
   0.51230000000000   0.58090000000000   0.21070000000000   N   (8d)
   0.01230000000000  −0.08090000000000   0.28930000000000   N   (8d)
   0.48770000000000   1.08090000000000   0.78930000000000   N   (8d)
   0.30680000000000   0.02160000000000   0.05000000000000   N   (8d)
   0.19320000000000  −0.02160000000000   0.55000000000000   N   (8d)
  −0.30680000000000   0.52160000000000  −0.05000000000000   N   (8d)
   0.80680000000000   0.47840000000000   0.45000000000000   N   (8d)
  −0.30680000000000  −0.02160000000000  −0.05000000000000   N   (8d)
   0.80680000000000   0.02160000000000   0.45000000000000   N   (8d)
   0.30680000000000   0.47840000000000   0.05000000000000   N   (8d)
   0.19320000000000   0.52160000000000   0.55000000000000   N   (8d)
   0.38880000000000   0.05930000000000   0.40900000000000   O   (8d)
   0.11120000000000  −0.05930000000000   0.90900000000000   O   (8d)
  −0.38880000000000   0.55930000000000  −0.40900000000000   O   (8d)
   0.88880000000000   0.44070000000000   0.09100000000000   O   (8d)
  −0.38880000000000  −0.05930000000000  −0.40900000000000   O   (8d)
   0.88880000000000   0.05930000000000   0.09100000000000   O   (8d)
   0.38880000000000   0.44070000000000   0.40900000000000   O   (8d)
   0.11120000000000   0.55930000000000   0.90900000000000   O   (8d)
```

$SbCl_5{\cdot}POCl_3$: A8BCD_oP44_62_4c2d_c_c_c - CIF

```
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_audit_creation_method FINDSYM

_chemical_name_mineral 'Cl8OPSb'
_chemical_formula_sum 'Cl8 O P Sb'

loop_
_publ_author_name
 'I. Lindqvist'
 'C.-I. Br\"{a}nd\\'{e}n'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 12
_journal_year 1959
_journal_page_first 642
_journal_page_last 645
_publ_Section_title
;
 The Crystal Structure of SbCl$_{5}$$\cdot$POCl$_{3}$
;

# Found in The Crystal Structure of SbCl$_{5}$.PO(CH$_{3}$)$_{3}$, 1961

_aflow_title 'SbCl$_{5}$$\cdot$POCl$_{3}$ Structure'
_aflow_proto 'A8BCD_oP44_62_4c2d_c_c_c'
_aflow_params 'a,b/a,c/a,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3},x_{4},z_{4
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      ↪ y_{9},z_{9}'
_aflow_params_values '16.42,0.490864799026,0.543848964677,0.2585,-0.0805
      ↪ ,0.0202,0.2048,0.2249,0.2938,0.4644,0.8699,0.0706,0.8774,0.0742
      ↪ ,0.7142,0.145,0.0797,0.1314,0.4442,0.6327,0.1405,0.5383,0.0597'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP44'

_symmetry_space_group_name_H-M "P 21/n 21/m 21/a"
_symmetry_Int_Tables_number 62

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_cell_length_b     8.06000
_cell_length_c     8.93000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

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_space_group_symop_operation_xyz
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2 x+1/2,-y+1/2,-z+1/2
3 -x,y+1/2,-z
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x,-y+1/2,z
8 x+1/2,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cl1  Cl  4  c  0.25850  0.25000  -0.08050  1.00000
Cl2  Cl  4  c  0.02020  0.25000  0.20480  1.00000
Cl3  Cl  4  c  0.22490  0.25000  0.29380  1.00000
Cl4  Cl  4  c  0.46440  0.25000  0.86990  1.00000
O1  O  4  c  0.07060  0.25000  0.87740  1.00000
P1  P  4  c  0.07420  0.25000  0.71420  1.00000
Sb1  Sb  4  c  0.14500  0.25000  0.07970  1.00000
Cl5  Cl  8  d  0.13140  0.44420  0.63270  1.00000
Cl6  Cl  8  d  0.14050  0.53830  0.05970  1.00000
```

$SbCl_5{\cdot}POCl_3$: A8BCD_oP44_62_4c2d_c_c_c - POSCAR

```
A8BCD_oP44_62_4c2d_c_c_c & a,b/a,c/a,x1,z1,x2,z2,x3,z3,x4,z4,x5,z5,x6,z6
      ↪ ,x7,z7,x8,y8,z8,x9,y9,z9 --params=16.42,0.490864799026,
      ↪ 0.543848964677,0.2585,-0.0805,0.0202,0.2048,0.2249,0.2938,
      ↪ 0.4644,0.8699,0.0706,0.8774,0.0742,0.7142,0.145,0.0797,0.1314,
      ↪ 0.4442,0.6327,0.1405,0.5383,0.0597 & Pnma D_{2h}^{16} #62 (c^7d
      ↪ ^2) & oP44 & None & Cl8OPSb & Cl8OPSb & I. Lindqvist and C.-I.
      ↪ Br\"{a}nd\'{e}n, Acta Cryst. 12, 642-645 (1959)
   1.00000000000000
  16.42000000000000   0.00000000000000   0.00000000000000
   0.00000000000000   8.06000000000000   0.00000000000000
   0.00000000000000   0.00000000000000   8.93000000000000
     Cl       O       P      Sb
     32       4       4       4
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   0.24150000000000   0.75000000000000   0.41950000000000   Cl   (4c)
  -0.25850000000000   0.75000000000000   0.08050000000000   Cl   (4c)
   0.75850000000000   0.25000000000000   0.58050000000000   Cl   (4c)
   0.02020000000000   0.25000000000000   0.20480000000000   Cl   (4c)
   0.47980000000000   0.75000000000000   0.70480000000000   Cl   (4c)
  -0.02020000000000   0.75000000000000  -0.20480000000000   Cl   (4c)
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   0.22490000000000   0.25000000000000   0.29380000000000   Cl   (4c)
   0.27510000000000   0.75000000000000   0.79380000000000   Cl   (4c)
  -0.22490000000000   0.75000000000000  -0.29380000000000   Cl   (4c)
   0.72490000000000   0.25000000000000   0.20620000000000   Cl   (4c)
   0.46440000000000   0.25000000000000   0.86990000000000   Cl   (4c)
   0.03560000000000   0.75000000000000   1.36990000000000   Cl   (4c)
  -0.46440000000000   0.75000000000000  -0.86990000000000   Cl   (4c)
   0.96440000000000   0.25000000000000  -0.36990000000000   Cl   (4c)
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   0.36860000000000  -0.44420000000000   1.13270000000000   Cl   (8d)
  -0.13140000000000   0.94420000000000  -0.63270000000000   Cl   (8d)
   0.63140000000000   0.05580000000000  -0.13270000000000   Cl   (8d)
  -0.13140000000000  -0.44420000000000  -0.63270000000000   Cl   (8d)
   0.63140000000000   0.44420000000000  -0.13270000000000   Cl   (8d)
   0.13140000000000   0.05580000000000   0.63270000000000   Cl   (8d)
   0.36860000000000   0.94420000000000   1.13270000000000   Cl   (8d)
   0.14050000000000   0.53830000000000   0.05970000000000   Cl   (8d)
   0.35950000000000  -0.53830000000000   0.55970000000000   Cl   (8d)
  -0.14050000000000   1.03830000000000  -0.05970000000000   Cl   (8d)
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   0.64050000000000   0.53830000000000   0.44030000000000   Cl   (8d)
   0.14050000000000  -0.03830000000000   0.05970000000000   Cl   (8d)
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   0.07060000000000   0.25000000000000   0.87740000000000    O   (4c)
   0.42940000000000   0.75000000000000   1.37740000000000    O   (4c)
  -0.07060000000000   0.75000000000000  -0.87740000000000    O   (4c)
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   0.07420000000000   0.25000000000000   0.71420000000000    P   (4c)
   0.42580000000000   0.75000000000000   1.21420000000000    P   (4c)
  -0.07420000000000   0.75000000000000  -0.71420000000000    P   (4c)
   0.57420000000000   0.25000000000000  -0.21420000000000    P   (4c)
   0.14500000000000   0.25000000000000   0.07970000000000   Sb   (4c)
   0.35500000000000   0.75000000000000   0.57970000000000   Sb   (4c)
  -0.14500000000000   0.75000000000000  -0.07970000000000   Sb   (4c)
   0.64500000000000   0.25000000000000   0.42030000000000   Sb   (4c)
```

Autunite $\{Ca[(UO_2)(PO_4)]_2(H_2O)_{11}\}$: AB22C23D2E2_oP200_62_c_11d_3c10d_d_d - CIF

```
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_audit_creation_method FINDSYM

_chemical_name_mineral 'Autunite'
_chemical_formula_sum 'Ca H22 O23 P2 U2'

loop_
_publ_author_name
 'A. J. Locock'
 'P. C. Burns'
_journal_name_full_name
;
 American Mineralogist
;
_journal_volume 88
_journal_year 2003
_journal_page_first 240
_journal_page_last 244
_publ_Section_title
;
 The crystal structure of synthetic autunite, Ca[(UO$_{2}$)(PO$_{4}$)]$_
      ↪ {2}$(H$_{2}$O)$_{11}$
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'Autunite \{Ca[(UO$_{2}$)(PO$_{4}$)]$_{2}$(H$_{2}$O)$_{11}$
      ↪ \} Structure'
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      ↪ y_{8},z_{8},x_{9},y_{9},z_{9},x_{10},y_{10},z_{10},x_{11},y_{11
      ↪ },z_{11},x_{12},y_{12},z_{12},x_{13},y_{13},z_{13},x_{14},y_{14
      ↪ },z_{14},x_{15},y_{15},z_{15},x_{16},y_{16},z_{16},x_{17},y_{17
      ↪ },z_{17},x_{18},y_{18},z_{18},x_{19},y_{19},z_{19},x_{20},y_{20
      ↪ },z_{20},x_{21},y_{21},z_{21},x_{22},y_{22},z_{22},x_{23},y_{23
      ↪ },z_{23},x_{24},y_{24},z_{24},x_{25},y_{25},z_{25},x_{26},y_{26
      ↪ },z_{26},x_{27},y_{27},z_{27}'
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      ↪ 0.722,0.697,0.133,-0.088,0.721,0.154,0.715,0.533,0.1191,0.835,
      ↪ 0.508,0.165,0.701,0.537,0.7851,0.733,0.63,0.2056,0.544,0.772,
      ↪ 0.174,0.558,0.8972,0.155,0.473,0.568,0.8819,0.575,0.6483,0.836,
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      ↪ 0.522,0.6235,-0.0449,0.2482,0.625,0.1278,0.2472,0.6252,0.044,
      ↪ 0.5733,0.5373,-0.044,0.7451,0.7118,-0.0443,0.7553,0.6206,0.0452
      ↪ ,-0.0774,0.7368,0.1647,0.851,0.5126,0.1645,0.8438,0.8257,0.1614
      ↪ ,0.4745,0.5759,0.8379,0.5243,0.6245,0.0003,0.7478,0.625,0.0412,
      ↪ 0.2498'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP200'

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_symmetry_Int_Tables_number 62

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_cell_length_c 6.99590
_cell_angle_alpha 90.00000
_cell_angle_beta 90.00000
_cell_angle_gamma 90.00000

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_space_group_symop_operation_xyz
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3 -x,y+1/2,-z
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x,-y+1/2,z
8 x+1/2,y,-z+1/2

loop_
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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
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O1 O 4 c 0.74090 0.25000 0.21490 1.00000
O2 O 4 c 0.49980 0.25000 0.21670 1.00000
O3 O 4 c 0.62630 0.25000 0.57570 1.00000
H1 H 8 d 0.22600 0.78300 0.72200 1.00000
H2 H 8 d 0.69700 0.13300 -0.08800 1.00000
H3 H 8 d 0.72100 0.15400 0.71500 1.00000
H4 H 8 d 0.53300 0.11910 0.83500 1.00000
H5 H 8 d 0.50800 0.16500 0.70100 1.00000
H6 H 8 d 0.53700 0.78510 0.73300 1.00000
H7 H 8 d 0.63000 0.20560 0.54400 1.00000
H8 H 8 d 0.77200 0.17400 0.55800 1.00000
H9 H 8 d 0.89720 0.15500 0.47300 1.00000
H10 H 8 d 0.56800 0.88190 0.57500 1.00000
H11 H 8 d 0.64830 0.83600 0.52200 1.00000
O4 O 8 d 0.62350 -0.04490 0.24820 1.00000
O5 O 8 d 0.62500 0.12780 0.24720 1.00000
O6 O 8 d 0.62520 0.04400 0.57330 1.00000
O7 O 8 d 0.53730 -0.04400 0.74510 1.00000
O8 O 8 d 0.71180 -0.04430 0.75530 1.00000
O9 O 8 d 0.62060 0.04520 -0.07740 1.00000
O10 O 8 d 0.73680 0.16470 0.85100 1.00000
O11 O 8 d 0.51260 0.16450 0.84380 1.00000
O12 O 8 d 0.82570 0.16140 0.47450 1.00000
O13 O 8 d 0.57590 0.83790 0.52430 1.00000
P1 P 8 d 0.62450 0.00030 0.74780 1.00000
U1 U 8 d 0.62500 0.04120 0.24980 1.00000
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Autunite $\{Ca[(UO_2)(PO_4)]_2(H_2O)_{11}\}$: AB22C23D2E2_oP200_62_c_11d_3c10d_d_d - POSCAR

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AB22C23D2E2_oP200_62_c_11d_3c10d_d_d & a,b/a,c/a,x1,z1,x2,z2,x3,z3,x4,z4
      ↪ ,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10,x11,
      ↪ y11,z11,x12,y12,z12,x13,y13,z13,x14,y14,z14,x15,y15,z15,x16,y16
      ↪ ,z16,x17,y17,z17,x18,y18,z18,x19,y19,z19,x20,y20,z20,x21,y21,
      ↪ z21,x22,y22,z22,x23,y23,z23,x24,y24,z24,x25,y25,z25,x26,y26,z26
      ↪ ,x27,y27,z27 --params=14.0135,1.47800335391,0.499225746601,
      ↪ 0.6251,-0.0499,0.7409,0.2149,0.4998,0.2167,0.6263,0.5757,0.226,
      ↪ 0.783,0.722,0.697,0.133,-0.088,0.721,0.154,0.715,0.533,0.1191,
      ↪ 0.835,0.508,0.165,0.701,0.537,0.7851,0.733,0.63,0.2056,0.544,
      ↪ 0.772,0.174,0.558,0.8972,0.155,0.473,0.568,0.8819,0.575,0.6483,
      ↪ 0.836,0.522,0.6235,-0.0449,0.2482,0.625,0.1278,0.2472,0.6252,
      ↪ 0.044,0.5733,0.5373,-0.044,0.7451,0.7118,-0.0443,0.7553,0.6206,
      ↪ 0.0452,-0.0774,0.7368,0.1647,0.851,0.5126,0.1645,0.8438,0.8257,
      ↪ 0.1614,0.4745,0.5759,0.8379,0.5243,0.6245,0.0003,0.7478,0.625,
      ↪ 0.0412,0.2498 & Pnma D_{2h}^{16} #62 (c^4d^23) & oP200 & None &
      ↪  CaH22O23P2U2 & Autunite & A. J. Locock and P. C. Burns, Am.
      ↪ Mineral. 88, 240-244 (2003)
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  14.01350000000000    0.00000000000000    0.00000000000000
   0.00000000000000   20.71200000000000    0.00000000000000
   0.00000000000000    0.00000000000000    6.99590000000000
    Ca    H    O    P    U
     4   88   92    8    8
Direct
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  -0.12510000000000    0.75000000000000    0.45010000000000   Ca   (4c)
  -0.62510000000000    0.75000000000000    0.04990000000000   Ca   (4c)
   1.12510000000000    0.25000000000000    0.54990000000000   Ca   (4c)
   0.22600000000000    0.78300000000000    0.72200000000000   H    (8d)
   0.27400000000000   -0.78300000000000    1.22200000000000   H    (8d)
  -0.22600000000000    1.28300000000000   -0.72200000000000   H    (8d)
   0.72600000000000   -0.28300000000000   -0.22200000000000   H    (8d)
  -0.22600000000000   -0.78300000000000   -0.72200000000000   H    (8d)
   0.72600000000000    0.78300000000000   -0.22200000000000   H    (8d)
   0.22600000000000   -0.28300000000000    0.72200000000000   H    (8d)
   0.27400000000000    1.28300000000000    1.22200000000000   H    (8d)
   0.69700000000000    0.13300000000000   -0.08800000000000   H    (8d)
  -0.19700000000000   -0.13300000000000    0.41200000000000   H    (8d)
  -0.69700000000000    0.63300000000000    0.08800000000000   H    (8d)
   1.19700000000000    0.36700000000000    0.58800000000000   H    (8d)
  -0.69700000000000   -0.13300000000000    0.08800000000000   H    (8d)
   1.19700000000000    0.13300000000000    0.58800000000000   H    (8d)
   0.69700000000000    0.36700000000000   -0.08800000000000   H    (8d)
  -0.19700000000000    0.63300000000000    0.41200000000000   H    (8d)
   0.72100000000000    0.15400000000000    0.71500000000000   H    (8d)
  -0.22100000000000   -0.15400000000000    1.21500000000000   H    (8d)
  -0.72100000000000    0.65400000000000   -0.71500000000000   H    (8d)
   1.22100000000000    0.34600000000000   -0.21500000000000   H    (8d)
  -0.72100000000000   -0.15400000000000   -0.71500000000000   H    (8d)
   1.22100000000000    0.15400000000000   -0.21500000000000   H    (8d)
   0.72100000000000    0.34600000000000    0.71500000000000   H    (8d)
  -0.22100000000000    0.65400000000000    1.21500000000000   H    (8d)
   0.53300000000000    0.11910000000000    0.83500000000000   H    (8d)
  -0.03300000000000   -0.11910000000000    1.33500000000000   H    (8d)
  -0.53300000000000    0.61910000000000   -0.83500000000000   H    (8d)
   1.03300000000000    0.38090000000000   -0.33500000000000   H    (8d)
  -0.53300000000000   -0.11910000000000   -0.83500000000000   H    (8d)
   1.03300000000000    0.11910000000000   -0.33500000000000   H    (8d)
   0.53300000000000    0.38090000000000    0.83500000000000   H    (8d)
  -0.03300000000000    0.61910000000000    1.33500000000000   H    (8d)
   0.50800000000000    0.16500000000000    0.70100000000000   H    (8d)
  -0.00800000000000   -0.16500000000000    1.20100000000000   H    (8d)
  -0.50800000000000    0.66500000000000   -0.70100000000000   H    (8d)
   1.00800000000000    0.33500000000000   -0.20100000000000   H    (8d)
  -0.50800000000000   -0.16500000000000   -0.70100000000000   H    (8d)
   1.00800000000000    0.16500000000000   -0.20100000000000   H    (8d)
   0.50800000000000    0.33500000000000    0.70100000000000   H    (8d)
  -0.00800000000000    0.66500000000000    1.20100000000000   H    (8d)
   0.53700000000000    0.78510000000000    0.73300000000000   H    (8d)
  -0.03700000000000   -0.78510000000000    1.23300000000000   H    (8d)
  -0.53700000000000    1.28510000000000   -0.73300000000000   H    (8d)
   1.03700000000000   -0.28510000000000   -0.23300000000000   H    (8d)
  -0.53700000000000   -0.78510000000000   -0.73300000000000   H    (8d)
   1.03700000000000    0.78510000000000   -0.23300000000000   H    (8d)
   0.53700000000000   -0.28510000000000    0.73300000000000   H    (8d)
  -0.03700000000000    1.28510000000000    1.23300000000000   H    (8d)
   0.63000000000000    0.20560000000000    0.54400000000000   H    (8d)
  -0.13000000000000   -0.20560000000000    1.04400000000000   H    (8d)
  -0.63000000000000    0.70560000000000   -0.54400000000000   H    (8d)
   1.13000000000000    0.29440000000000   -0.04400000000000   H    (8d)
  -0.63000000000000   -0.20560000000000   -0.54400000000000   H    (8d)
   1.13000000000000    0.20560000000000   -0.04400000000000   H    (8d)
   0.63000000000000    0.29440000000000    0.54400000000000   H    (8d)
  -0.13000000000000    0.70560000000000    1.04400000000000   H    (8d)
   0.77200000000000    0.17400000000000    0.55800000000000   H    (8d)
  -0.27200000000000   -0.17400000000000    1.05800000000000   H    (8d)
  -0.77200000000000    0.67400000000000   -0.55800000000000   H    (8d)
   1.27200000000000    0.32600000000000   -0.05800000000000   H    (8d)
  -0.77200000000000   -0.17400000000000   -0.55800000000000   H    (8d)
   1.27200000000000    0.17400000000000   -0.05800000000000   H    (8d)
   0.77200000000000    0.32600000000000    0.55800000000000   H    (8d)
  -0.27200000000000    0.67400000000000    1.05800000000000   H    (8d)
   0.89720000000000    0.15500000000000    0.47300000000000   H    (8d)
  -0.39720000000000   -0.15500000000000    0.97300000000000   H    (8d)
  -0.89720000000000    0.65500000000000   -0.47300000000000   H    (8d)
   1.39720000000000    0.34500000000000    0.02700000000000   H    (8d)
  -0.89720000000000   -0.15500000000000   -0.47300000000000   H    (8d)
   1.39720000000000    0.15500000000000    0.02700000000000   H    (8d)
   0.89720000000000    0.34500000000000    0.47300000000000   H    (8d)
  -0.39720000000000    0.65500000000000    0.97300000000000   H    (8d)
   0.56800000000000    0.88190000000000    0.57500000000000   H    (8d)
  -0.06800000000000   -0.88190000000000    1.07500000000000   H    (8d)
  -0.56800000000000    1.38190000000000   -0.57500000000000   H    (8d)
   1.06800000000000   -0.38190000000000   -0.07500000000000   H    (8d)
  -0.56800000000000   -0.88190000000000   -0.57500000000000   H    (8d)
   1.06800000000000    0.88190000000000   -0.07500000000000   H    (8d)
   0.56800000000000   -0.38190000000000    0.57500000000000   H    (8d)
  -0.06800000000000    1.38190000000000    1.07500000000000   H    (8d)
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  -0.14830000000000   -0.83600000000000    1.02200000000000   H    (8d)
  -0.64830000000000    1.33600000000000   -0.52200000000000   H    (8d)
   1.14830000000000   -0.33600000000000   -0.02200000000000   H    (8d)
  -0.64830000000000   -0.83600000000000   -0.52200000000000   H    (8d)
   1.14830000000000    0.83600000000000   -0.02200000000000   H    (8d)
   0.64830000000000   -0.33600000000000    0.52200000000000   H    (8d)
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  -0.74090000000000    0.75000000000000   -0.21490000000000   O    (4c)
   1.24090000000000    0.25000000000000    0.28510000000000   O    (4c)
   0.49980000000000    0.25000000000000    0.21670000000000   O    (4c)
   0.00020000000000    0.75000000000000    0.71670000000000   O    (4c)
  -0.49980000000000    0.75000000000000   -0.21670000000000   O    (4c)
   0.99980000000000    0.25000000000000    0.28330000000000   O    (4c)
   0.62630000000000    0.25000000000000    0.57570000000000   O    (4c)
  -0.12630000000000    0.75000000000000    1.07570000000000   O    (4c)
  -0.62630000000000    0.75000000000000   -0.57570000000000   O    (4c)
   1.12630000000000    0.25000000000000   -0.07570000000000   O    (4c)
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   0.62500000000000    0.12780000000000    0.24720000000000   O    (8d)
  -0.12500000000000   -0.12780000000000    0.74720000000000   O    (8d)
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  -0.62500000000000   -0.12780000000000   -0.24720000000000   O    (8d)
```

```
  1.12500000000000   0.12780000000000   0.25280000000000   O   (8d)
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  0.62520000000000   0.04400000000000   0.57330000000000   O   (8d)
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  0.73680000000000   0.33530000000000   0.85100000000000   O   (8d)
 -0.23680000000000   0.66470000000000   1.35100000000000   O   (8d)
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  0.57590000000000  -0.33790000000000   0.52430000000000   O   (8d)
 -0.07590000000000   1.33790000000000   1.02430000000000   O   (8d)
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  0.62450000000000   0.49970000000000   0.74780000000000   P   (8d)
 -0.12450000000000   0.50030000000000   1.24780000000000   P   (8d)
  0.62500000000000   0.04120000000000   0.24980000000000   U   (8d)
 -0.12500000000000  -0.04120000000000   0.74980000000000   U   (8d)
 -0.62500000000000   0.54120000000000  -0.24980000000000   U   (8d)
  1.12500000000000   0.45880000000000   0.25020000000000   U   (8d)
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  0.62500000000000   0.45880000000000   0.24980000000000   U   (8d)
 -0.12500000000000   0.54120000000000   0.74980000000000   U   (8d)
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Atacamite ($Cu_2(OH)_3Cl$): AB2C3D3_oP36_62_c_ac_cd_cd - CIF

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_audit_creation_method FINDSYM

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loop_
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 'J. B. Parise'
 'B. G. Hyde'
_journal_name_full_name
;
 Acta Crystallographica Section C: Structural Chemistry
;
_journal_volume 42
_journal_year 1986
_journal_page_first 1277
_journal_page_last 1280
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;
 The structure of atacamite and its relationship to spinel
;

# Found in The American Mineralogist Crystal Structure Database, 2003

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_aflow_proto 'AB2C3D3_oP36_62_c_ac_cd_cd'
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_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP36'

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_symmetry_Int_Tables_number 62

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_cell_angle_gamma 90.00000

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_space_group_symop_operation_xyz
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3 -x,y+1/2,-z
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x,-y+1/2,z
8 x+1/2,y,-z+1/2

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_atom_site_type_symbol
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_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
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Cl1 Cl 4 c 0.85180 0.25000 0.55560 1.00000
Cu2 Cu 4 c 0.30940 0.25000 0.24470 1.00000
H1 H 4 c 0.19510 0.25000 0.51480 1.00000
O1 O 4 c 0.35020 0.25000 0.50180 1.00000
H2 H 8 d -0.06690 0.46660 0.72790 1.00000
O2 O 8 d -0.05940 0.56510 0.78790 1.00000
```

Atacamite ($Cu_2(OH)_3Cl$): AB2C3D3_oP36_62_c_ac_cd_cd - POSCAR

```
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    ↪ x7,y7,z7 --params=6.03,1.13847429519,1.51243781095,0.8518,
    ↪ 0.5556,0.3094,0.2447,0.1951,0.5148,0.3502,0.5018,-0.0669,0.4666
    ↪ ,0.7279,-0.0594,0.5651,0.7879 & Pnma D_{2h}^{16} #62 (ac^4d^2)
    ↪ & oP36 & None & ClCu2H3O3 & Atacamite & J. B. Parise and B. G.
    ↪ Hyde, Acta Crystallogr. C 42, 1277-1280 (1986)
   1.00000000000000
   6.03000000000000   0.00000000000000   0.00000000000000
   0.00000000000000   6.86500000000000   0.00000000000000
   0.00000000000000   0.00000000000000   9.12000000000000
     Cl     Cu     H     O
      4      8     12    12
Direct
   0.85180000000000   0.25000000000000   0.55560000000000   Cl   (4c)
  -0.35180000000000   0.75000000000000   1.05560000000000   Cl   (4c)
  -0.85180000000000   0.75000000000000  -0.55560000000000   Cl   (4c)
   1.35180000000000   0.25000000000000  -0.05560000000000   Cl   (4c)
   0.00000000000000   0.00000000000000   0.00000000000000   Cu   (4a)
   0.50000000000000   0.00000000000000   0.50000000000000   Cu   (4a)
   0.00000000000000   0.50000000000000   0.00000000000000   Cu   (4a)
   0.50000000000000   0.50000000000000   0.50000000000000   Cu   (4a)
   0.30940000000000   0.25000000000000   0.24470000000000   Cu   (4c)
   0.19060000000000   0.75000000000000   0.74470000000000   Cu   (4c)
  -0.30940000000000   0.75000000000000  -0.24470000000000   Cu   (4c)
   0.80940000000000   0.25000000000000   0.25530000000000   Cu   (4c)
   0.19510000000000   0.25000000000000   0.51480000000000   H    (4c)
   0.30490000000000   0.75000000000000   1.01480000000000   H    (4c)
  -0.19510000000000   0.75000000000000  -0.51480000000000   H    (4c)
   0.69510000000000   0.25000000000000  -0.01480000000000   H    (4c)
  -0.06690000000000   0.46660000000000   0.72790000000000   H    (8d)
   0.56690000000000  -0.46660000000000   1.22790000000000   H    (8d)
   0.06690000000000   0.96660000000000  -0.72790000000000   H    (8d)
   0.43310000000000   0.03340000000000  -0.22790000000000   H    (8d)
   0.06690000000000  -0.46660000000000  -0.72790000000000   H    (8d)
   0.43310000000000   0.46660000000000  -0.22790000000000   H    (8d)
  -0.06690000000000   0.03340000000000   0.72790000000000   H    (8d)
   0.56690000000000   0.96660000000000   1.22790000000000   H    (8d)
   0.35020000000000   0.25000000000000   0.50180000000000   O    (4c)
   0.14980000000000   0.75000000000000   1.00180000000000   O    (4c)
  -0.35020000000000   0.75000000000000  -0.50180000000000   O    (4c)
   0.85020000000000   0.25000000000000  -0.00180000000000   O    (4c)
  -0.05940000000000   0.56510000000000   0.78790000000000   O    (8d)
   0.55940000000000  -0.56510000000000   1.28790000000000   O    (8d)
   0.05940000000000   1.06510000000000  -0.78790000000000   O    (8d)
   0.44060000000000  -0.06510000000000  -0.28790000000000   O    (8d)
   0.05940000000000  -0.56510000000000  -0.78790000000000   O    (8d)
   0.44060000000000   0.56510000000000  -0.28790000000000   O    (8d)
```

```
  −0.05940000000000   −0.06510000000000    0.78790000000000    O    (8d)
   0.55940000000000    1.06510000000000    1.28790000000000    O    (8d)
```

NH_4CdCl_3 ($E2_4$): AB3C_oP20_62_c_3c_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'CdCl3(NH4)'
_chemical_formula_sum 'Cd Cl3 (NH4)'

loop_
_publ_author_name
 'H. Brasseur'
 'L. Pauling'
_journal_name_full_name
;
 Journal of the American Chemical Society
;
_journal_volume 60
_journal_year 1938
_journal_page_first 2886
_journal_page_last 2890
_publ_Section_title
;
 The Crystal Structure of Ammonium Cadmium Chloride, NH$_{4}$CdCl$_{3}$
;

_aflow_title 'NH$_{4}$CdCl$_{3}$ ($E2_{4}$) Structure'
_aflow_proto 'AB3C_oP20_62_c_3c_c'
_aflow_params 'a,b/a,c/a,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3},x_{4},z_{4
      ↪ },x_{5},z_{5}'
_aflow_params_values '8.96,0.443080357143,1.65959821429,0.665,-0.054,
      ↪ 0.784,0.785,0.667,0.504,0.526,0.102,-0.07,0.18'
_aflow_Strukturbericht '$E2_{4}$'
_aflow_Pearson 'oP20'

_symmetry_space_group_name_H-M "P 21/n 21/m 21/a"
_symmetry_Int_Tables_number 62

_cell_length_a      8.96000
_cell_length_b      3.97000
_cell_length_c      14.87000
_cell_angle_alpha   90.00000
_cell_angle_beta    90.00000
_cell_angle_gamma   90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x,y+1/2,-z
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x,-y+1/2,z
8 x+1/2,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cd1 Cd 4 c 0.66500 0.25000 -0.05400 1.00000
Cl1 Cl 4 c 0.78400 0.25000 0.78500 1.00000
Cl2 Cl 4 c 0.66700 0.25000 0.50400 1.00000
Cl3 Cl 4 c 0.52600 0.25000 0.10200 1.00000
NH41 NH4 4 c -0.07000 0.25000 0.18000 1.00000
```

NH_4CdCl_3 ($E2_4$): AB3C_oP20_62_c_3c_c - POSCAR

```
AB3C_oP20_62_c_3c_c & a,b/a,c/a,x1,z1,x2,z2,x3,z3,x4,z4,x5,z5 --params=
      ↪ 8.96,0.443080357143,1.65959821429,0.665,-0.054,0.784,0.785,
      ↪ 0.667,0.504,0.526,0.102,-0.07,0.18 & Pnma D_{2h}^{16} #62 (c^5)
      ↪ & oP20 & $E2_{4}$ & CdCl3(NH4) & CdCl3(NH4) & H. Brasseur and
      ↪ L. Pauling, J. Am. Chem. Soc. 60, 2886-2890 (1938)
   1.00000000000000
   8.96000000000000    0.00000000000000    0.00000000000000
   0.00000000000000    3.97000000000000    0.00000000000000
   0.00000000000000    0.00000000000000   14.87000000000000
     Cd     Cl    NH4
      4     12      4
Direct
   0.66500000000000    0.25000000000000   -0.05400000000000   Cd    (4c)
  -0.16500000000000    0.75000000000000    0.44600000000000   Cd    (4c)
  -0.66500000000000    0.75000000000000    0.05400000000000   Cd    (4c)
   1.16500000000000    0.25000000000000    0.55400000000000   Cd    (4c)
   0.78400000000000    0.25000000000000    0.78500000000000   Cl    (4c)
  -0.28400000000000    0.75000000000000    1.28500000000000   Cl    (4c)
  -0.78400000000000    0.75000000000000   -0.78500000000000   Cl    (4c)
   1.28400000000000    0.25000000000000   -0.28500000000000   Cl    (4c)
   0.66700000000000    0.25000000000000    0.50400000000000   Cl    (4c)
  -0.16700000000000    0.75000000000000    1.00400000000000   Cl    (4c)
  -0.66700000000000    0.75000000000000   -0.50400000000000   Cl    (4c)
   1.16700000000000    0.25000000000000   -0.00400000000000   Cl    (4c)
   0.52600000000000    0.25000000000000    0.10200000000000   Cl    (4c)
  -0.02600000000000    0.75000000000000    0.60200000000000   Cl    (4c)
  -0.52600000000000    0.75000000000000   -0.10200000000000   Cl    (4c)
   1.02600000000000    0.25000000000000    0.39800000000000   Cl    (4c)
  -0.07000000000000    0.25000000000000    0.18000000000000   NH4   (4c)
   0.57000000000000    0.75000000000000    0.68000000000000   NH4   (4c)
   0.07000000000000    0.75000000000000   -0.18000000000000   NH4   (4c)
   0.43000000000000    0.25000000000000    0.32000000000000   NH4   (4c)
```

Berthierite ($FeSb_2S_4$, $E3_3$): AB4C2_oP28_62_c_4c_2c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Berthierite'
_chemical_formula_sum 'Fe S4 Sb2'

loop_
_publ_author_name
 'M. J. Buerger'
 'T. Hahn'
_journal_year 1953
_publ_Section_title
;
 The Crystal Structure of Berthierite, FeSb$_{2}$S$_{4}$
;

_aflow_title 'Berthierite (FeSb$_{2}$S$_{4}$, $E3_{3}$) Structure'
_aflow_proto 'AB4C2_oP28_62_c_4c_2c'
_aflow_params 'a,b/a,c/a,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3},x_{4},z_{4
      ↪ },x_{5},z_{5},x_{6},z_{6},x_{7},z_{7}'
_aflow_params_values '11.44,0.328671328671,1.23426573427,0.184,0.334,
      ↪ 0.695,0.728,0.076,0.184,0.274,0.492,-0.049,0.595,0.355,0.062,
      ↪ 0.537,0.614'
_aflow_Strukturbericht '$E3_{3}$'
_aflow_Pearson 'oP28'

_symmetry_space_group_name_H-M "P 21/n 21/m 21/a"
_symmetry_Int_Tables_number 62

_cell_length_a      11.44000
_cell_length_b      3.76000
_cell_length_c      14.12000
_cell_angle_alpha   90.00000
_cell_angle_beta    90.00000
_cell_angle_gamma   90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x,y+1/2,-z
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x,-y+1/2,z
8 x+1/2,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Fe1 Fe 4 c 0.18400 0.25000 0.33400 1.00000
S1 S 4 c 0.69500 0.25000 0.72800 1.00000
S2 S 4 c 0.07600 0.25000 0.18400 1.00000
S3 S 4 c 0.27400 0.25000 0.49200 1.00000
S4 S 4 c -0.04900 0.25000 0.59500 1.00000
Sb1 Sb 4 c 0.35500 0.25000 0.06200 1.00000
Sb2 Sb 4 c 0.53700 0.25000 0.61400 1.00000
```

Berthierite ($FeSb_2S_4$, $E3_3$): AB4C2_oP28_62_c_4c_2c - POSCAR

```
AB4C2_oP28_62_c_4c_2c & a,b/a,c/a,x1,z1,x2,z2,x3,z3,x4,z4,x5,z5,x6,z6,x7
      ↪ ,z7 --params=11.44,0.328671328671,1.23426573427,0.184,0.334,
      ↪ 0.695,0.728,0.076,0.184,0.274,0.492,-0.049,0.595,0.355,0.062,
      ↪ 0.537,0.614 & Pnma D_{2h}^{16} #62 (c^7) & oP28 & $E3_{3}$ &
      ↪ FeS4Sb2 & Berthierite & M. J. Buerger and T. Hahn, (1953)
   1.00000000000000
  11.44000000000000    0.00000000000000    0.00000000000000
   0.00000000000000    3.76000000000000    0.00000000000000
   0.00000000000000    0.00000000000000   14.12000000000000
     Fe      S     Sb
      4     16      8
Direct
   0.18400000000000    0.25000000000000    0.33400000000000   Fe    (4c)
   0.31600000000000    0.75000000000000    0.83400000000000   Fe    (4c)
  -0.18400000000000    0.75000000000000   -0.33400000000000   Fe    (4c)
   0.68400000000000    0.25000000000000    0.16600000000000   Fe    (4c)
   0.69500000000000    0.25000000000000    0.72800000000000   S     (4c)
  -0.19500000000000    0.75000000000000    1.22800000000000   S     (4c)
  -0.69500000000000    0.75000000000000   -0.72800000000000   S     (4c)
   1.19500000000000    0.25000000000000   -0.22800000000000   S     (4c)
   0.07600000000000    0.25000000000000    0.18400000000000   S     (4c)
   0.42400000000000    0.75000000000000    0.68400000000000   S     (4c)
  -0.07600000000000    0.75000000000000   -0.18400000000000   S     (4c)
   0.57600000000000    0.25000000000000    0.31600000000000   S     (4c)
   0.27400000000000    0.25000000000000    0.49200000000000   S     (4c)
   0.22600000000000    0.75000000000000    0.99200000000000   S     (4c)
  -0.27400000000000    0.75000000000000   -0.49200000000000   S     (4c)
   0.77400000000000    0.25000000000000    0.00800000000000   S     (4c)
  -0.04900000000000    0.25000000000000    0.59500000000000   S     (4c)
   0.54900000000000    0.75000000000000    1.09500000000000   S     (4c)
```

```
   0.04900000000000   0.75000000000000  −0.59500000000000   S   (4c)
   0.45100000000000   0.25000000000000  −0.09500000000000   S   (4c)
   0.35500000000000   0.25000000000000   0.06200000000000  Sb   (4c)
   0.14500000000000   0.75000000000000   0.56200000000000  Sb   (4c)
  −0.35500000000000   0.75000000000000  −0.06200000000000  Sb   (4c)
   0.85500000000000   0.25000000000000   0.43800000000000  Sb   (4c)
   0.53700000000000   0.25000000000000   0.61400000000000  Sb   (4c)
  −0.03700000000000   0.75000000000000   1.11400000000000  Sb   (4c)
  −0.53700000000000   0.75000000000000  −0.61400000000000  Sb   (4c)
   1.03700000000000   0.25000000000000  −0.11400000000000  Sb   (4c)
```

Chalcocyanite ($CuSO_4$): AB4C_oP24_62_a_2cd_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Chalcocyanite'
_chemical_formula_sum 'Cu O4 S'

loop_
_publ_author_name
 'M. Wildner'
 'G. Giester'
_journal_name_full_name
;
 Mineralogy and Petrology
;
_journal_volume 39
_journal_year 1988
_journal_page_first 201
_journal_page_last 209
_publ_Section_title
;
 Crystal structure refinements of synthetic chalcocyanite (CuSO$_{4}$)
     ↪ and zincosite (ZnSO$_{4}$)
;

_aflow_title 'Chalcocyanite (CuSO$_{4}$) Structure'
_aflow_proto 'AB4C_oP24_62_a_2cd_c'
_aflow_params 'a,b/a,c/a,x_{2},z_{2},x_{3},z_{3},x_{4},z_{4},x_{5},y_{5
    ↪ },z_{5}'
_aflow_params_values '8.409,0.797835652277,0.574741348555,0.1293,0.7353,
    ↪ 0.3646,0.4385,0.18363,0.44979,0.1328,0.0674,0.3083'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP24'

_symmetry_space_group_name_H-M "P 21/n 21/m 21/a"
_symmetry_Int_Tables_number 62

_cell_length_a     8.40900
_cell_length_b     6.70900
_cell_length_c     4.83300
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x,y+1/2,-z
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x,-y+1/2,z
8 x+1/2,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cu1 Cu 4 a 0.00000 0.00000 0.00000 1.00000
O1 O 4 c 0.12930 0.25000 0.73530 1.00000
O2 O 4 c 0.36460 0.25000 0.43850 1.00000
S1 S 4 c 0.18363 0.25000 0.44979 1.00000
O3 O 8 d 0.13280 0.06740 0.30830 1.00000
```

Chalcocyanite ($CuSO_4$): AB4C_oP24_62_a_2cd_c - POSCAR

```
AB4C_oP24_62_a_2cd_c & a,b/a,c/a,x2,z2,x3,z3,x4,z4,x5,y5,z5 --params=
    ↪ 8.409,0.797835652277,0.574741348555,0.1293,0.7353,0.3646,0.4385
    ↪ ,0.18363,0.44979,0.1328,0.0674,0.3083 & Pnma D_{2h}^{16} #62 (
    ↪ ac^3d) & oP24 & None & CuO4S & Chalcocyanite & M. Wildner and
    ↪ G. Giester, Mineral. Petrol. 39, 201-209 (1988)
   1.00000000000000
   8.40900000000000   0.00000000000000   0.00000000000000
   0.00000000000000   6.70900000000000   0.00000000000000
   0.00000000000000   0.00000000000000   4.83300000000000
     Cu    O    S
      4   16    4
Direct
   0.00000000000000   0.00000000000000   0.00000000000000   Cu   (4a)
   0.50000000000000   0.00000000000000   0.50000000000000   Cu   (4a)
   0.00000000000000   0.50000000000000   0.00000000000000   Cu   (4a)
   0.50000000000000   0.50000000000000   0.50000000000000   Cu   (4a)
   0.12930000000000   0.25000000000000   0.73530000000000    O   (4c)
   0.37070000000000   0.75000000000000   1.23530000000000    O   (4c)
  -0.12930000000000   0.75000000000000  -0.73530000000000    O   (4c)
   0.62930000000000   0.25000000000000  -0.23530000000000    O   (4c)
   0.36460000000000   0.25000000000000   0.43850000000000    O   (4c)
   0.13540000000000   0.75000000000000   0.93850000000000    O   (4c)
  -0.36460000000000   0.75000000000000  -0.43850000000000    O   (4c)
   0.86460000000000   0.25000000000000   0.06150000000000    O   (4c)
   0.13280000000000   0.06740000000000   0.30830000000000    O   (8d)
   0.36720000000000  -0.06740000000000   0.80830000000000    O   (8d)
  -0.13280000000000   0.56740000000000  -0.30830000000000    O   (8d)
   0.63280000000000   0.43260000000000   0.19170000000000    O   (8d)
  -0.13280000000000  -0.06740000000000  -0.30830000000000    O   (8d)
   0.63280000000000   0.06740000000000   0.19170000000000    O   (8d)
   0.13280000000000   0.43260000000000   0.30830000000000    O   (8d)
   0.36720000000000   0.56740000000000   0.80830000000000    O   (8d)
   0.18363000000000   0.25000000000000   0.44979000000000    S   (4c)
   0.31637000000000   0.75000000000000   0.94979000000000    S   (4c)
  -0.18363000000000   0.75000000000000  -0.44979000000000    S   (4c)
   0.68363000000000   0.25000000000000   0.05021000000000    S   (4c)
```

Rynersonite (Orthorhombic $CaTa_2O_6$): AB6C2_oP36_62_c_2c2d_d - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Rynersonite'
_chemical_formula_sum 'Ca O6 Ta2'

loop_
_publ_author_name
 'L. Jahnberg'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 71
_journal_year 1963
_journal_page_first 2548
_journal_page_last 2559
_publ_Section_title
;
 Crystal Structure of Orthorhombic CaTa$_{2}$O$_{6}$
;

_aflow_title 'Rynersonite (Orthorhombic CaTa$_{2}$O$_{6}$) Structure'
_aflow_proto 'AB6C2_oP36_62_c_2c2d_d'
_aflow_params 'a,b/a,c/a,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3},x_{4},y_{4
    ↪ },z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{6}'
_aflow_params_values '11.068,0.678080954102,0.485905312613,0.042,0.54,
    ↪ 0.146,-0.033,0.878,0.838,-0.024,0.035,0.225,0.213,0.049,0.383,
    ↪ 0.1412,-0.0056,0.0376'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP36'

_symmetry_space_group_name_H-M "P 21/n 21/m 21/a"
_symmetry_Int_Tables_number 62

_cell_length_a     11.06800
_cell_length_b     7.50500
_cell_length_c     5.37800
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x,y+1/2,-z
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x,-y+1/2,z
8 x+1/2,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ca1 Ca 4 c 0.04200 0.25000 0.54000 1.00000
O1 O 4 c 0.14600 0.25000 -0.03300 1.00000
O2 O 4 c 0.87800 0.25000 0.83800 1.00000
O3 O 8 d -0.02400 0.03500 0.22500 1.00000
O4 O 8 d 0.21300 0.04900 0.38300 1.00000
Ta1 Ta 8 d 0.14120 -0.00560 0.03760 1.00000
```

Rynersonite (Orthorhombic $CaTa_2O_6$): AB6C2_oP36_62_c_2c2d_d - POSCAR

```
AB6C2_oP36_62_c_2c2d_d & a,b/a,c/a,x1,z1,x2,z2,x3,z3,x4,y4,z4,x5,y5,z5,
    ↪ x6,y6,z6 --params=11.068,0.678080954102,0.485905312613,0.042,
    ↪ 0.54,0.146,-0.033,0.878,0.838,-0.024,0.035,0.225,0.213,0.049,
    ↪ 0.383,0.1412,-0.0056,0.0376 & Pnma D_{2h}^{16} #62 (c^3d^3) &
    ↪ oP36 & None & CaO6Ta2 & Rynersonite & L. Jahnberg, Acta Chem.
    ↪ Scand. 71, 2548-2559 (1963)
   1.00000000000000
  11.06800000000000   0.00000000000000   0.00000000000000
   0.00000000000000   7.50500000000000   0.00000000000000
   0.00000000000000   0.00000000000000   5.37800000000000
     Ca    O   Ta
      4   24    8
```

```
Direct
    0.04200000000000    0.25000000000000    0.54000000000000   Ca    (4c)
    0.45800000000000    0.75000000000000    1.04000000000000   Ca    (4c)
   -0.04200000000000    0.75000000000000   -0.54000000000000   Ca    (4c)
    0.54200000000000    0.25000000000000   -0.04000000000000   Ca    (4c)
    0.14600000000000    0.25000000000000   -0.03300000000000   O     (4c)
    0.35400000000000    0.75000000000000    0.46700000000000   O     (4c)
   -0.14600000000000    0.75000000000000    0.03300000000000   O     (4c)
    0.64600000000000    0.25000000000000    0.53300000000000   O     (4c)
    0.87800000000000    0.25000000000000    0.83800000000000   O     (4c)
   -0.37800000000000    0.75000000000000    1.33800000000000   O     (4c)
   -0.87800000000000    0.75000000000000   -0.83800000000000   O     (4c)
    1.37800000000000    0.25000000000000   -0.33800000000000   O     (4c)
   -0.02400000000000    0.03500000000000    0.22500000000000   O     (8d)
    0.52400000000000   -0.03500000000000    0.72500000000000   O     (8d)
    0.02400000000000    0.53500000000000   -0.22500000000000   O     (8d)
    0.47600000000000    0.46500000000000    0.27500000000000   O     (8d)
    0.02400000000000   -0.03500000000000   -0.22500000000000   O     (8d)
    0.47600000000000    0.03500000000000    0.27500000000000   O     (8d)
   -0.02400000000000    0.46500000000000    0.22500000000000   O     (8d)
    0.52400000000000    0.53500000000000    0.72500000000000   O     (8d)
    0.21300000000000    0.04900000000000    0.38300000000000   O     (8d)
    0.28700000000000   -0.04900000000000    0.88300000000000   O     (8d)
   -0.21300000000000    0.54900000000000   -0.38300000000000   O     (8d)
    0.71300000000000    0.45100000000000    0.11700000000000   O     (8d)
   -0.21300000000000   -0.04900000000000   -0.38300000000000   O     (8d)
    0.71300000000000    0.04900000000000    0.11700000000000   O     (8d)
    0.21300000000000    0.45100000000000    0.38300000000000   O     (8d)
    0.28700000000000    0.54900000000000    0.88300000000000   O     (8d)
    0.14120000000000   -0.00560000000000    0.03760000000000   Ta    (8d)
    0.35880000000000    0.00560000000000    0.53760000000000   Ta    (8d)
   -0.14120000000000    0.49440000000000   -0.03760000000000   Ta    (8d)
    0.64120000000000    0.50560000000000    0.46240000000000   Ta    (8d)
   -0.14120000000000    0.00560000000000   -0.03760000000000   Ta    (8d)
    0.64120000000000   -0.00560000000000    0.46240000000000   Ta    (8d)
    0.14120000000000    0.50560000000000    0.03760000000000   Ta    (8d)
    0.35880000000000    0.49440000000000    0.53760000000000   Ta    (8d)
```

Copper (II) Azide [$Cu(N_3)_2$]: AB6_oP28_62_c_6c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Copper (ii) azide'
_chemical_formula_sum 'Cu N6'

loop_
_publ_author_name
 'I. Agrell'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 21
_journal_year 1967
_journal_page_first 2647
_journal_page_last 2658
_publ_Section_title
;
 The Crystal Structure of Cu(N$_{3}$)$_{2}$
;

_aflow_title 'Copper (II) Azide [Cu(N$_{3}$)$_{2}$] Structure'
_aflow_proto 'AB6_oP28_62_c_6c'
_aflow_params 'a,b/a,c/a,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3},x_{4},z_{4
    ↪ },x_{5},z_{5},x_{6},z_{6},x_{7},z_{7}'
_aflow_params_values '13.481,0.228766411987,0.673243824642,0.39692,
    ↪ 0.58022,0.81158,0.66343,0.7465,0.59084,0.6767,0.50317,0.52408,
    ↪ 0.35102,0.47071,0.24672,0.42026,0.15152'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP28'

_symmetry_space_group_name_H-M "P 21/n 21/m 21/a"
_symmetry_Int_Tables_number 62

_cell_length_a    13.48100
_cell_length_b    3.08400
_cell_length_c    9.07600
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x,y+1/2,-z
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x,-y+1/2,z
8 x+1/2,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cu1 Cu 4 c 0.39692 0.25000 0.58022 1.00000
N1 N 4 c 0.81158 0.25000 0.66343 1.00000
N2 N 4 c 0.74650 0.25000 0.59084 1.00000
N3 N 4 c 0.67670 0.25000 0.50317 1.00000
N4 N 4 c 0.52408 0.25000 0.35102 1.00000
N5 N 4 c 0.47071 0.25000 0.24672 1.00000
N6 N 4 c 0.42026 0.25000 0.15152 1.00000
```

Copper (II) Azide [$Cu(N_3)_2$]: AB6_oP28_62_c_6c - POSCAR

```
AB6_oP28_62_c_6c & a,b/a,c/a,x1,z1,x2,z2,x3,z3,x4,z4,x5,z5,x6,z6,x7,z7
    ↪ --params=13.481,0.228766411987,0.673243824642,0.39692,0.58022,
    ↪ 0.81158,0.66343,0.7465,0.59084,0.6767,0.50317,0.52408,0.35102,
    ↪ 0.47071,0.24672,0.42026,0.15152 & Pnma D_{2h}^{16} #62 (c^7) &
    ↪ oP28 & None & CuN6 & Copper (ii) azide & I. Agrell, Acta Chem.
    ↪ Scand. 21, 2647-2658 (1967)
   1.00000000000000
  13.48100000000000    0.00000000000000    0.00000000000000
   0.00000000000000    3.08400000000000    0.00000000000000
   0.00000000000000    0.00000000000000    9.07600000000000
     Cu      N
      4     24
Direct
    0.39692000000000    0.25000000000000    0.58022000000000   Cu    (4c)
    0.10308000000000    0.75000000000000    1.08022000000000   Cu    (4c)
   -0.39692000000000    0.75000000000000   -0.58022000000000   Cu    (4c)
    0.89692000000000    0.25000000000000   -0.08022000000000   Cu    (4c)
    0.81158000000000    0.25000000000000    0.66343000000000   N     (4c)
   -0.31158000000000    0.75000000000000    1.16343000000000   N     (4c)
   -0.81158000000000    0.75000000000000   -0.66343000000000   N     (4c)
    1.31158000000000    0.25000000000000   -0.16343000000000   N     (4c)
    0.74650000000000    0.25000000000000    0.59084000000000   N     (4c)
   -0.24650000000000    0.75000000000000    1.09084000000000   N     (4c)
   -0.74650000000000    0.75000000000000   -0.59084000000000   N     (4c)
    1.24650000000000    0.25000000000000   -0.09084000000000   N     (4c)
    0.67670000000000    0.25000000000000    0.50317000000000   N     (4c)
   -0.17670000000000    0.75000000000000    1.00317000000000   N     (4c)
   -0.67670000000000    0.75000000000000   -0.50317000000000   N     (4c)
    1.17670000000000    0.25000000000000   -0.00317000000000   N     (4c)
    0.52408000000000    0.25000000000000    0.35102000000000   N     (4c)
   -0.02408000000000    0.75000000000000    0.85102000000000   N     (4c)
   -0.52408000000000    0.75000000000000   -0.35102000000000   N     (4c)
    1.02408000000000    0.25000000000000    0.14898000000000   N     (4c)
    0.47071000000000    0.25000000000000    0.24672000000000   N     (4c)
    0.02929000000000    0.75000000000000    0.74672000000000   N     (4c)
   -0.47071000000000    0.75000000000000   -0.24672000000000   N     (4c)
    0.97071000000000    0.25000000000000    0.25328000000000   N     (4c)
    0.42026000000000    0.25000000000000    0.15152000000000   N     (4c)
    0.07974000000000    0.75000000000000    0.65152000000000   N     (4c)
   -0.42026000000000    0.75000000000000   -0.15152000000000   N     (4c)
    0.92026000000000    0.25000000000000    0.34848000000000   N     (4c)
```

Diaspore (AlOOH, $E0_2$): ABC2_oP16_62_c_c_2c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Diaspore'
_chemical_formula_sum 'Al H O2'

loop_
_publ_author_name
 'R. J. Hill'
_journal_name_full_name
;
 Physics and Chemistry of Minerals
;
_journal_volume 5
_journal_year 1979
_journal_page_first 179
_journal_page_last 200
_publ_Section_title
;
 Crystal Structure Refinement and Electron Density Distribution in
    ↪ Diaspore
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'Diaspore (AlOOH, $E0_{2}$) Structure'
_aflow_proto 'ABC2_oP16_62_c_c_2c'
_aflow_params 'a,b/a,c/a,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3},x_{4},z_{4
    ↪ }'
_aflow_params_values '9.4253,0.301868375542,0.466902910252,0.1445,-
    ↪ 0.04472,0.80108,0.28766,-0.05338,0.80286,-0.0876,0.5905'
_aflow_Strukturbericht '$E0_{2}$'
_aflow_Pearson 'oP16'

_symmetry_space_group_name_H-M "P 21/n 21/m 21/a"
_symmetry_Int_Tables_number 62

_cell_length_a    9.42530
_cell_length_b    2.84520
_cell_length_c    4.40070
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x,y+1/2,-z
4 -x+1/2,-y,z+1/2
```

```
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x,-y+1/2,z
8 x+1/2,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Al1 Al 4 c 0.14450 0.25000 -0.04472 1.00000
H1 H 4 c 0.80108 0.25000 0.28766 1.00000
O1 O 4 c -0.05338 0.25000 0.80286 1.00000
O2 O 4 c -0.08760 0.25000 0.59050 1.00000
```

Diaspore (AlOOH, $E0_2$): ABC2_oP16_62_c_c_2c - POSCAR

```
ABC2_oP16_62_c_c_2c & a,b/a,c/a,x1,z1,x2,z2,x3,z3,x4,z4 --params=9.4253,
    ↪ 0.301868375542,0.466902910252,0.1445,-0.04472,0.80108,0.28766,-
    ↪ 0.05338,0.80286,-0.0876,0.5905 & Pnma D_{2h}^{16} #62 (c^4) &
    ↪ oP16 & $E0_{2}$ & AlHO2 & Diaspore & R. J. Hill, Phys. Chem.
    ↪ Miner. 5, 179-200 (1979)
   1.00000000000000
   9.42530000000000   0.00000000000000   0.00000000000000
   0.00000000000000   2.84520000000000   0.00000000000000
   0.00000000000000   0.00000000000000   4.40070000000000
     Al    H    O
      4    4    8
Direct
   0.14450000000000   0.25000000000000  -0.04472000000000   Al   (4c)
   0.35550000000000   0.75000000000000   0.45528000000000   Al   (4c)
  -0.14450000000000   0.75000000000000   0.04472000000000   Al   (4c)
   0.64450000000000   0.25000000000000   0.54472000000000   Al   (4c)
   0.80108000000000   0.25000000000000   0.28766000000000   H    (4c)
  -0.30108000000000   0.75000000000000   0.78766000000000   H    (4c)
  -0.80108000000000   0.75000000000000  -0.28766000000000   H    (4c)
   1.30108000000000   0.25000000000000   0.21234000000000   H    (4c)
  -0.05338000000000   0.25000000000000   0.80286000000000   O    (4c)
   0.55338000000000   0.75000000000000   1.30286000000000   O    (4c)
   0.05338000000000   0.75000000000000  -0.80286000000000   O    (4c)
   0.44662000000000   0.25000000000000  -0.30286000000000   O    (4c)
  -0.08760000000000   0.25000000000000   0.59050000000000   O    (4c)
   0.58760000000000   0.75000000000000   1.09050000000000   O    (4c)
   0.08760000000000   0.75000000000000  -0.59050000000000   O    (4c)
   0.41240000000000   0.25000000000000  -0.09050000000000   O    (4c)
```

α-Potassium Nitrate (KNO_3) I: ABC3_oP20_62_c_c_cd - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'KNO3'
_chemical_formula_sum 'K N O3'

loop_
_publ_author_name
 'J. K. Nimmo'
 'B. W. Lucas'
_journal_name_full_name
;
 Journal of Physics C: Solid State Physics
;
_journal_volume 6
_journal_year 1973
_journal_page_first 201
_journal_page_last 211
_publ_Section_title
;
A neutron diffraction determination of the crystal structure of $\
    ↪ alpha$-phase potassium nitrate at 25$^{\circ}$C and 100$^{\
    ↪ circ}$C
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title '$\alpha$-Potassium Nitrate (KNO$_{3}$) I Structure'
_aflow_proto 'ABC3_oP20_62_c_c_cd'
_aflow_params 'a,b/a,c/a,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3},x_{4},y_{4
    ↪ },z_{4}'
_aflow_params_values '6.4213,0.842804416551,1.425988507,0.7568,0.4166,-
    ↪ 0.0848,0.7548,-0.0893,0.8902,-0.0849,0.4492,0.6866'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP20'

_symmetry_space_group_name_H-M "P 21/n 21/m 21/a"
_symmetry_Int_Tables_number 62

_cell_length_a    6.42130
_cell_length_b    5.41190
_cell_length_c    9.15670
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x,y+1/2,-z
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x,-y+1/2,z
8 x+1/2,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
K1 K 4 c 0.75680 0.25000 0.41660 1.00000
N1 N 4 c -0.08480 0.25000 0.75480 1.00000
O1 O 4 c -0.08930 0.25000 0.89020 1.00000
O2 O 8 d -0.08490 0.44920 0.68660 1.00000
```

α-Potassium Nitrate (KNO_3) I: ABC3_oP20_62_c_c_cd - POSCAR

```
ABC3_oP20_62_c_c_cd & a,b/a,c/a,x1,z1,x2,z2,x3,z3,x4,y4,z4 --params=
    ↪ 6.4213,0.842804416551,1.425988507,0.7568,0.4166,-0.0848,0.7548
    ↪ ,-0.0893,0.8902,-0.0849,0.4492,0.6866 & Pnma D_{2h}^{16} #62 (c
    ↪ ^3d) & oP20 & None & KNO3 & KNO3 & J. K. Nimmo and B. W. Lucas,
    ↪ J. Phys. C: Solid State Phys. 6, 201-211 (1973)
   1.00000000000000
   6.42130000000000   0.00000000000000   0.00000000000000
   0.00000000000000   5.41190000000000   0.00000000000000
   0.00000000000000   0.00000000000000   9.15670000000000
     K    N    O
     4    4   12
Direct
   0.75680000000000   0.25000000000000   0.41660000000000   K   (4c)
  -0.25680000000000   0.75000000000000   0.91660000000000   K   (4c)
  -0.75680000000000   0.75000000000000  -0.41660000000000   K   (4c)
   1.25680000000000   0.25000000000000   0.08340000000000   K   (4c)
  -0.08480000000000   0.25000000000000   0.75480000000000   N   (4c)
   0.58480000000000   0.75000000000000   1.25480000000000   N   (4c)
   0.08480000000000   0.75000000000000  -0.75480000000000   N   (4c)
   0.41520000000000   0.25000000000000  -0.25480000000000   N   (4c)
  -0.08930000000000   0.25000000000000   0.89020000000000   O   (4c)
   0.58930000000000   0.75000000000000   1.39020000000000   O   (4c)
   0.08930000000000   0.75000000000000  -0.89020000000000   O   (4c)
   0.41070000000000   0.25000000000000  -0.39020000000000   O   (4c)
  -0.08490000000000   0.44920000000000   0.68660000000000   O   (8d)
   0.58490000000000  -0.44920000000000   1.18660000000000   O   (8d)
   0.08490000000000   0.94920000000000  -0.68660000000000   O   (8d)
   0.41510000000000   0.05080000000000  -0.18660000000000   O   (8d)
   0.08490000000000  -0.44920000000000  -0.68660000000000   O   (8d)
   0.41510000000000   0.44920000000000  -0.18660000000000   O   (8d)
  -0.08490000000000   0.05080000000000   0.68660000000000   O   (8d)
   0.58490000000000   0.94920000000000   1.18660000000000   O   (8d)
```

NH_4NO_3 III ($G0_{10}$): ABC3_oP20_62_c_c_cd - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'N(NH4)O3'
_chemical_formula_sum 'N (NH4) O3'

loop_
_publ_author_name
 'T. H. Goodwin'
 'J. Whetstone'
_journal_name_full_name
;
 Journal of the Chemical Society
;
_journal_volume
_journal_year 1947
_journal_page_first 1455
_journal_page_last 1461
_publ_Section_title
;
 The crystal structure of ammonium nitrate III, and atomic scattering
    ↪ factors in ionic crystals
;

_aflow_title 'NH$_{4}$NO$_{3}$ III ($G0_{10}$) Structure'
_aflow_proto 'ABC3_oP20_62_c_c_cd'
_aflow_params 'a,b/a,c/a,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3},x_{4},y_{4
    ↪ },z_{4}'
_aflow_params_values '7.65,0.762091503268,0.933333333333,0.145,0.14,0.02
    ↪ ,0.68,0.085,0.3,0.175,0.06,0.06'
_aflow_Strukturbericht '$G0_{10}$'
_aflow_Pearson 'oP20'

_symmetry_space_group_name_H-M "P 21/n 21/m 21/a"
_symmetry_Int_Tables_number 62

_cell_length_a    7.65000
_cell_length_b    5.83000
_cell_length_c    7.14000
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
```

```
2 x+1/2,-y+1/2,-z+1/2
3 -x,y+1/2,-z
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x,-y+1/2,z
8 x+1/2,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
N1 N 4 c 0.14500 0.25000 0.14000 1.00000
NH41 NH4 4 c 0.02000 0.25000 0.68000 1.00000
O1 O 4 c 0.08500 0.25000 0.30000 1.00000
O2 O 8 d 0.17500 0.06000 0.06000 1.00000
```

NH_4NO_3 III ($G0_{10}$): ABC3_oP20_62_c_c_cd - POSCAR

```
ABC3_oP20_62_c_c_cd & a,b/a,c/a,x1,z1,x2,z2,x3,z3,x4,y4,z4 --params=7.65
    ↪ ,0.762091503268,0.933333333333,0.145,0.14,0.02,0.68,0.085,0.3,
    ↪ 0.175,0.06,0.06 & Pnma D_{2h}^{16} #62 (c^3d) & oP20 & $G0_{10}
    ↪ $ & N(NH4)O3 & N(NH4)O3 & T. H. Goodwin and J. Whetstone, J.
    ↪ Chem. Soc., 1455-1461 (1947)
   1.00000000000000
   7.65000000000000   0.00000000000000   0.00000000000000
   0.00000000000000   5.83000000000000   0.00000000000000
   0.00000000000000   0.00000000000000   7.14000000000000
     N    NH4     O
     4     4    12
Direct
   0.14500000000000   0.25000000000000   0.14000000000000   N    (4c)
   0.35500000000000   0.75000000000000   0.64000000000000   N    (4c)
  -0.14500000000000   0.75000000000000  -0.14000000000000   N    (4c)
   0.64500000000000   0.25000000000000   0.36000000000000   N    (4c)
   0.02000000000000   0.25000000000000   0.68000000000000   NH4  (4c)
   0.48000000000000   0.75000000000000   1.18000000000000   NH4  (4c)
  -0.02000000000000   0.75000000000000  -0.68000000000000   NH4  (4c)
   0.52000000000000   0.25000000000000  -0.18000000000000   NH4  (4c)
   0.08500000000000   0.25000000000000   0.30000000000000   O    (4c)
   0.41500000000000   0.75000000000000   0.80000000000000   O    (4c)
  -0.08500000000000   0.75000000000000  -0.30000000000000   O    (4c)
   0.58500000000000   0.25000000000000   0.20000000000000   O    (4c)
   0.17500000000000   0.06000000000000   0.06000000000000   O    (8d)
   0.32500000000000  -0.06000000000000   0.56000000000000   O    (8d)
  -0.17500000000000   0.56000000000000  -0.06000000000000   O    (8d)
   0.67500000000000   0.44000000000000   0.44000000000000   O    (8d)
  -0.17500000000000  -0.06000000000000  -0.06000000000000   O    (8d)
   0.67500000000000   0.06000000000000   0.44000000000000   O    (8d)
   0.17500000000000   0.44000000000000   0.06000000000000   O    (8d)
   0.32500000000000   0.56000000000000   0.56000000000000   O    (8d)
```

Aragonite ($CaCO_3$, $G0_2$): ABC3_oP20_62_c_c_cd - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Aragonite'
_chemical_formula_sum 'C Ca O3'

loop_
_publ_author_name
 'J. P. R. {de Villiers}'
_journal_name_full_name
;
 American Mineralogist
;
_journal_volume 56
_journal_year 1971
_journal_page_first 758
_journal_page_last 767
_publ_Section_title
;
 Crystal Structures of Aragonite, Strontianite, and Witherite
;

# Found in Aragonite, {Mineral Database}

_aflow_title 'Aragonite (CaCO$_{3}$, $G0_{2}$) Structure'
_aflow_proto 'ABC3_oP20_62_c_c_cd'
_aflow_params 'a,b/a,c/a,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3},x_{4},y_{4
    ↪ },z_{4}'
_aflow_params_values '5.7404,0.864295171068,1.38789979792,0.4138,0.2622,
    ↪ 0.2597,-0.085,0.4038,0.4225,0.414,0.4736,0.181'
_aflow_Strukturbericht '$G0_{2}$'
_aflow_Pearson 'oP20'

_symmetry_space_group_name_H-M "P 21/n 21/m 21/a"
_symmetry_Int_Tables_number 62

_cell_length_a    5.74040
_cell_length_b    4.96140
_cell_length_c    7.96710
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x,y+1/2,-z
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x,-y+1/2,z
8 x+1/2,y,-z+1/2

loop_
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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
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_atom_site_fract_z
_atom_site_occupancy
C1 C 4 c 0.41380 0.25000 0.26220 1.00000
Ca1 Ca 4 c 0.25970 0.25000 -0.08500 1.00000
O1 O 4 c 0.40380 0.25000 0.42250 1.00000
O2 O 8 d 0.41400 0.47360 0.18100 1.00000
```

Aragonite ($CaCO_3$, $G0_2$): ABC3_oP20_62_c_c_cd - POSCAR

```
ABC3_oP20_62_c_c_cd & a,b/a,c/a,x1,z1,x2,z2,x3,z3,x4,y4,z4 --params=
    ↪ 5.7404,0.864295171068,1.38789979792,0.4138,0.2622,0.2597,-0.085
    ↪ ,0.4038,0.4225,0.414,0.4736,0.181 & Pnma D_{2h}^{16} #62 (c^3d)
    ↪ & oP20 & $G0_{2}$ & CaCO3 & Aragonite & J. P. R. {de Villiers
    ↪ }, Am. Mineral. 56, 758-767 (1971)
   1.00000000000000
   5.74040000000000   0.00000000000000   0.00000000000000
   0.00000000000000   4.96140000000000   0.00000000000000
   0.00000000000000   0.00000000000000   7.96710000000000
     C    Ca     O
     4     4    12
Direct
   0.41380000000000   0.25000000000000   0.26220000000000   C    (4c)
   0.08620000000000   0.75000000000000   0.76220000000000   C    (4c)
  -0.41380000000000   0.75000000000000  -0.26220000000000   C    (4c)
   0.91380000000000   0.25000000000000   0.23780000000000   C    (4c)
   0.25970000000000   0.25000000000000  -0.08500000000000   Ca   (4c)
   0.24030000000000   0.75000000000000   0.41500000000000   Ca   (4c)
  -0.25970000000000   0.75000000000000   0.08500000000000   Ca   (4c)
   0.75970000000000   0.25000000000000   0.58500000000000   Ca   (4c)
   0.40380000000000   0.25000000000000   0.42250000000000   O    (4c)
   0.09620000000000   0.75000000000000   0.92250000000000   O    (4c)
  -0.40380000000000   0.75000000000000  -0.42250000000000   O    (4c)
   0.90380000000000   0.25000000000000   0.07750000000000   O    (4c)
   0.41400000000000   0.47360000000000   0.18100000000000   O    (8d)
   0.08600000000000  -0.47360000000000   0.68100000000000   O    (8d)
  -0.41400000000000   0.97360000000000  -0.18100000000000   O    (8d)
   0.91400000000000   0.02640000000000   0.31900000000000   O    (8d)
  -0.41400000000000  -0.47360000000000  -0.18100000000000   O    (8d)
   0.91400000000000   0.47360000000000   0.31900000000000   O    (8d)
   0.41400000000000   0.02640000000000   0.18100000000000   O    (8d)
   0.08600000000000   0.97360000000000   0.68100000000000   O    (8d)
```

Epididymite ($BeHNaO_8Si_3$, $S4_7$): ABCD8E3_oP112_62_d_2c_d_4c6d_3d - CIF

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_chemical_formula_sum 'Be H Na O8 Si3'

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_publ_author_name
 'G. {Diego Gatta}'
 'N. Rotiroti'
 'G. J. {McIntyre}'
 'A. Guastoni'
 'F. Nestola'
_journal_name_full_name
;
 American Mineralogist
;
_journal_volume 93
_journal_year 2008
_journal_page_first 1158
_journal_page_last 1165
_publ_Section_title
;
 New insights into the crystal chemistry of epididymite and eudidymite
    ↪ from Malosa, Malawi: A single-crystal neutron diffraction
    ↪ study
;

# Found in The American Mineralogist Crystal Structure Database, 2003

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    ↪ x_{9},y_{9},z_{9},x_{10},y_{10},z_{10},x_{11},y_{11},z_{11},x_{
    ↪ 12},y_{12},z_{12},x_{13},y_{13},z_{13},x_{14},y_{14},z_{14},x_{
    ↪ 15},y_{15},z_{15},x_{16},y_{16},z_{16},x_{17},y_{17},z_{17}'
_aflow_params_values '12.7334,1.07039753719,0.576962947838,0.00647,
    ↪ 0.58561,0.0134,0.38486,0.12005,1e-05,0.37761,0.76529,0.37,
    ↪ 0.26883,0.05713,0.48974,0.49472,-0.00042,0.34273,0.10285,
    ↪ 0.06934,0.50192,0.21704,0.1224,0.21004,0.06068,0.06612,0.00658,
    ↪ 0.23777,0.11405,0.84938,0.43489,0.06457,0.78946,0.30497,0.1234,
```

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      ↪ 0.52074 ,0.41552 ,0.063 ,0.22624 ,0.15908 ,0.1378 ,0.01543 ,0.3426 ,
      ↪ 0.13635 ,0.73115 ,0.33173 ,0.13741 ,0.30525'
_aflow_Strukturbericht '$S4_{7}$'
_aflow_Pearson 'oP112'

_symmetry_space_group_name_H-M "P 21/n 21/m 21/a"
_symmetry_Int_Tables_number 62

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_cell_length_c      7.34670
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_space_group_symop_operation_xyz
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2 x+1/2,-y+1/2,-z+1/2
3 -x,y+1/2,-z
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x,-y+1/2,z
8 x+1/2,y,-z+1/2

loop_
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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
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H2 H 4 c 0.01340 0.25000 0.38486 1.00000
O1 O 4 c 0.12005 0.25000 0.00001 1.00000
O2 O 4 c 0.37761 0.25000 0.76529 1.00000
O3 O 4 c 0.37000 0.25000 0.26883 1.00000
O4 O 4 c 0.05713 0.25000 0.48974 1.00000
Be1 Be 8 d 0.49472 -0.00042 0.34273 1.00000
Na1 Na 8 d 0.10285 0.06934 0.50192 1.00000
O5 O 8 d 0.21704 0.12240 0.21004 1.00000
O6 O 8 d 0.06068 0.06612 0.00658 1.00000
O7 O 8 d 0.23777 0.11405 0.84938 1.00000
O8 O 8 d 0.43489 0.06457 0.78946 1.00000
O9 O 8 d 0.30497 0.12340 0.52074 1.00000
O10 O 8 d 0.41552 0.06300 0.22624 1.00000
Si1 Si 8 d 0.15908 0.13780 0.01543 1.00000
Si2 Si 8 d 0.34260 0.13635 0.73115 1.00000
Si3 Si 8 d 0.33173 0.13741 0.30525 1.00000
```

Epididymite ($BeHNaO_8Si_3$, $S4_7$): ABCD8E3_oP112_62_d_2c_d_4c6d_3d - POSCAR

```
ABCD8E3_oP112_62_d_2c_d_4c6d_3d & a,b/a,c/a,x1,z1,x2,z2,x3,z3,x4,z4,x5,
      ↪ z5,x6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10,x11,y11,z11,x12
      ↪ ,y12,z12,x13,y13,z13,x14,y14,z14,x15,y15,z15,x16,y16,z16,x17,
      ↪ y17,z17 --params=12.7334,1.07039753719,0.576962947838,0.00647,
      ↪ 0.58561,0.0134,0.38486,0.12005,1e-05,0.37761,0.76529,0.37,
      ↪ 0.26883,0.05713,0.48974,0.49472,-0.00042,0.34273,0.10285,
      ↪ 0.06934,0.50192,0.21704,0.1224,0.21004,0.06068,0.06612,0.00658,
      ↪ 0.23777,0.11405,0.84938,0.43489,0.06457,0.78946,0.30497,0.1234,
      ↪ 0.52074,0.41552,0.063,0.22624,0.15908,0.1378,0.01543,0.3426,
      ↪ 0.13635,0.73115,0.33173,0.13741,0.30525 & Pnma D_{2h}^{16} #62
      ↪ (c^6d^11) & oP112 & $S4_{7}$ & BeHNaO8Si3 & Epididymite & G. {
      ↪ Diego Gatta} et al., Am. Mineral. 93, 1158-1165 (2008)
   1.00000000000000
  12.73340000000000    0.00000000000000    0.00000000000000
   0.00000000000000   13.62980000000000    0.00000000000000
   0.00000000000000    0.00000000000000    7.34670000000000
     Be      H      Na      O      Si
      8      8       8      64     24
Direct
   0.49472000000000   -0.00042000000000    0.34273000000000   Be    (8d)
   0.00528000000000    0.00042000000000    0.84273000000000   Be    (8d)
  -0.49472000000000    0.49958000000000   -0.34273000000000   Be    (8d)
   0.99472000000000    0.50042000000000    0.15727000000000   Be    (8d)
  -0.49472000000000    0.00042000000000   -0.34273000000000   Be    (8d)
   0.99472000000000   -0.00042000000000    0.15727000000000   Be    (8d)
   0.49472000000000    0.50042000000000    0.34273000000000   Be    (8d)
   0.00528000000000    0.49958000000000    0.84273000000000   Be    (8d)
   0.00647000000000    0.25000000000000    0.58561000000000   H     (4c)
   0.49353000000000    0.75000000000000    1.08561000000000   H     (4c)
  -0.00647000000000    0.75000000000000   -0.58561000000000   H     (4c)
   0.50647000000000    0.25000000000000   -0.08561000000000   H     (4c)
   0.01340000000000    0.25000000000000    0.38486000000000   H     (4c)
   0.48660000000000    0.75000000000000    0.88486000000000   H     (4c)
  -0.01340000000000    0.75000000000000   -0.38486000000000   H     (4c)
   0.51340000000000    0.25000000000000    0.11514000000000   H     (4c)
   0.10285000000000    0.06934000000000    0.50192000000000   Na    (8d)
   0.39715000000000   -0.06934000000000    1.00192000000000   Na    (8d)
  -0.10285000000000    0.56934000000000   -0.50192000000000   Na    (8d)
   0.60285000000000    0.43066000000000   -0.00192000000000   Na    (8d)
  -0.10285000000000   -0.06934000000000   -0.50192000000000   Na    (8d)
   0.60285000000000    0.06934000000000   -0.00192000000000   Na    (8d)
   0.10285000000000    0.43066000000000    0.50192000000000   Na    (8d)
   0.39715000000000    0.56934000000000    1.00192000000000   Na    (8d)
   0.12005000000000    0.25000000000000    0.00001000000000   O     (4c)
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   0.62005000000000    0.25000000000000    0.49999000000000   O     (4c)
   0.37761000000000    0.25000000000000    0.76529000000000   O     (4c)
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   0.87761000000000    0.25000000000000   -0.26529000000000   O     (4c)
   0.37000000000000    0.25000000000000    0.26883000000000   O     (4c)
   0.13000000000000    0.75000000000000    0.76883000000000   O     (4c)
  -0.37000000000000    0.75000000000000   -0.26883000000000   O     (4c)
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   0.05713000000000    0.25000000000000    0.48974000000000   O     (4c)
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   0.55713000000000    0.25000000000000    0.01026000000000   O     (4c)
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   0.28296000000000   -0.12240000000000    0.71004000000000   O     (8d)
  -0.21704000000000    0.62240000000000   -0.21004000000000   O     (8d)
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   0.21704000000000    0.37760000000000    0.21004000000000   O     (8d)
   0.28296000000000    0.62240000000000    0.71004000000000   O     (8d)
   0.06068000000000    0.06612000000000    0.00658000000000   O     (8d)
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  -0.06068000000000    0.56612000000000   -0.00658000000000   O     (8d)
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  -0.06068000000000   -0.06612000000000   -0.00658000000000   O     (8d)
   0.56068000000000    0.06612000000000    0.49342000000000   O     (8d)
   0.06068000000000    0.43388000000000    0.00658000000000   O     (8d)
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   0.73777000000000    0.38595000000000   -0.34938000000000   O     (8d)
  -0.23777000000000   -0.11405000000000   -0.84938000000000   O     (8d)
   0.73777000000000    0.11405000000000   -0.34938000000000   O     (8d)
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   0.43489000000000    0.06457000000000    0.78946000000000   O     (8d)
   0.06511000000000   -0.06457000000000    1.28946000000000   O     (8d)
  -0.43489000000000    0.56457000000000   -0.78946000000000   O     (8d)
   0.93489000000000    0.43543000000000   -0.28946000000000   O     (8d)
  -0.43489000000000   -0.06457000000000   -0.78946000000000   O     (8d)
   0.93489000000000    0.06457000000000   -0.28946000000000   O     (8d)
   0.43489000000000    0.43543000000000    0.78946000000000   O     (8d)
   0.06511000000000    0.56457000000000    1.28946000000000   O     (8d)
   0.30497000000000    0.12340000000000    0.52074000000000   O     (8d)
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  -0.30497000000000   -0.12340000000000   -0.52074000000000   O     (8d)
   0.80497000000000    0.12340000000000   -0.02074000000000   O     (8d)
   0.30497000000000    0.37660000000000    0.52074000000000   O     (8d)
   0.19503000000000    0.62340000000000    1.02074000000000   O     (8d)
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   0.08448000000000   -0.06300000000000    0.72624000000000   O     (8d)
  -0.41552000000000    0.56300000000000   -0.22624000000000   O     (8d)
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  -0.41552000000000   -0.06300000000000   -0.22624000000000   O     (8d)
   0.91552000000000    0.06300000000000    0.27376000000000   O     (8d)
   0.41552000000000    0.43700000000000    0.22624000000000   O     (8d)
   0.08448000000000    0.56300000000000    0.72624000000000   O     (8d)
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   0.34092000000000   -0.13780000000000    0.51543000000000   Si    (8d)
  -0.15908000000000    0.63780000000000   -0.01543000000000   Si    (8d)
   0.65908000000000    0.36220000000000    0.48457000000000   Si    (8d)
  -0.15908000000000   -0.13780000000000   -0.01543000000000   Si    (8d)
   0.65908000000000    0.13780000000000    0.48457000000000   Si    (8d)
   0.15908000000000    0.36220000000000    0.01543000000000   Si    (8d)
   0.34092000000000    0.63780000000000    0.51543000000000   Si    (8d)
   0.34260000000000    0.13635000000000    0.73115000000000   Si    (8d)
   0.15740000000000   -0.13635000000000    1.23115000000000   Si    (8d)
  -0.34260000000000    0.63635000000000   -0.73115000000000   Si    (8d)
   0.84260000000000    0.36365000000000   -0.23115000000000   Si    (8d)
  -0.34260000000000   -0.13635000000000   -0.73115000000000   Si    (8d)
   0.84260000000000    0.13635000000000   -0.23115000000000   Si    (8d)
   0.34260000000000    0.36365000000000    0.73115000000000   Si    (8d)
   0.15740000000000    0.63635000000000    1.23115000000000   Si    (8d)
   0.33173000000000    0.13741000000000    0.30525000000000   Si    (8d)
   0.16827000000000   -0.13741000000000    0.80525000000000   Si    (8d)
  -0.33173000000000    0.63741000000000   -0.30525000000000   Si    (8d)
   0.83173000000000    0.36259000000000    0.19475000000000   Si    (8d)
  -0.33173000000000   -0.13741000000000   -0.30525000000000   Si    (8d)
   0.83173000000000    0.13741000000000    0.19475000000000   Si    (8d)
   0.33173000000000    0.36259000000000    0.30525000000000   Si    (8d)
   0.16827000000000    0.63741000000000    0.80525000000000   Si    (8d)
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NH_4ClBrI ($F5_{14}$): ABCD_oP16_62_c_c_c_c - CIF

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 'R. C. L. Mooney'
_journal_name_full_name
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 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 98
_journal_year 1938
_journal_page_first 324
_journal_page_last 333
_publ_Section_title
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 The Crystal Structure of Ammonium Chlorobromoiodide and the
      ↪ Configuration of the Chlorobromoiodide Group
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# Found in Strukturbericht Band V 1937, 1940

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_aflow_Strukturbericht '$F5_{14}$'
_aflow_Pearson 'oP16'

_symmetry_space_group_name_H-M "P 21/n 21/m 21/a"
_symmetry_Int_Tables_number 62

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_cell_length_c     8.50000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

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_space_group_symop_operation_xyz
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2 x+1/2,-y+1/2,-z+1/2
3 -x,y+1/2,-z
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x,-y+1/2,z
8 x+1/2,y,-z+1/2

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_atom_site_type_symbol
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_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Br1 Br 4 c 0.57600 0.25000 0.74800 1.00000
Cl1 Cl 4 c 0.19200 0.25000 0.38100 1.00000
I1 I 4 c 0.37600 0.25000 0.56100 1.00000
NH41 NH4 4 c 0.37200 0.25000 0.07400 1.00000
```

NH_4ClBrI ($F5_{14}$): ABCD_oP16_62_c_c_c_c - POSCAR

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      ↪ 0.561,0.372,0.074 & Pnma D_{2h}^{16} #62 (c^4) & oP16 & $F5_{14
      ↪ }$ & ClBrI(NH4) & ClBrI(NH4) & R. C. L. Mooney, Zeitschrift f{"
      ↪ u}r Kristallographie - Crystalline Materials 98, 324-333 (1938)
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   9.94000000000000    0.00000000000000    0.00000000000000
   0.00000000000000    6.13000000000000    0.00000000000000
   0.00000000000000    0.00000000000000    8.50000000000000
    Br    Cl     I    NH4
     4     4     4     4
Direct
   0.57600000000000    0.25000000000000    0.74800000000000   Br    (4c)
  -0.07600000000000    0.75000000000000    1.24800000000000   Br    (4c)
  -0.57600000000000    0.75000000000000   -0.74800000000000   Br    (4c)
   1.07600000000000    0.25000000000000   -0.24800000000000   Br    (4c)
   0.19200000000000    0.25000000000000    0.38100000000000   Cl    (4c)
   0.30800000000000    0.75000000000000    0.88100000000000   Cl    (4c)
  -0.19200000000000    0.75000000000000   -0.38100000000000   Cl    (4c)
   0.69200000000000    0.25000000000000    0.11900000000000   Cl    (4c)
   0.37600000000000    0.25000000000000    0.56100000000000    I    (4c)
   0.12400000000000    0.75000000000000    1.06100000000000    I    (4c)
  -0.37600000000000    0.75000000000000   -0.56100000000000    I    (4c)
   0.87600000000000    0.25000000000000   -0.06100000000000    I    (4c)
   0.37200000000000    0.25000000000000    0.07400000000000   NH4   (4c)
   0.12800000000000    0.75000000000000    0.57400000000000   NH4   (4c)
  -0.37200000000000    0.75000000000000   -0.07400000000000   NH4   (4c)
   0.87200000000000    0.25000000000000    0.42600000000000   NH4   (4c)
```

MnCuP: ABC_oP12_62_c_c_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'CuMnP'
_chemical_formula_sum 'Cu Mn P'

loop_
_publ_author_name
 'J. M\"{u}ndelein'
 'H.-U. Schuster'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Naturforschung B
;
_journal_volume 47
_journal_year 1992
_journal_page_first 925
_journal_page_last 928
_publ_Section_title
;
 Darstellung und Kristallstruktur der Verbindungen MnCu$X$ ($X$ = P, As,
      ↪ P$_{x}$As$_{1-x}$)
;

# Found in Room-temperature antiferromagnetism in CuMnAs, 2012

_aflow_title 'MnCuP Structure'
_aflow_proto 'ABC_oP12_62_c_c_c'
_aflow_params 'a,b/a,c/a,x_{1},z_{1},x_{2},z_{2},x_{3},z_{3}'
_aflow_params_values '6.3187,0.58934590976,1.12178137908,0.6282,0.061,
      ↪ 0.5376,0.6709,0.2492,0.1245'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oP12'

_symmetry_space_group_name_H-M "P 21/n 21/m 21/a"
_symmetry_Int_Tables_number 62

_cell_length_a     6.31870
_cell_length_b     3.72390
_cell_length_c     7.08820
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x,y+1/2,-z
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x,-y+1/2,z
8 x+1/2,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cu1 Cu 4 c 0.62820 0.25000 0.06100 1.00000
Mn1 Mn 4 c 0.53760 0.25000 0.67090 1.00000
P1 P 4 c 0.24920 0.25000 0.12450 1.00000
```

MnCuP: ABC_oP12_62_c_c_c - POSCAR

```
ABC_oP12_62_c_c_c & a,b/a,c/a,x1,z1,x2,z2,x3,z3 --params=6.3187,
      ↪ 0.58934590976,1.12178137908,0.6282,0.061,0.5376,0.6709,0.2492,
      ↪ 0.1245 & Pnma D_{2h}^{16} #62 (c^3) & oP12 & None & CuMnP &
      ↪ CuMnP & J. M\"{u}ndelein and H.-U. Schuster, Z. Naturforsch. B
      ↪ 47, 925-928 (1992)
   1.00000000000000
   6.31870000000000    0.00000000000000    0.00000000000000
   0.00000000000000    3.72390000000000    0.00000000000000
   0.00000000000000    0.00000000000000    7.08820000000000
    Cu    Mn     P
     4     4     4
Direct
   0.62820000000000    0.25000000000000    0.06100000000000   Cu    (4c)
  -0.12820000000000    0.75000000000000    0.56100000000000   Cu    (4c)
  -0.62820000000000    0.75000000000000   -0.06100000000000   Cu    (4c)
   1.12820000000000    0.25000000000000    0.43900000000000   Cu    (4c)
   0.53760000000000    0.25000000000000    0.67090000000000   Mn    (4c)
  -0.03760000000000    0.75000000000000    1.17090000000000   Mn    (4c)
  -0.53760000000000    0.75000000000000   -0.67090000000000   Mn    (4c)
   1.03760000000000    0.25000000000000   -0.17090000000000   Mn    (4c)
   0.24920000000000    0.25000000000000    0.12450000000000    P    (4c)
   0.25080000000000    0.75000000000000    0.62450000000000    P    (4c)
  -0.24920000000000    0.75000000000000   -0.12450000000000    P    (4c)
   0.74920000000000    0.25000000000000    0.37550000000000    P    (4c)
```

η-NiSi (B_d): AB_oP8_62_c_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral '$\eta$-nisi'
_chemical_formula_sum 'Ni Si'

loop_
_publ_author_name
 'K. Toman'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 4
_journal_year 1951
_journal_page_first 462
_journal_page_last 464
_publ_Section_title
;
 The structure of NiSi
;

_aflow_title '$\eta$-NiSi ($B_{d}$) Structure'
_aflow_proto 'AB_oP8_62_c_c'
_aflow_params 'a,b/a,c/a,x_{1},z_{1},x_{2},z_{2}'
```

```
_aflow_params_values '5.18,0.644787644788,1.08494208494,0.006,0.184,0.67
    ↪ ,-0.08'
_aflow_Strukturbericht '$B_{d}$'
_aflow_Pearson 'oP8'

_symmetry_space_group_name_H-M "P 21/n 21/m 21/a"
_symmetry_Int_Tables_number 62

_cell_length_a 5.18000
_cell_length_b 3.34000
_cell_length_c 5.62000
_cell_angle_alpha 90.00000
_cell_angle_beta 90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x,y+1/2,-z
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x+1/2,y+1/2,z+1/2
7 x,-y+1/2,z
8 x+1/2,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ni1 Ni 4 c 0.00600 0.25000 0.18400 1.00000
Si1 Si 4 c 0.67000 0.25000 -0.08000 1.00000
```

η-NiSi (B_d): AB_oP8_62_c_c - POSCAR

```
AB_oP8_62_c_c & a,b/a,c/a,x1,z1,x2,z2 --params=5.18,0.644787644788,
    ↪ 1.08494208494,0.006,0.184,0.67,-0.08 & Pnma D_{2h}^{16} #62 (c^
    ↪ 2) & oP8 & $B_{d}$ & NiSi & $\eta$-nisi & K. Toman, Acta Cryst.
    ↪ 4, 462-464 (1951)
   1.00000000000000
   5.18000000000000   0.00000000000000   0.00000000000000
   0.00000000000000   3.34000000000000   0.00000000000000
   0.00000000000000   0.00000000000000   5.62000000000000
    Ni  Si
     4   4
Direct
   0.00600000000000   0.25000000000000   0.18400000000000  Ni  (4c)
   0.49400000000000   0.75000000000000   0.68400000000000  Ni  (4c)
  -0.00600000000000   0.75000000000000  -0.18400000000000  Ni  (4c)
   0.50600000000000   0.25000000000000   0.31600000000000  Ni  (4c)
   0.67000000000000   0.25000000000000  -0.08000000000000  Si  (4c)
  -0.17000000000000   0.75000000000000   0.42000000000000  Si  (4c)
  -0.67000000000000   0.75000000000000   0.08000000000000  Si  (4c)
   1.17000000000000   0.25000000000000   0.58000000000000  Si  (4c)
```

$Cu_2Pb(SeO_3)_2Br_2$: A2B2C6DE2_oC52_63_g_e_fh_c_f - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Br2Cu2O6PbSe2'
_chemical_formula_sum 'Br2 Cu2 O6 Pb Se2'

loop_
_publ_author_name
 'O. I. Siidra'
 'M. S. Kozin'
 'W. Depmeier'
 'R. A. Kayukov'
 'V. M. Kovrugin'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 74
_journal_year 2018
_journal_page_first 712
_journal_page_last 724
_publ_Section_title
;
 Copper-lead selenite bromides: a new large family of compounds partly
    ↪ having Cu$^{2+}$ substructures derivable from kagome nets
;

_aflow_title 'Cu$_{2}$Pb(SeO$_{3}$)$_{2}$Br$_{2}$ Structure'
_aflow_proto 'A2B2C6DE2_oC52_63_g_e_fh_c_f'
_aflow_params 'a,b/a,c/a,y_{1},x_{2},y_{3},z_{3},y_{4},z_{4},x_{5},y_{5
    ↪ },x_{6},y_{6},z_{6}'
_aflow_params_values '8.275,1.12422960725,1.43697885196,0.33496,0.1862,
    ↪ 0.1287,0.5177,0.31221,0.54036,0.22773,0.07717,0.1531,0.3602,
    ↪ 0.4531'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oC52'

_symmetry_space_group_name_H-M "C 2/m 2/c 21/m"
_symmetry_Int_Tables_number 63

_cell_length_a 8.27500
_cell_length_b 9.30300
_cell_length_c 11.89100
_cell_angle_alpha 90.00000
_cell_angle_beta 90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z+1/2
4 -x,-y,z+1/2
5 -x,-y,-z
6 -x,y,z
7 x,-y,z+1/2
8 x,y,-z+1/2
9 x+1/2,y+1/2,z
10 x+1/2,-y+1/2,-z
11 -x+1/2,y+1/2,-z+1/2
12 -x+1/2,-y+1/2,z+1/2
13 -x+1/2,-y+1/2,-z
14 -x+1/2,y+1/2,z
15 x+1/2,-y+1/2,z+1/2
16 x+1/2,y+1/2,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Pb1 Pb 4 c 0.00000 0.33496 0.25000 1.00000
Cu1 Cu 8 e 0.18620 0.00000 0.00000 1.00000
O1 O 8 f 0.00000 0.12870 0.51770 1.00000
Se1 Se 8 f 0.00000 0.31221 0.54036 1.00000
Br1 Br 8 g 0.22773 0.07717 0.25000 1.00000
O2 O 16 h 0.15310 0.36020 0.45310 1.00000
```

$Cu_2Pb(SeO_3)_2Br_2$: A2B2C6DE2_oC52_63_g_e_fh_c_f - POSCAR

```
A2B2C6DE2_oC52_63_g_e_fh_c_f & a,b/a,c/a,y1,x2,y3,z3,y4,z4,x5,y5,x6,y6,
    ↪ z6 --params=8.275,1.12422960725,1.43697885196,0.33496,0.1862,
    ↪ 0.1287,0.5177,0.31221,0.54036,0.22773,0.07717,0.1531,0.3602,
    ↪ 0.4531 & Cmcm D_{2h}^{17} #63 (cef^2gh) & oC52 & None &
    ↪ Br2Cu2O6PbSe2 & Br2Cu2O6PbSe2 & O. I. Siidra et al., Acta
    ↪ Crystallogr. Sect. B Struct. Sci. 74, 712-724 (2018)
   1.00000000000000
   4.13750000000000  -4.65150000000000   0.00000000000000
   4.13750000000000   4.65150000000000   0.00000000000000
   0.00000000000000   0.00000000000000  11.89100000000000
    Br  Cu  O  Pb  Se
     4   4 12   2   4
Direct
   0.15056000000000   0.30490000000000   0.25000000000000  Br  (8g)
  -0.15056000000000  -0.30490000000000   0.75000000000000  Br  (8g)
  -0.30490000000000  -0.15056000000000   0.25000000000000  Br  (8g)
   0.30490000000000   0.15056000000000   0.75000000000000  Br  (8g)
   0.18620000000000   0.18620000000000   0.00000000000000  Cu  (8e)
  -0.18620000000000  -0.18620000000000   0.50000000000000  Cu  (8e)
  -0.18620000000000  -0.18620000000000   0.00000000000000  Cu  (8e)
   0.18620000000000   0.18620000000000   0.50000000000000  Cu  (8e)
  -0.12870000000000   0.12870000000000   0.51770000000000  O  (8f)
   0.12870000000000  -0.12870000000000   1.01770000000000  O  (8f)
  -0.12870000000000   0.12870000000000  -0.01770000000000  O  (8f)
   0.12870000000000  -0.12870000000000  -0.51770000000000  O  (8f)
  -0.20710000000000   0.51330000000000   0.45310000000000  O  (16h)
   0.20710000000000  -0.51330000000000   0.95310000000000  O  (16h)
  -0.51330000000000   0.20710000000000   0.04690000000000  O  (16h)
   0.51330000000000  -0.20710000000000  -0.45310000000000  O  (16h)
   0.20710000000000  -0.51330000000000  -0.45310000000000  O  (16h)
  -0.20710000000000   0.51330000000000   0.04690000000000  O  (16h)
   0.51330000000000  -0.20710000000000   0.95310000000000  O  (16h)
  -0.51330000000000   0.20710000000000   0.45310000000000  O  (16h)
  -0.33496000000000   0.33496000000000   0.25000000000000  Pb  (4c)
   0.33496000000000  -0.33496000000000   0.75000000000000  Pb  (4c)
  -0.31221000000000   0.31221000000000   0.54036000000000  Se  (8f)
   0.31221000000000  -0.31221000000000   1.04036000000000  Se  (8f)
  -0.31221000000000   0.31221000000000  -0.04036000000000  Se  (8f)
   0.31221000000000  -0.31221000000000  -0.54036000000000  Se  (8f)
```

Pseudobrookite (Fe_2TiO_5, $E4_1$): A2B5C_oC32_63_f_c2f_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Pseudobrookite'
_chemical_formula_sum 'Fe2 O5 Ti'

loop_
_publ_author_name
 'W. Q. Guo'
 'S. Malus'
 'D. H. Ryan'
 'Z. Altounian'
_journal_name_full_name
;
 Journal of Physics: Condensed Matter
;
_journal_volume 11
_journal_year 1999
```

```
_journal_page_first 6337
_journal_page_last 6346
_publ_Section_title
;
 Crystal structure and cation distributions in the FeTi$_{2}$O$_{5}
      ↪ $-Fe$_{2}$TiO$_{5}$ solid solution series
;

_aflow_title 'Pseudobrookite (Fe$_{2}$TiO$_{5}$, $E4_{1}$) Structure'
_aflow_proto 'A2B5C_oC32_63_f_c2f_c'
_aflow_params 'a,b/a,c/a,y_{1},y_{2},y_{3},z_{3},y_{4},z_{4},y_{5},z_{5
      ↪ }'
_aflow_params_values '3.7318,2.62428318774,2.67393751005,0.7617,0.18848,
      ↪ 0.1367,0.5649,0.048,0.1167,0.3108,0.0709'
_aflow_Strukturbericht '$E4_{1}$'
_aflow_Pearson 'oC32'

_symmetry_space_group_name_H-M "C 2/m 2/c 21/m"
_symmetry_Int_Tables_number 63

_cell_length_a    3.73180
_cell_length_b    9.79330
_cell_length_c    9.97860
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z+1/2
4 -x,-y,z+1/2
5 -x,-y,-z
6 -x,y,z
7 x,-y,z+1/2
8 x,y,-z+1/2
9 x+1/2,y+1/2,z
10 x+1/2,-y+1/2,-z
11 -x+1/2,y+1/2,-z+1/2
12 -x+1/2,-y+1/2,z+1/2
13 -x+1/2,-y+1/2,-z
14 -x+1/2,y+1/2,z
15 x+1/2,-y+1/2,z+1/2
16 x+1/2,y+1/2,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 4 c 0.00000 0.76170 0.25000 1.00000
Ti1 Ti 4 c 0.00000 0.18848 0.25000 1.00000
Fe1 Fe 8 f 0.00000 0.13670 0.56490 1.00000
O2 O 8 f 0.00000 0.04800 0.11670 1.00000
O3 O 8 f 0.00000 0.31080 0.07090 1.00000
```

Pseudobrookite (Fe_2TiO_5, $E4_1$): A2B5C_oC32_63_f_c2f_c - POSCAR

```
A2B5C_oC32_63_f_c2f_c & a,b/a,c/a,y1,y2,y3,z3,y4,z4,y5,z5 --params=
      ↪ 3.7318,2.62428318774,2.67393751005,0.7617,0.18848,0.1367,0.5649
      ↪ ,0.048,0.1167,0.3108,0.0709 & Cmcm D_{2h}^{17} #63 (c^2f^3) &
      ↪ oC32 & $E4_{1}$ & Fe2O5Ti & Pseudobrookite & W. Q. Guo et al.,
      ↪ J. Phys.: Condens. Matter 11, 6337-6346 (1999)
   1.00000000000000
   1.86590000000000  -4.89665000000000   0.00000000000000
   1.86590000000000   4.89665000000000   0.00000000000000
   0.00000000000000   0.00000000000000   9.97860000000000
    Fe    O    Ti
     4   10    2
Direct
  -0.13670000000000   0.13670000000000   0.56490000000000  Fe   (8f)
   0.13670000000000  -0.13670000000000   1.06490000000000  Fe   (8f)
  -0.13670000000000   0.13670000000000  -0.06490000000000  Fe   (8f)
   0.13670000000000  -0.13670000000000  -0.56490000000000  Fe   (8f)
  -0.76170000000000   0.76170000000000   0.25000000000000  O    (4c)
   0.76170000000000  -0.76170000000000   0.75000000000000  O    (4c)
  -0.04800000000000   0.04800000000000   0.11670000000000  O    (8f)
   0.04800000000000  -0.04800000000000   0.61670000000000  O    (8f)
  -0.04800000000000   0.04800000000000   0.38330000000000  O    (8f)
   0.04800000000000  -0.04800000000000  -0.11670000000000  O    (8f)
  -0.31080000000000   0.31080000000000   0.07090000000000  O    (8f)
   0.31080000000000  -0.31080000000000   0.57090000000000  O    (8f)
  -0.31080000000000   0.31080000000000   0.42910000000000  O    (8f)
   0.31080000000000  -0.31080000000000  -0.07090000000000  O    (8f)
  -0.18848000000000   0.18848000000000   0.25000000000000  Ti   (4c)
   0.18848000000000  -0.18848000000000   0.75000000000000  Ti   (4c)
```

$MgCuAl_2$ ($E1_a$): A2BC_oC16_63_f_c_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Al2CuMg'
_chemical_formula_sum 'Al2 Cu Mg'

loop_
_publ_author_name
 'B. Heying'
 'R.-D. Hoffmann'
 'R. P\"{o}ttgen'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Naturforschung B
;
_journal_volume 60
_journal_year 2005
_journal_page_first 491
_journal_page_last 494
_publ_Section_title
;
 Structure Refinement of the S-Phase Precipitate MgCuAl$_{2}$
;

_aflow_title 'MgCuAl$_{2}$ ($E1_{a}$) Structure'
_aflow_proto 'A2BC_oC16_63_f_c_c'
_aflow_params 'a,b/a,c/a,y_{1},y_{2},y_{3},z_{3}'
_aflow_params_values '4.0119,2.3093795957,1.77571724121,0.7801,0.0651,
      ↪ 0.3558,0.0556'
_aflow_Strukturbericht '$E1_{a}$'
_aflow_Pearson 'oC16'

_symmetry_space_group_name_H-M "C 2/m 2/c 21/m"
_symmetry_Int_Tables_number 63

_cell_length_a    4.01190
_cell_length_b    9.26500
_cell_length_c    7.12400
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z+1/2
4 -x,-y,z+1/2
5 -x,-y,-z
6 -x,y,z
7 x,-y,z+1/2
8 x,y,-z+1/2
9 x+1/2,y+1/2,z
10 x+1/2,-y+1/2,-z
11 -x+1/2,y+1/2,-z+1/2
12 -x+1/2,-y+1/2,z+1/2
13 -x+1/2,-y+1/2,-z
14 -x+1/2,y+1/2,z
15 x+1/2,-y+1/2,z+1/2
16 x+1/2,y+1/2,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cu1 Cu 4 c 0.00000 0.78010 0.25000 1.00000
Mg1 Mg 4 c 0.00000 0.06510 0.25000 1.00000
Al1 Al 8 f 0.00000 0.35580 0.05560 1.00000
```

$MgCuAl_2$ ($E1_a$): A2BC_oC16_63_f_c_c - POSCAR

```
A2BC_oC16_63_f_c_c & a,b/a,c/a,y1,y2,y3,z3 --params=4.0119,2.3093795957,
      ↪ 1.77571724121,0.7801,0.0651,0.3558,0.0556 & Cmcm D_{2h}^{17} #
      ↪ 63 (c^2f) & oC16 & $E1_{a}$ & Al2CuMg & Al2CuMg & B. Heying and
      ↪ R.-D. Hoffmann and R. P\"{o}ttgen, Z. Naturforsch. B 60,
      ↪ 491-494 (2005)
   1.00000000000000
   2.00595000000000  -4.63250000000000   0.00000000000000
   2.00595000000000   4.63250000000000   0.00000000000000
   0.00000000000000   0.00000000000000   7.12400000000000
    Al    Cu    Mg
     4     2     2
Direct
  -0.35580000000000   0.35580000000000   0.05560000000000  Al   (8f)
   0.35580000000000  -0.35580000000000   0.55560000000000  Al   (8f)
  -0.35580000000000   0.35580000000000   0.44440000000000  Al   (8f)
   0.35580000000000  -0.35580000000000  -0.05560000000000  Al   (8f)
  -0.78010000000000   0.78010000000000   0.25000000000000  Cu   (4c)
   0.78010000000000  -0.78010000000000   0.75000000000000  Cu   (4c)
  -0.06510000000000   0.06510000000000   0.25000000000000  Mg   (4c)
   0.06510000000000  -0.06510000000000   0.75000000000000  Mg   (4c)
```

$S0_4$ (Staurolite, $Fe(OH)_2Al_4Si_2O_{10}$) (*obsolete*): A4BC12D2_oC76_63_eg_c_f3gh_g - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Staurolite'
_chemical_formula_sum 'Al4 Fe O12 Si2'

loop_
_publ_author_name
 'C. Hermann'
 'O. Lohrmann'
 'H. Philipp'
_journal_year 1937
_publ_Section_title
```

```
;
 Strukturbericht Band II 1928-1932
;

_aflow_title '$S0_{4}$ (Staurolite, Fe(OH)$_{2}$Al$_{4}$Si$_{2}$O$_{10}$
      ↪ ) ({\em{obsolete}}) Structure'
_aflow_proto 'A4BC12D2_oC76_63_eg_c_f3gh_g'
_aflow_params 'a,b/a,c/a,y_{1},x_{2},y_{3},z_{3},x_{4},y_{4},x_{5},y_{5
      ↪ },x_{6},y_{6},x_{7},y_{7},x_{8},y_{8},x_{9},y_{9},z_{9}'
_aflow_params_values '16.52,0.473365617433,0.340799031477,0.611,0.328,
      ↪ 0.778,0.0,-0.092,0.25,0.094,0.0,0.25,0.0,0.406,0.0,0.667,0.375,
      ↪ 0.658,0.173,0.0'
_aflow_Strukturbericht '$S0_{4}$'
_aflow_Pearson 'oC76'

_symmetry_space_group_name_H-M "C 2/m 2/c 21/m"
_symmetry_Int_Tables_number 63

_cell_length_a     16.52000
_cell_length_b     7.82000
_cell_length_c     5.63000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z+1/2
4 -x,-y,z+1/2
5 -x,-y,-z
6 -x,y,z
7 x,-y,z+1/2
8 x,y,-z+1/2
9 x+1/2,y+1/2,z
10 x+1/2,-y+1/2,-z
11 -x+1/2,y+1/2,-z+1/2
12 -x+1/2,-y+1/2,z+1/2
13 -x+1/2,-y+1/2,-z
14 -x+1/2,y+1/2,z
15 x+1/2,-y+1/2,z+1/2
16 x+1/2,y+1/2,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Fe1 Fe 4 c 0.00000 0.61100 0.25000 1.00000
Al1 Al 8 e 0.32800 0.00000 0.00000 1.00000
O1 O 8 f 0.00000 0.77800 0.00000 1.00000
Al2 Al 8 g -0.09200 0.25000 0.25000 1.00000
O2 O 8 g 0.09400 0.00000 0.25000 1.00000
O3 O 8 g 0.25000 0.00000 0.25000 1.00000
O4 O 8 g 0.40600 0.00000 0.25000 1.00000
Si1 Si 8 g 0.66700 0.37500 0.25000 1.00000
O5 O 16 h 0.65800 0.17300 0.00000 1.00000
```

$S0_4$ (Staurolite, $Fe(OH)_2Al_4Si_2O_{10}$) (*obsolete*): A4BC12D2_oC76_63_eg_c_f3gh_g - POSCAR

```
A4BC12D2_oC76_63_eg_c_f3gh_g & a,b/a,c/a,y1,x2,y3,z3,x4,y4,x5,y5,x6,y6,
      ↪ x7,y7,x8,y8,x9,y9,z9 --params=16.52,0.473365617433,
      ↪ 0.340799031477,0.611,0.328,0.778,0.0,-0.092,0.25,0.094,0.0,0.25
      ↪ ,0.0,0.406,0.0,0.667,0.375,0.658,0.173,0.0 & Cmcm D_{2h}^{17} #
      ↪ 63 (cefg^5h) & oC76 & $S0_{4}$ & Al4FeO12Si2 & Staurolite & C.
      ↪ Hermann and O. Lohrmann and H. Philipp, (1937)
   1.00000000000000
   8.26000000000000  -3.91000000000000   0.00000000000000
   8.26000000000000   3.91000000000000   0.00000000000000
   0.00000000000000   0.00000000000000   5.63000000000000
     Al     Fe     O     Si
      8      2     24     4
Direct
   0.32800000000000   0.32800000000000   0.00000000000000   Al   (8e)
  -0.32800000000000  -0.32800000000000   0.50000000000000   Al   (8e)
  -0.32800000000000  -0.32800000000000   0.00000000000000   Al   (8e)
   0.32800000000000   0.32800000000000   0.50000000000000   Al   (8e)
  -0.34200000000000   0.15800000000000   0.25000000000000   Al   (8g)
   0.34200000000000  -0.15800000000000   0.75000000000000   Al   (8g)
  -0.15800000000000   0.34200000000000   0.25000000000000   Al   (8g)
   0.15800000000000  -0.34200000000000   0.75000000000000   Al   (8g)
  -0.61100000000000   0.61100000000000   0.25000000000000   Fe   (4c)
   0.61100000000000  -0.61100000000000   0.75000000000000   Fe   (4c)
  -0.77800000000000   0.77800000000000   0.00000000000000   O    (8f)
   0.77800000000000  -0.77800000000000   0.50000000000000   O    (8f)
  -0.77800000000000   0.77800000000000   0.50000000000000   O    (8f)
   0.77800000000000  -0.77800000000000   0.00000000000000   O    (8f)
   0.09400000000000   0.09400000000000   0.25000000000000   O    (8g)
  -0.09400000000000  -0.09400000000000   0.75000000000000   O    (8g)
  -0.09400000000000  -0.09400000000000   0.25000000000000   O    (8g)
   0.09400000000000   0.09400000000000   0.75000000000000   O    (8g)
   0.25000000000000   0.25000000000000   0.25000000000000   O    (8g)
  -0.25000000000000  -0.25000000000000   0.75000000000000   O    (8g)
  -0.25000000000000  -0.25000000000000   0.25000000000000   O    (8g)
   0.25000000000000   0.25000000000000   0.75000000000000   O    (8g)
   0.40600000000000   0.40600000000000   0.25000000000000   O    (8g)
  -0.40600000000000  -0.40600000000000   0.75000000000000   O    (8g)
  -0.40600000000000  -0.40600000000000   0.25000000000000   O    (8g)
   0.40600000000000   0.40600000000000   0.75000000000000   O    (8g)
   0.48500000000000   0.83100000000000   0.00000000000000   O    (16h)
  -0.48500000000000  -0.83100000000000   0.50000000000000   O    (16h)
  -0.83100000000000  -0.48500000000000   0.50000000000000   O    (16h)
   0.83100000000000   0.48500000000000   0.00000000000000   O    (16h)
  -0.48500000000000  -0.83100000000000   0.00000000000000   O    (16h)
   0.48500000000000   0.83100000000000   0.50000000000000   O    (16h)
   0.83100000000000   0.48500000000000   0.50000000000000   O    (16h)
  -0.83100000000000  -0.48500000000000   0.00000000000000   O    (16h)
   0.29200000000000   1.04200000000000   0.25000000000000   Si   (8g)
  -0.29200000000000  -1.04200000000000   0.75000000000000   Si   (8g)
  -1.04200000000000  -0.29200000000000   0.25000000000000   Si   (8g)
   1.04200000000000   0.29200000000000   0.75000000000000   Si   (8g)
```

Pd_5Pu_3: A5B3_oC32_63_cfg_ce - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Pd5Pu3'
_chemical_formula_sum 'Pd5 Pu3'

loop_
_publ_author_name
 'D. T. Cromer'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 32
_journal_year 1976
_journal_page_first 1930
_journal_page_last 1932
_publ_Section_title
;
 Plutonium-palladium Pu$_{3}$Pd$_{5}$
;

# Found in New $R$$_{3}$Pd$_{5}$ Compounds ($R$ = Sc, Y, Gd-Lu):
      ↪ Formation and Stability, Crystal Structure, and
      ↪ Antiferromagnetism, 2016

_aflow_title 'Pd$_{5}$Pu$_{3}$ Structure'
_aflow_proto 'A5B3_oC32_63_cfg_ce'
_aflow_params 'a,b/a,c/a,y_{1},y_{2},x_{3},y_{4},z_{4},x_{5},y_{5}'
_aflow_params_values '9.20101,0.781979369656,1.0619486339,0.0254,0.6251,
      ↪ 0.2018,0.3147,0.451,0.2219,0.2863'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oC32'

_symmetry_space_group_name_H-M "C 2/m 2/c 21/m"
_symmetry_Int_Tables_number 63

_cell_length_a     9.20101
_cell_length_b     7.19500
_cell_length_c     9.77100
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z+1/2
4 -x,-y,z+1/2
5 -x,-y,-z
6 -x,y,z
7 x,-y,z+1/2
8 x,y,-z+1/2
9 x+1/2,y+1/2,z
10 x+1/2,-y+1/2,-z
11 -x+1/2,y+1/2,-z+1/2
12 -x+1/2,-y+1/2,z+1/2
13 -x+1/2,-y+1/2,-z
14 -x+1/2,y+1/2,z
15 x+1/2,-y+1/2,z+1/2
16 x+1/2,y+1/2,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Pd1 Pd 4 c 0.00000 0.02540 0.25000 1.00000
Pu1 Pu 4 c 0.00000 0.62510 0.25000 1.00000
Pu2 Pu 8 e 0.20180 0.00000 0.00000 1.00000
Pd2 Pd 8 f 0.00000 0.31470 0.45100 1.00000
Pd3 Pd 8 g 0.22190 0.28630 0.25000 1.00000
```

Pd_5Pu_3: A5B3_oC32_63_cfg_ce - POSCAR

```
A5B3_oC32_63_cfg_ce & a,b/a,c/a,y1,y2,x3,y4,z4,x5,y5 --params=9.20101,
      ↪ 0.781979369656,1.0619486339,0.0254,0.6251,0.2018,0.3147,0.451,
      ↪ 0.2219,0.2863 & Cmcm D_{2h}^{17} #63 (c^2efg) & oC32 & None &
      ↪ Pd5Pu3 & Pd5Pu3 & D. T. Cromer, Acta Crystallogr. Sect. B
      ↪ Struct. Sci. 32, 1930-1932 (1976)
   1.00000000000000
   4.60050500000000  -3.59750000000000   0.00000000000000
   4.60050500000000   3.59750000000000   0.00000000000000
```

```
    0.00000000000000    0.00000000000000    9.77100000000000
     Pd      Pu
     10       6
Direct
  -0.02540000000000    0.02540000000000    0.25000000000000    Pd    (4c)
   0.02540000000000   -0.02540000000000    0.75000000000000    Pd    (4c)
  -0.31470000000000    0.31470000000000    0.45100000000000    Pd    (8f)
   0.31470000000000   -0.31470000000000    0.95100000000000    Pd    (8f)
  -0.31470000000000    0.31470000000000    0.04900000000000    Pd    (8f)
   0.31470000000000   -0.31470000000000   -0.45100000000000    Pd    (8f)
  -0.06440000000000    0.50820000000000    0.25000000000000    Pd    (8g)
   0.06440000000000   -0.50820000000000    0.75000000000000    Pd    (8g)
  -0.50820000000000    0.06440000000000    0.25000000000000    Pd    (8g)
   0.50820000000000   -0.06440000000000    0.75000000000000    Pd    (8g)
  -0.62510000000000    0.62510000000000    0.25000000000000    Pu    (4c)
   0.62510000000000   -0.62510000000000    0.75000000000000    Pu    (4c)
   0.20180000000000    0.20180000000000    0.00000000000000    Pu    (8e)
  -0.20180000000000   -0.20180000000000    0.50000000000000    Pu    (8e)
  -0.20180000000000   -0.20180000000000    0.00000000000000    Pu    (8e)
   0.20180000000000    0.20180000000000    0.50000000000000    Pu    (8e)
```

$ZrTe_5$: A5B_oC24_63_c2f_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Te5Zr'
_chemical_formula_sum 'Te5 Zr'

loop_
_publ_author_name
 'H. Fjellv{\aa}g'
 'A. Kjekshus'
_journal_name_full_name
;
 Solid State Communications
;
_journal_volume 60
_journal_year 1986
_journal_page_first 91
_journal_page_last 93
_publ_Section_title
;
 Structural Properties of ZrTe$_{5}$ and HfTe$_{5}$ as Seen by Power
     ↪ Diffraction
;

_aflow_title 'ZrTe$_{5}$ Structure'
_aflow_proto 'A5B_oC24_63_c2f_c'
_aflow_params 'a,b/a,c/a,y_{1},y_{2},y_{3},z_{3},y_{4},z_{4}'
_aflow_params_values '3.9875,3.64388714734,3.4407523511,0.316,0.633,-
    ↪ 0.067,0.151,0.209,0.434'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oC24'

_symmetry_space_group_name_H-M "C 2/m 2/c 21/m"
_symmetry_Int_Tables_number 63

_cell_length_a    3.98750
_cell_length_b    14.53000
_cell_length_c    13.72000
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z+1/2
4 -x,-y,z+1/2
5 -x,-y,-z
6 -x,y,z
7 x,-y,z+1/2
8 x,y,-z+1/2
9 x+1/2,y+1/2,z
10 x+1/2,-y+1/2,-z
11 -x+1/2,y+1/2,-z+1/2
12 -x+1/2,-y+1/2,z+1/2
13 -x+1/2,-y+1/2,-z
14 -x+1/2,y+1/2,z
15 x+1/2,-y+1/2,z+1/2
16 x+1/2,y+1/2,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Te1 Te 4 c 0.00000 0.31600 0.25000 1.00000
Zr1 Zr 4 c 0.00000 0.63300 0.25000 1.00000
Te2 Te 8 f 0.00000 -0.06700 0.15100 1.00000
Te3 Te 8 f 0.00000 0.20900 0.43400 1.00000
```

$ZrTe_5$: A5B_oC24_63_c2f_c - POSCAR

```
A5B_oC24_63_c2f_c & a,b/a,c/a,y1,y2,y3,z3,y4,z4 --params=3.9875,
    ↪ 3.64388714734,3.4407523511,0.316,0.633,-0.067,0.151,0.209,0.434
    ↪ & Cmcm D_{2h}^{17} #63 (c^2f^2) & oC24 & None & Te5Zr & Te5Zr
    ↪ & H. Fjellv{\aa}g and A. Kjekshus, Solid State Commun. 60,
    ↪ 91-93 (1986)
   1.00000000000000
   1.99375000000000   -7.26500000000000    0.00000000000000
   1.99375000000000    7.26500000000000    0.00000000000000
   0.00000000000000    0.00000000000000   13.72000000000000
    Te      Zr
    10       2
Direct
  -0.31600000000000    0.31600000000000    0.25000000000000    Te    (4c)
   0.31600000000000   -0.31600000000000    0.75000000000000    Te    (4c)
   0.06700000000000   -0.06700000000000    0.15100000000000    Te    (8f)
  -0.06700000000000    0.06700000000000    0.65100000000000    Te    (8f)
   0.06700000000000   -0.06700000000000    0.34900000000000    Te    (8f)
  -0.06700000000000    0.06700000000000   -0.15100000000000    Te    (8f)
  -0.20900000000000    0.20900000000000    0.43400000000000    Te    (8f)
   0.20900000000000   -0.20900000000000    0.93400000000000    Te    (8f)
  -0.20900000000000    0.20900000000000    0.06600000000000    Te    (8f)
   0.20900000000000   -0.20900000000000   -0.43400000000000    Te    (8f)
  -0.63300000000000    0.63300000000000    0.25000000000000    Zr    (4c)
   0.63300000000000   -0.63300000000000    0.75000000000000    Zr    (4c)
```

Lepidocrocite (γ-FeO(OH), $E0_4$): AB2C2_oC20_63_c_f_2c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Lepidocrocite'
_chemical_formula_sum 'Fe H2 O2'

loop_
_publ_author_name
 'A. {N{\o}rlund Christensen}'
 'M. S. Lehmann'
 'P. Convert'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 36a
_journal_year 1982
_journal_page_first 303
_journal_page_last 308
_publ_Section_title
;
 Deuteration of Crystalline Hydroxides, Hydrogen Bonds of $\gamma$-AlOO(
     ↪ H,D) and $\gamma$-FeOO(H,D)
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'Lepidocrocite ($\gamma$-FeO(OH), $E0_{4}$) Structure'
_aflow_proto 'AB2C2_oC20_63_c_f_2c'
_aflow_params 'a,b/a,c/a,y_{1},y_{2},y_{3},y_{4},z_{4}'
_aflow_params_values '3.07,4.08143322476,1.22475570033,-0.3137,0.2842,
    ↪ 0.0724,0.0143,0.3663'
_aflow_Strukturbericht '$E0_{4}$'
_aflow_Pearson 'oC20'

_symmetry_space_group_name_H-M "C 2/m 2/c 21/m"
_symmetry_Int_Tables_number 63

_cell_length_a    3.07000
_cell_length_b    12.53000
_cell_length_c    3.76000
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z+1/2
4 -x,-y,z+1/2
5 -x,-y,-z
6 -x,y,z
7 x,-y,z+1/2
8 x,y,-z+1/2
9 x+1/2,y+1/2,z
10 x+1/2,-y+1/2,-z
11 -x+1/2,y+1/2,-z+1/2
12 -x+1/2,-y+1/2,z+1/2
13 -x+1/2,-y+1/2,-z
14 -x+1/2,y+1/2,z
15 x+1/2,-y+1/2,z+1/2
16 x+1/2,y+1/2,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Fe1 Fe 4 c 0.00000 -0.31370 0.25000 1.00000
O1 O 4 c 0.00000 0.28420 0.25000 1.00000
O2 O 4 c 0.00000 0.07240 0.25000 1.00000
H1 H 8 f 0.00000 0.01430 0.36630 0.47100
```

Lepidocrocite (γ-FeO(OH), $E0_4$): AB2C2_oC20_63_c_f_2c - POSCAR

```
AB2C2_oC20_63_c_f_2c & a,b/a,c/a,y1,y2,y3,y4,z4 --params=3.07,
    ↪ 4.08143322476,1.22475570033,-0.3137,0.2842,0.0724,0.0143,0.3663
    ↪ & Cmcm D_{2h}^{17} #63 (c^3f) & oC20 & $E0_{4}$ & FeHO2 &
    ↪ Lepidocrocite & A. {N{\o}rlund Christensen} and M. S. Lehmann
    ↪ and P. Convert, Acta Chem. Scand. 36a, 303-308 (1982)
   1.00000000000000
   1.53500000000000  -6.26500000000000   0.00000000000000
   1.53500000000000   6.26500000000000   0.00000000000000
   0.00000000000000   0.00000000000000   3.76000000000000
    Fe    H    O
     2    4    4
Direct
   0.31370000000000  -0.31370000000000   0.25000000000000  Fe   (4c)
  -0.31370000000000   0.31370000000000   0.75000000000000  Fe   (4c)
  -0.01430000000000   0.01430000000000   0.36630000000000  H    (8f)
   0.01430000000000  -0.01430000000000   0.86630000000000  H    (8f)
  -0.01430000000000   0.01430000000000   0.13370000000000  H    (8f)
   0.01430000000000  -0.01430000000000  -0.36630000000000  H    (8f)
  -0.28420000000000   0.28420000000000   0.25000000000000  O    (4c)
   0.28420000000000  -0.28420000000000   0.75000000000000  O    (4c)
  -0.07240000000000   0.07240000000000   0.25000000000000  O    (4c)
   0.07240000000000  -0.07240000000000   0.75000000000000  O    (4c)
```

Na_2CrO_4 ($H1_8$): AB2C4_oC28_63_c_bc_fg - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'CrNa2O4'
_chemical_formula_sum 'Cr Na2 O4'

loop_
_publ_author_name
 'A. Niggli'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 7
_journal_year 1954
_journal_page_first 776
_journal_page_last 776
_publ_Section_title
;
 Die Raumgruppe von Na$_{2}$CrO$_{4}$
;

_aflow_title 'Na$_{2}$CrO$_{4}$ ($H1_{8}$) Structure'
_aflow_proto 'AB2C4_oC28_63_c_bc_fg'
_aflow_params 'a,b/a,c/a,y_{2},y_{3},y_{4},z_{4},x_{5},y_{5}'
_aflow_params_values '5.91,1.56175972927,1.21827411168,0.8472,0.2,0.75,
    ↪ 0.0694,0.2778,0.45'
_aflow_Strukturbericht '$H1_{8}$'
_aflow_Pearson 'oC28'

_symmetry_space_group_name_H-M "C 2/m 2/c 21/m"
_symmetry_Int_Tables_number 63

_cell_length_a    5.91000
_cell_length_b    9.23000
_cell_length_c    7.20000
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z+1/2
4 -x,-y,z+1/2
5 -x,-y,-z
6 -x,y,z
7 x,-y,z+1/2
8 x,y,-z+1/2
9 x+1/2,y+1/2,z
10 x+1/2,-y+1/2,-z
11 -x+1/2,y+1/2,-z+1/2
12 -x+1/2,-y+1/2,z+1/2
13 -x+1/2,-y+1/2,-z
14 -x+1/2,y+1/2,z
15 x+1/2,-y+1/2,z+1/2
16 x+1/2,y+1/2,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Na1 Na 4 b 0.00000 0.50000 0.00000 1.00000
Cr1 Cr 4 c 0.00000 0.84720 0.25000 1.00000
Na2 Na 4 c 0.00000 0.20000 0.25000 1.00000
O1 O 8 f 0.00000 0.75000 0.06940 1.00000
O2 O 8 g 0.27780 0.45000 0.25000 1.00000
```

Na_2CrO_4 ($H1_8$): AB2C4_oC28_63_c_bc_fg - POSCAR

```
AB2C4_oC28_63_c_bc_fg & a,b/a,c/a,y2,y3,y4,z4,x5,y5 --params=5.91,
    ↪ 1.56175972927,1.21827411168,0.8472,0.2,0.75,0.0694,0.2778,0.45
    ↪ & Cmcm D_{2h}^{17} #63 (bc^2fg) & oC28 & $H1_{8}$ & CrNa2O4 &
    ↪ CrNa2O4 & A. Niggli, Acta Cryst. 7, 776(1954)
   1.00000000000000
   2.95500000000000  -4.61500000000000   0.00000000000000
   2.95500000000000   4.61500000000000   0.00000000000000
   0.00000000000000   0.00000000000000   7.20000000000000
    Cr    Na    O
     2     4     8
Direct
  -0.84720000000000   0.84720000000000   0.25000000000000  Cr   (4c)
   0.84720000000000  -0.84720000000000   0.75000000000000  Cr   (4c)
   0.50000000000000   0.50000000000000   0.00000000000000  Na   (4b)
   0.50000000000000   0.50000000000000   0.50000000000000  Na   (4b)
  -0.20000000000000   0.20000000000000   0.25000000000000  Na   (4c)
   0.20000000000000  -0.20000000000000   0.75000000000000  Na   (4c)
  -0.75000000000000   0.75000000000000   0.06940000000000  O    (8f)
   0.75000000000000  -0.75000000000000   0.56940000000000  O    (8f)
  -0.75000000000000   0.75000000000000   0.43060000000000  O    (8f)
   0.75000000000000  -0.75000000000000  -0.06940000000000  O    (8f)
  -0.17220000000000   0.72780000000000   0.25000000000000  O    (8g)
   0.17220000000000  -0.72780000000000   0.75000000000000  O    (8g)
  -0.72780000000000   0.17220000000000   0.25000000000000  O    (8g)
   0.72780000000000  -0.17220000000000   0.75000000000000  O    (8g)
```

$ThFe_2SiC$: AB2CD_oC20_63_b_f_c_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'CFe2SiTh'
_chemical_formula_sum 'C Fe2 Si Th'

loop_
_publ_author_name
 'A. M. Witte'
 'W. Jeitschko'
_journal_name_full_name
;
 Journal of Solid State Chemistry
;
_journal_volume 112
_journal_year 1994
_journal_page_first 232
_journal_page_last 236
_publ_Section_title
;
 Carbides with Filled Re$_{3}$B-Type Structure
;

_aflow_title 'ThFe$_{2}$SiC Structure'
_aflow_proto 'AB2CD_oC20_63_b_f_c_c'
_aflow_params 'a,b/a,c/a,y_{2},y_{3},y_{4},z_{4}'
_aflow_params_values '3.863,2.79730779187,1.79911985503,0.7719,0.05275,
    ↪ 0.66585,0.5611'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oC20'

_symmetry_space_group_name_H-M "C 2/m 2/c 21/m"
_symmetry_Int_Tables_number 63

_cell_length_a    3.86300
_cell_length_b    10.80600
_cell_length_c    6.95000
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z+1/2
4 -x,-y,z+1/2
5 -x,-y,-z
6 -x,y,z
7 x,-y,z+1/2
8 x,y,-z+1/2
9 x+1/2,y+1/2,z
10 x+1/2,-y+1/2,-z
11 -x+1/2,y+1/2,-z+1/2
12 -x+1/2,-y+1/2,z+1/2
13 -x+1/2,-y+1/2,-z
14 -x+1/2,y+1/2,z
15 x+1/2,-y+1/2,z+1/2
16 x+1/2,y+1/2,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
C1 C 4 b 0.00000 0.50000 0.00000 1.00000
Si1 Si 4 c 0.00000 0.77190 0.25000 1.00000
Th1 Th 4 c 0.00000 0.05275 0.25000 1.00000
Fe1 Fe 8 f 0.00000 0.66585 0.56110 1.00000
```

$ThFe_2SiC$: AB2CD_oC20_63_b_f_c_c - POSCAR

```
AB2CD_oC20_63_b_f_c_c & a,b/a,c/a,y2,y3,y4,z4 --params=3.863,
    ↪ 2.79730779187,1.79911985503,0.7719,0.05275,0.66585,0.5611 &
    ↪ Cmcm D_{2h}^{17} #63 (bc^2f) & oC20 & None & CFe2SiTh &
    ↪ CFe2SiTh & A. M. Witte and W. Jeitschko, J. Solid State Chem.
    ↪ 112, 232-236 (1994)
  1.00000000000000
  1.93150000000000  -5.40300000000000   0.00000000000000
  1.93150000000000   5.40300000000000   0.00000000000000
  0.00000000000000   0.00000000000000   6.95000000000000
    C    Fe    Si    Th
    2     4     2     2
Direct
   0.50000000000000   0.50000000000000   0.00000000000000   C    (4b)
   0.50000000000000   0.50000000000000   0.50000000000000   C    (4b)
  -0.66585000000000   0.66585000000000   0.56110000000000   Fe   (8f)
   0.66585000000000  -0.66585000000000   1.06110000000000   Fe   (8f)
  -0.66585000000000   0.66585000000000  -0.06110000000000   Fe   (8f)
   0.66585000000000  -0.66585000000000  -0.56110000000000   Fe   (8f)
  -0.77190000000000   0.77190000000000   0.25000000000000   Si   (4c)
   0.77190000000000  -0.77190000000000   0.75000000000000   Si   (4c)
  -0.05275000000000   0.05275000000000   0.25000000000000   Th   (4c)
   0.05275000000000  -0.05275000000000   0.75000000000000   Th   (4c)
```

Mn_3As ($D0_d$): AB3_oC16_63_c_3c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'AsMn3'
_chemical_formula_sum 'As Mn3'

loop_
_publ_author_name
 'W. {Carrillo-Cabrera}'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 37a
_journal_year 1983
_journal_page_first 93
_journal_page_last 98
_publ_Section_title
;
 The Crystal Structure of TiCu$_{2}$P and Its Relationship to the
    ↪ Structure of Mn$_{3}$As
;

_aflow_title 'Mn$_{3}$As ($D0_{d}$) Structure'
_aflow_proto 'AB3_oC16_63_c_3c'
_aflow_params 'a,b/a,c/a,y_{1},y_{2},y_{3},y_{4}'
_aflow_params_values '3.78779,4.28999759754,1.0,0.159,-0.0565,0.5565,
    ↪ 0.3155'
_aflow_Strukturbericht '$D0_{d}$'
_aflow_Pearson 'oC16'

_symmetry_space_group_name_H-M "C 2/m 2/c 21/m"
_symmetry_Int_Tables_number 63

_cell_length_a     3.78779
_cell_length_b     16.24961
_cell_length_c     3.78779
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z+1/2
4 -x,-y,z+1/2
5 -x,-y,-z
6 -x,y,z
7 x,-y,z+1/2
8 x,y,-z+1/2
9 x+1/2,y+1/2,z
10 x+1/2,-y+1/2,-z
11 -x+1/2,y+1/2,-z+1/2
12 -x+1/2,-y+1/2,z+1/2
13 -x+1/2,-y+1/2,-z
14 -x+1/2,y+1/2,z
15 x+1/2,-y+1/2,z+1/2
16 x+1/2,y+1/2,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
As1 As 4 c 0.00000 0.15900 0.25000 1.00000
Mn1 Mn 4 c 0.00000 -0.05650 0.25000 1.00000
Mn2 Mn 4 c 0.00000 0.55650 0.25000 1.00000
Mn3 Mn 4 c 0.00000 0.31550 0.25000 1.00000
```

Mn_3As ($D0_d$): AB3_oC16_63_c_3c - POSCAR

```
AB3_oC16_63_c_3c & a,b/a,c/a,y1,y2,y3,y4 --params=3.78779,4.28999759754,
    ↪ 1.0,0.159,-0.0565,0.5565,0.3155 & Cmcm D_{2h}^{17} #63 (c^4) &
    ↪ oC16 & $D0_{d}$ & AsMn3 & AsMn3 & W. {Carrillo-Cabrera}, Acta
    ↪ Chem. Scand. 37a, 93-98 (1983)
  1.00000000000000
  1.89389500000000  -8.12480500000000   0.00000000000000
  1.89389500000000   8.12480500000000   0.00000000000000
  0.00000000000000   0.00000000000000   3.78779000000000
    As    Mn
     2     6
Direct
  -0.15900000000000   0.15900000000000   0.25000000000000   As   (4c)
   0.15900000000000  -0.15900000000000   0.75000000000000   As   (4c)
   0.05650000000000  -0.05650000000000   0.25000000000000   Mn   (4c)
  -0.05650000000000   0.05650000000000   0.75000000000000   Mn   (4c)
  -0.55650000000000   0.55650000000000   0.25000000000000   Mn   (4c)
   0.55650000000000  -0.55650000000000   0.75000000000000   Mn   (4c)
  -0.31550000000000   0.31550000000000   0.25000000000000   Mn   (4c)
   0.31550000000000  -0.31550000000000   0.75000000000000   Mn   (4c)
```

Re_3B: AB3_oC16_63_c_cf - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'BRe3'
_chemical_formula_sum 'B Re3'

loop_
_publ_author_name
 'B. Aronsson'
 'M. B\"{a}ckman'
 'S. Rundqvist'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 14
_journal_year 1960
_journal_page_first 1001
_journal_page_last 1005
_publ_Section_title
;
 The Crystal Structure of Re$_{3}$B
;

_aflow_title 'Re$_{3}$B Structure'
_aflow_proto 'AB3_oC16_63_c_cf'
_aflow_params 'a,b/a,c/a,y_{1},y_{2},y_{3},z_{3}'
_aflow_params_values '2.89,3.22249134948,2.51141868512,0.744,0.4262,
    ↪ 0.1345,0.062'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oC16'

_symmetry_space_group_name_H-M "C 2/m 2/c 21/m"
_symmetry_Int_Tables_number 63

_cell_length_a     2.89000
_cell_length_b     9.31300
_cell_length_c     7.25800
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z+1/2
4 -x,-y,z+1/2
5 -x,-y,-z
6 -x,y,z
7 x,-y,z+1/2
8 x,y,-z+1/2
9 x+1/2,y+1/2,z
10 x+1/2,-y+1/2,-z
11 -x+1/2,y+1/2,-z+1/2
12 -x+1/2,-y+1/2,z+1/2
13 -x+1/2,-y+1/2,-z
14 -x+1/2,y+1/2,z
15 x+1/2,-y+1/2,z+1/2
16 x+1/2,y+1/2,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
B1 B 4 c 0.00000 0.74400 0.25000 1.00000
Re1 Re 4 c 0.00000 0.42620 0.25000 1.00000
Re2 Re 8 f 0.00000 0.13450 0.06200 1.00000
```

Re_3B: AB3_oC16_63_c_cf - POSCAR

```
AB3_oC16_63_c_cf & a,b/a,c/a,y1,y2,y3,z3 --params=2.89,3.22249134948,
    ↪ 2.51141868512,0.744,0.4262,0.1345,0.062 & Cmcm D_{2h}^{17} #63
    ↪ (c^2f) & oC16 & None & BRe3 & BRe3 & B. Aronsson and M. B\"{a}
    ↪ ckman and S. Rundqvist, Acta Chem. Scand. 14, 1001-1005 (1960)
  1.00000000000000
```

```
    1.44500000000000   -4.65650000000000    0.00000000000000
    1.44500000000000    4.65650000000000    0.00000000000000
    0.00000000000000    0.00000000000000    7.25800000000000
      B     Re
      2      6
Direct
  -0.74400000000000    0.74400000000000    0.25000000000000    B    (4c)
   0.74400000000000   -0.74400000000000    0.75000000000000    B    (4c)
  -0.42620000000000    0.42620000000000    0.25000000000000    Re   (4c)
   0.42620000000000   -0.42620000000000    0.75000000000000    Re   (4c)
  -0.13450000000000    0.13450000000000    0.06200000000000    Re   (8f)
   0.13450000000000   -0.13450000000000    0.56200000000000    Re   (8f)
  -0.13450000000000    0.13450000000000    0.43800000000000    Re   (8f)
   0.13450000000000   -0.13450000000000   -0.06200000000000    Re   (8f)
```

Ta_2NiS_5: AB5C2_oC32_63_c_c2f_f - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'NiS5Ta2'
_chemical_formula_sum 'Ni S5 Ta2'

loop_
_publ_author_name
 'S. A. Sunshine'
 'J. A. Ibers'
_journal_name_full_name
;
 Inorganic Chemistry
;
_journal_volume 24
_journal_year 1985
_journal_page_first 3611
_journal_page_last 3614
_publ_Section_title
;
 Structure and physical properties of the new layered ternary
      ↪ chalcogenides tantalum nickel sulfide (Ta$_{2}$NiS$_{5}$) and
      ↪ tantalum nickel selenide (Ta$_{2}$NiSe$_{5}$)
;

# Found in Physical and structural properties of the new layered
      ↪ compounds Ta$_{2}$NiS$_{5}$ and Ta$_{2}$NiSe$_{5}$, 1986

_aflow_title 'Ta$_{2}$NiS$_{5}$ Structure'
_aflow_proto 'AB5C2_oC32_63_c_c2f_f'
_aflow_params 'a,b/a,c/a,y_{1},y_{2},y_{3},z_{3},y_{4},z_{4},y_{5},z_{5
      ↪ }'
_aflow_params_values '3.415,3.55666471449,4.42079062958,0.69692,0.31968,
      ↪ 0.58282,0.13527,0.1485,-0.05026,0.22082,0.10879'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oC32'

_symmetry_space_group_name_H-M "C 2/m 2/c 21/m"
_symmetry_Int_Tables_number 63

_cell_length_a     3.41500
_cell_length_b     12.14601
_cell_length_c     15.09700
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z+1/2
4 -x,-y,z+1/2
5 -x,-y,-z
6 -x,y,z
7 x,-y,z+1/2
8 x,y,-z+1/2
9 x+1/2,y+1/2,z
10 x+1/2,-y+1/2,-z
11 -x+1/2,y+1/2,-z+1/2
12 -x+1/2,-y+1/2,z+1/2
13 -x+1/2,-y+1/2,-z
14 -x+1/2,y+1/2,z
15 x+1/2,-y+1/2,z+1/2
16 x+1/2,y+1/2,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ni1 Ni 4 c 0.00000 0.69692 0.25000 1.00000
S1 S 4 c 0.00000 0.31968 0.25000 1.00000
S2 S 8 f 0.00000 0.58282 0.13527 1.00000
S3 S 8 f 0.00000 0.14850 -0.05026 1.00000
Ta1 Ta 8 f 0.00000 0.22082 0.10879 1.00000
```

Ta_2NiS_5: AB5C2_oC32_63_c_c2f_f - POSCAR

```
AB5C2_oC32_63_c_c2f_f & a,b/a,c/a,y1,y2,y3,z3,y4,z4,y5,z5 --params=3.415
      ↪ ,3.55666471449,4.42079062958,0.69692,0.31968,0.58282,0.13527,
      ↪ 0.1485,-0.05026,0.22082,0.10879 & Cmcm D_{2h}^{17} #63 (c^2f^3)
      ↪ & oC32 & None & NiS5Ta2 & NiS5Ta2 & S. A. Sunshine and J. A.
      ↪ Ibers, Inorg. Chem. 24, 3611-3614 (1985)
   1.00000000000000
   1.70750000000000   -6.07300500000000    0.00000000000000
   1.70750000000000    6.07300500000000    0.00000000000000
   0.00000000000000    0.00000000000000   15.09700000000000
     Ni      S      Ta
      2     10       4
Direct
  -0.69692000000000    0.69692000000000    0.25000000000000    Ni   (4c)
   0.69692000000000   -0.69692000000000    0.75000000000000    Ni   (4c)
  -0.31968000000000    0.31968000000000    0.25000000000000    S    (4c)
   0.31968000000000   -0.31968000000000    0.75000000000000    S    (4c)
  -0.58282000000000    0.58282000000000    0.13527000000000    S    (8f)
   0.58282000000000   -0.58282000000000    0.63527000000000    S    (8f)
  -0.58282000000000    0.58282000000000    0.36473000000000    S    (8f)
   0.58282000000000   -0.58282000000000   -0.13527000000000    S    (8f)
  -0.14850000000000    0.14850000000000   -0.05026000000000    S    (8f)
   0.14850000000000   -0.14850000000000    0.44974000000000    S    (8f)
  -0.14850000000000    0.14850000000000    0.55026000000000    S    (8f)
   0.14850000000000   -0.14850000000000    0.05026000000000    S    (8f)
  -0.22082000000000    0.22082000000000    0.10879000000000    Ta   (8f)
   0.22082000000000   -0.22082000000000    0.60879000000000    Ta   (8f)
  -0.22082000000000    0.22082000000000    0.39121000000000    Ta   (8f)
   0.22082000000000   -0.22082000000000   -0.10879000000000    Ta   (8f)
```

V_3AsC: ABC3_oC20_63_c_b_cf - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'AsCV3'
_chemical_formula_sum 'As C V3'

loop_
_publ_author_name
 'H. Boller'
 'H. Nowotny'
_journal_name_full_name
;
 Monatshefte f{\"u}r Chemie und verwandte Teile anderer Wissenschaften
;
_journal_volume 98
_journal_year 1967
_journal_page_first 2127
_journal_page_last 2132
_publ_Section_title
;
 Zum Dreistoff: Vanadin-Arsen-Kohlenstoff
;

# Found in Carbides with Filled Re$_{3}$B-Type Structure, 1994

_aflow_title 'V$_{3}$AsC Structure'
_aflow_proto 'ABC3_oC20_63_c_b_cf'
_aflow_params 'a,b/a,c/a,y_{2},y_{3},y_{4},z_{4}'
_aflow_params_values '3.128,3.24168797954,2.46131713555,0.756,0.452,
      ↪ 0.1283,0.052'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oC20'

_symmetry_space_group_name_H-M "C 2/m 2/c 21/m"
_symmetry_Int_Tables_number 63

_cell_length_a     3.12800
_cell_length_b     10.14000
_cell_length_c     7.69900
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z+1/2
4 -x,-y,z+1/2
5 -x,-y,-z
6 -x,y,z
7 x,-y,z+1/2
8 x,y,-z+1/2
9 x+1/2,y+1/2,z
10 x+1/2,-y+1/2,-z
11 -x+1/2,y+1/2,-z+1/2
12 -x+1/2,-y+1/2,z+1/2
13 -x+1/2,-y+1/2,-z
14 -x+1/2,y+1/2,z
15 x+1/2,-y+1/2,z+1/2
16 x+1/2,y+1/2,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
C1 C 4 b 0.00000 0.50000 0.00000 1.00000
As1 As 4 c 0.00000 0.75600 0.25000 1.00000
V1 V 4 c 0.00000 0.45200 0.25000 1.00000
V2 V 8 f 0.00000 0.12830 0.05200 1.00000
```

V_3AsC: ABC3_oC20_63_c_b_cf - POSCAR

```
ABC3_oC20_63_c_b_cf & a,b/a,c/a,y2,y3,y4,z4 --params=3.128,3.24168797954
    ↪ ,2.46131713555,0.756,0.452,0.1283,0.052 & Cmcm D_{2h}^{17} #63
    ↪ (bc^2f) & oC20 & None & AsCV3 & AsCV3 & H. Boller and H.
    ↪ Nowotny, Monatsh. Chem. Verw. Teile Anderer Wiss. 98, 2127-2132
    ↪ (1967)
   1.00000000000000
   1.56400000000000  -5.07000000000000   0.00000000000000
   1.56400000000000   5.07000000000000   0.00000000000000
   0.00000000000000   0.00000000000000   7.69900000000000
    As    C    V
     2    2    6
Direct
  -0.75600000000000   0.75600000000000   0.25000000000000  As   (4c)
   0.75600000000000  -0.75600000000000   0.75000000000000  As   (4c)
   0.50000000000000   0.50000000000000   0.00000000000000  C    (4b)
   0.50000000000000   0.50000000000000   0.50000000000000  C    (4b)
  -0.45200000000000   0.45200000000000   0.25000000000000  V    (4c)
   0.45200000000000  -0.45200000000000   0.75000000000000  V    (4c)
  -0.12830000000000   0.12830000000000   0.05200000000000  V    (8f)
   0.12830000000000  -0.12830000000000   0.55200000000000  V    (8f)
  -0.12830000000000   0.12830000000000   0.44800000000000  V    (8f)
   0.12830000000000  -0.12830000000000  -0.05200000000000  V    (8f)
```

Si_{24} Clathrate: A_oC24_63_3f - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Clathrate'
_chemical_formula_sum 'Si'

loop_
_publ_author_name
 'D. Y. Kim'
 'S. Stefanoski'
 'O. O. Kurakevych'
 'T. A. Strobel'
_journal_name_full_name
;
 Nature Materials
;
_journal_volume 14
_journal_year 2015
_journal_page_first 169
_journal_page_last 173
_publ_Section_title
;
 Synthesis of an open-framework allotrope of silicon
;

_aflow_title 'Si${24}$ Clathrate Structure'
_aflow_proto 'A_oC24_63_3f'
_aflow_params 'a,b/a,c/a,y_{1},z_{1},y_{2},z_{2},y_{3},z_{3}'
_aflow_params_values '3.82236,2.79950083195,3.30314256114,0.0284,0.5903,
    ↪ 0.2435,0.5551,0.5705,0.3412'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oC24'

_symmetry_space_group_name_H-M "C 2/m 2/c 21/m"
_symmetry_Int_Tables_number 63

_cell_length_a    3.82236
_cell_length_b    10.70070
_cell_length_c    12.62580
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z+1/2
4 -x,-y,z+1/2
5 -x,-y,-z
6 -x,y,z
7 x,-y,z+1/2
8 x,y,-z+1/2
9 x+1/2,y+1/2,z
10 x+1/2,-y+1/2,-z
11 -x+1/2,y+1/2,-z+1/2
12 -x+1/2,-y+1/2,z+1/2
13 -x+1/2,-y+1/2,-z
14 -x+1/2,y+1/2,z
15 x+1/2,-y+1/2,z+1/2
16 x+1/2,y+1/2,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Si1 Si 8 f 0.00000 0.02840 0.59030 1.00000
Si2 Si 8 f 0.00000 0.24350 0.55510 1.00000
Si3 Si 8 f 0.00000 0.57050 0.34120 1.00000
```

Si_{24} Clathrate: A_oC24_63_3f - POSCAR

```
A_oC24_63_3f & a,b/a,c/a,y1,z1,y2,z2,y3,z3 --params=3.82236,
    ↪ 2.79950083195,3.30314256114,0.0284,0.5903,0.2435,0.5551,0.5705,
    ↪ 0.3412 & Cmcm D_{2h}^{17} #63 (f^3) & oC24 & None & Si24 &
    ↪ Clathrate & D. Y. Kim et al., Nat. Mater. 14, 169-173 (2015)
   1.00000000000000
   1.91118000000000  -5.35035000000000   0.00000000000000
   1.91118000000000   5.35035000000000   0.00000000000000
   0.00000000000000   0.00000000000000  12.62580000000000
    Si
    12
Direct
  -0.02840000000000   0.02840000000000   0.59030000000000  Si   (8f)
   0.02840000000000  -0.02840000000000   1.09030000000000  Si   (8f)
  -0.02840000000000   0.02840000000000  -0.09030000000000  Si   (8f)
   0.02840000000000  -0.02840000000000  -0.59030000000000  Si   (8f)
  -0.24350000000000   0.24350000000000   0.55510000000000  Si   (8f)
   0.24350000000000  -0.24350000000000   1.05510000000000  Si   (8f)
  -0.24350000000000   0.24350000000000  -0.05510000000000  Si   (8f)
   0.24350000000000  -0.24350000000000  -0.55510000000000  Si   (8f)
  -0.57050000000000   0.57050000000000   0.34120000000000  Si   (8f)
   0.57050000000000  -0.57050000000000   0.84120000000000  Si   (8f)
  -0.57050000000000   0.57050000000000   0.15880000000000  Si   (8f)
   0.57050000000000  -0.57050000000000  -0.34120000000000  Si   (8f)
```

Base-centered orthorhombic $Sr_4Ru_3O_{10}$: A10B3C4_oC68_64_2dfg_ad_2d - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'O10Ru3Sr4'
_chemical_formula_sum 'O10 Ru3 Sr4'

loop_
_publ_author_name
 'M. K. Crawford'
 'R. L. Harlow'
 'W. Marshall'
 'Z. Li'
 'G. Cao'
 'R. L. Lindstrom'
 'Q. Huang'
 'J. W. Lynn'
_journal_name_full_name
;
 Physical Review B
;
_journal_volume 65
_journal_year 2002
_journal_page_first 214412
_journal_page_last 214412
_publ_Section_title
;
 Structure and magnetism of single crystal Sr$_{4}$Ru$_{3}$O$_{10}$: A
    ↪ ferromagnetic triple-layer ruthenate
;

_aflow_title 'Base-centered orthorhombic Sr$_{4}$Ru$_{3}$O$_{10}$
    ↪ Structure'
_aflow_proto 'A10B3C4_oC68_64_2dfg_ad_2d'
_aflow_params 'a,b/a,c/a,x_{2},x_{3},x_{4},x_{5},x_{6},y_{7},z_{7},x_{8
    ↪ },y_{8},z_{8}'
_aflow_params_values '28.573,0.13649599272,0.13649599272,-0.0695,0.787,
    ↪ 0.8598,0.7039,0.5699,0.7971,0.2972,0.6392,0.2271,0.2279'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oC68'

_symmetry_space_group_name_H-M "C 2/m 2/c 21/a"
_symmetry_Int_Tables_number 64

_cell_length_a    28.57300
_cell_length_b    3.90010
_cell_length_c    3.90010
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x+1/2,y,-z+1/2
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x,y,z
7 x+1/2,-y,z+1/2
8 x+1/2,y,-z+1/2
9 x+1/2,y+1/2,z
10 x+1/2,-y+1/2,-z
11 -x,y+1/2,-z+1/2
12 -x,-y+1/2,z+1/2
13 -x+1/2,-y+1/2,-z
14 -x+1/2,y+1/2,z
15 x,-y+1/2,z+1/2
16 x,y+1/2,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
```

```
_atom_site_occupancy
Ru1 Ru 4 a 0.00000 0.00000 0.00000 1.00000
O1 O 8 d -0.06950 0.00000 0.00000 1.00000
O2 O 8 d 0.78700 0.00000 0.00000 1.00000
Ru2 Ru 8 d 0.85980 0.00000 0.00000 1.00000
Sr1 Sr 8 d 0.70390 0.00000 0.00000 1.00000
Sr2 Sr 8 d 0.56990 0.00000 0.00000 1.00000
O3 O 8 f 0.00000 0.79710 0.29720 1.00000
O4 O 16 g 0.63920 0.22710 0.22790 1.00000
```

Base-centered orthorhombic $Sr_4Ru_3O_{10}$: A10B3C4_oC68_64_2dfg_ad_2d - POSCAR

```
A10B3C4_oC68_64_2dfg_ad_2d & a,b/a,c/a,x2,x3,x4,x5,x6,y7,z7,x8,y8,z8 --
    ↪ params=28.573,0.13649599272,0.13649599272,-0.0695,0.787,0.8598,
    ↪ 0.7039,0.5699,0.7971,0.2972,0.6392,0.2271,0.2279 & Cmca D_{2h
    ↪ }^{18} #64 (ad^5fg) & oC68 & None & O10Ru3Sr4 & O10Ru3Sr4 & M.
    ↪ K. Crawford et al., Phys. Rev. B 65, 214412(2002)
   1.00000000000000
  14.28650000000000  -1.95005000000000   0.00000000000000
  14.28650000000000   1.95005000000000   0.00000000000000
   0.00000000000000   0.00000000000000   3.90010000000000
     O     Ru     Sr
    20      6      8
Direct
  -0.06950000000000  -0.06950000000000   0.00000000000000   O    (8d)
   0.56950000000000   0.56950000000000   0.50000000000000   O    (8d)
   0.06950000000000   0.06950000000000   0.00000000000000   O    (8d)
   0.43050000000000   0.43050000000000   0.50000000000000   O    (8d)
   0.78700000000000   0.78700000000000   0.00000000000000   O    (8d)
  -0.28700000000000  -0.28700000000000   0.50000000000000   O    (8d)
  -0.78700000000000  -0.78700000000000   0.00000000000000   O    (8d)
   1.28700000000000   1.28700000000000   0.50000000000000   O    (8d)
  -0.79710000000000   0.79710000000000   0.29720000000000   O    (8f)
   1.29710000000000  -0.29710000000000   0.79720000000000   O    (8f)
  -0.29710000000000   1.29710000000000   0.20280000000000   O    (8f)
   0.79710000000000  -0.79710000000000  -0.29720000000000   O    (8f)
   0.41210000000000   0.86630000000000   0.22790000000000   O   (16g)
   0.08790000000000  -0.36630000000000   0.72790000000000   O   (16g)
  -0.36630000000000   0.08790000000000   0.27210000000000   O   (16g)
   0.86630000000000   0.41210000000000  -0.22790000000000   O   (16g)
  -0.41210000000000  -0.86630000000000  -0.22790000000000   O   (16g)
   0.91210000000000   1.36630000000000   0.27210000000000   O   (16g)
   1.36630000000000   0.91210000000000   0.72790000000000   O   (16g)
  -0.86630000000000  -0.41210000000000   0.22790000000000   O   (16g)
   0.00000000000000   0.00000000000000   0.00000000000000   Ru   (4a)
   0.50000000000000   0.50000000000000   0.50000000000000   Ru   (4a)
   0.85980000000000   0.85980000000000   0.00000000000000   Ru   (8d)
  -0.35980000000000  -0.35980000000000   0.50000000000000   Ru   (8d)
  -0.85980000000000  -0.85980000000000   0.00000000000000   Ru   (8d)
   1.35980000000000   1.35980000000000   0.50000000000000   Ru   (8d)
   0.70390000000000   0.70390000000000   0.00000000000000   Sr   (8d)
  -0.20390000000000  -0.20390000000000   0.50000000000000   Sr   (8d)
  -0.70390000000000  -0.70390000000000   0.00000000000000   Sr   (8d)
   1.20390000000000   1.20390000000000   0.50000000000000   Sr   (8d)
   0.56990000000000   0.56990000000000   0.00000000000000   Sr   (8d)
  -0.06990000000000  -0.06990000000000   0.50000000000000   Sr   (8d)
  -0.56990000000000  -0.56990000000000   0.00000000000000   Sr   (8d)
   1.06990000000000   1.06990000000000   0.50000000000000   Sr   (8d)
```

$Na_2Mo_2O_7$: A2B2C7_oC88_64_ef_df_3f2g - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Mo2Na2O7'
_chemical_formula_sum 'Mo2 Na2 O7'

loop_
_publ_author_name
 'I. Lindqvist'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 4
_journal_year 1950
_journal_page_first 1066
_journal_page_last 1074
_publ_Section_title
;
 Crystal Structure Studies on Anhydrous Sodium Molybdates and Tungstates
;

_aflow_title 'Na$_{2}$Mo$_{2}$O$_{7}$ Structure'
_aflow_proto 'A2B2C7_oC88_64_ef_df_3f2g'
_aflow_params 'a,b/a,c/a,x_{1},y_{2},y_{3},z_{3},y_{4},z_{4},y_{5},z_{5
    ↪ },y_{6},z_{6},y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9}'
_aflow_params_values '7.17,1.64993026499,2.05020920502,0.25,0.08,0.25,
    ↪ 0.08,0.36,0.3,0.08,0.27,0.41,0.13,0.35,0.49,0.21,0.21,0.16,0.21
    ↪ ,0.49,0.34'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oC88'

_symmetry_space_group_name_H-M "C 2/m 2/c 21/a"
_symmetry_Int_Tables_number 64

_cell_length_a     7.17000
_cell_length_b     11.83000
_cell_length_c     14.70000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x+1/2,y,-z+1/2
4 -x+1/2,-y,z+1/2
5 -x,-y,-z
6 -x,y,z
7 x+1/2,-y,z+1/2
8 x+1/2,y,-z+1/2
9 x+1/2,y+1/2,z
10 x+1/2,-y+1/2,-z
11 -x,y+1/2,-z+1/2
12 -x,-y+1/2,z+1/2
13 -x+1/2,-y+1/2,-z
14 -x+1/2,y+1/2,z
15 x,-y+1/2,z+1/2
16 x,y+1/2,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Na1 Na 8 d 0.25000 0.00000 0.00000 1.00000
Mo1 Mo 8 e 0.25000 0.08000 0.25000 1.00000
Mo2 Mo 8 f 0.00000 0.25000 0.08000 1.00000
Na2 Na 8 f 0.00000 0.36000 0.30000 1.00000
O1 O 8 f 0.00000 0.08000 0.27000 1.00000
O2 O 8 f 0.00000 0.41000 0.13000 1.00000
O3 O 8 f 0.00000 0.35000 0.49000 1.00000
O4 O 16 g 0.21000 0.21000 0.16000 1.00000
O5 O 16 g 0.21000 0.49000 0.34000 1.00000
```

$Na_2Mo_2O_7$: A2B2C7_oC88_64_ef_df_3f2g - POSCAR

```
A2B2C7_oC88_64_ef_df_3f2g & a,b/a,c/a,x1,y2,y3,z3,y4,z4,y5,z5,y6,z6,y7,
    ↪ z7,x8,y8,z8,x9,y9,z9 --params=7.17,1.64993026499,2.05020920502,
    ↪ 0.25,0.08,0.25,0.08,0.36,0.3,0.08,0.27,0.41,0.13,0.35,0.49,0.21
    ↪ ,0.21,0.16,0.21,0.49,0.34 & Cmca D_{2h}^{18} #64 (def^5g^2) &
    ↪ oC88 & None & Mo2Na2O7 & Mo2Na2O7 & I. Lindqvist, Acta Chem.
    ↪ Scand. 4, 1066-1074 (1950)
   1.00000000000000
   3.58500000000000  -5.91500000000000   0.00000000000000
   3.58500000000000   5.91500000000000   0.00000000000000
   0.00000000000000   0.00000000000000  14.70000000000000
     Mo     Na      O
      8      8     28
Direct
   0.17000000000000   0.33000000000000   0.25000000000000   Mo    (8e)
   0.33000000000000   0.17000000000000   0.75000000000000   Mo    (8e)
   0.83000000000000   0.67000000000000   0.75000000000000   Mo    (8e)
   0.67000000000000   0.83000000000000   0.25000000000000   Mo    (8e)
  -0.25000000000000   0.25000000000000   0.08000000000000   Mo    (8f)
   0.75000000000000   0.25000000000000   0.58000000000000   Mo    (8f)
   0.25000000000000   0.75000000000000   0.42000000000000   Mo    (8f)
   0.25000000000000  -0.25000000000000  -0.08000000000000   Mo    (8f)
   0.25000000000000   0.25000000000000   0.00000000000000   Na    (8d)
   0.25000000000000   0.25000000000000   0.50000000000000   Na    (8d)
  -0.25000000000000  -0.25000000000000   0.00000000000000   Na    (8d)
   0.75000000000000   0.75000000000000   0.50000000000000   Na    (8d)
  -0.36000000000000   0.36000000000000   0.30000000000000   Na    (8f)
   0.86000000000000   0.14000000000000   0.80000000000000   Na    (8f)
   0.14000000000000   0.86000000000000   0.20000000000000   Na    (8f)
   0.36000000000000  -0.36000000000000  -0.30000000000000   Na    (8f)
  -0.08000000000000   0.08000000000000   0.27000000000000    O    (8f)
   0.58000000000000   0.42000000000000   0.77000000000000    O    (8f)
   0.42000000000000   0.58000000000000   0.23000000000000    O    (8f)
   0.08000000000000  -0.08000000000000  -0.27000000000000    O    (8f)
  -0.41000000000000   0.41000000000000   0.13000000000000    O    (8f)
   0.91000000000000   0.09000000000000   0.63000000000000    O    (8f)
   0.09000000000000   0.91000000000000   0.37000000000000    O    (8f)
   0.41000000000000  -0.41000000000000  -0.13000000000000    O    (8f)
  -0.35000000000000   0.35000000000000   0.49000000000000    O    (8f)
   0.85000000000000   0.15000000000000   0.99000000000000    O    (8f)
   0.15000000000000   0.85000000000000   0.01000000000000    O    (8f)
   0.35000000000000  -0.35000000000000  -0.49000000000000    O    (8f)
   0.00000000000000   0.42000000000000   0.16000000000000    O   (16g)
   0.50000000000000   0.08000000000000   0.66000000000000    O   (16g)
   0.08000000000000   0.50000000000000   0.34000000000000    O   (16g)
   0.42000000000000   0.00000000000000  -0.16000000000000    O   (16g)
   0.00000000000000  -0.42000000000000  -0.16000000000000    O   (16g)
   0.50000000000000   0.92000000000000   0.34000000000000    O   (16g)
   0.92000000000000   0.50000000000000   0.66000000000000    O   (16g)
  -0.42000000000000   0.00000000000000   0.16000000000000    O   (16g)
  -0.28000000000000   0.70000000000000   0.34000000000000    O   (16g)
   0.78000000000000  -0.20000000000000   0.84000000000000    O   (16g)
  -0.20000000000000   0.78000000000000   0.16000000000000    O   (16g)
   0.70000000000000  -0.28000000000000  -0.34000000000000    O   (16g)
   0.28000000000000  -0.70000000000000  -0.34000000000000    O   (16g)
   0.22000000000000   1.20000000000000   0.16000000000000    O   (16g)
   1.20000000000000   0.22000000000000   0.84000000000000    O   (16g)
  -0.70000000000000   0.28000000000000   0.34000000000000    O   (16g)
```

Li_2PrO_3: A2B3C_oC12_65_h_bh_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Li2O3Pr'
```

```
_chemical_formula_sum 'Li2 O3 Pr'

loop_
_publ_author_name
 'Y. Hinatsu'
 'Y. Doi'
_journal_name_full_name
;
 Journal of Alloys and Compounds
;
_journal_volume 418
_journal_year 2006
_journal_page_first 155
_journal_page_last 160
_publ_Section_title
;
 Crystal structures and magnetic properties of alkali-metal lanthanide
     ↪ oxides $A_{2}$$Ln$O$_{3}$ ($A$ = Li, Na; $Ln$ = Ce, Pr, Tb)
;

_aflow_title 'Li$_{2}$PrO$_{3}$ Structure'
_aflow_proto 'A2B3C_oC12_65_h_bh_a'
_aflow_params 'a,b/a,c/a,x_{3},x_{4}'
_aflow_params_values '9.5198,0.467551839324,0.36581650875,0.6556,0.8557'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oC12'

_symmetry_space_group_name_H-M "C 2/m 2/m 2/m"
_symmetry_Int_Tables_number 65

_cell_length_a      9.51980
_cell_length_b      4.45100
_cell_length_c      3.48250
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 -x,-y,-z
6 -x,y,z
7 x,-y,z
8 x,y,-z
9 x+1/2,y+1/2,z
10 x+1/2,-y+1/2,-z
11 -x+1/2,y+1/2,-z
12 -x+1/2,-y+1/2,z
13 -x+1/2,-y+1/2,-z
14 -x+1/2,y+1/2,z
15 x+1/2,-y+1/2,z
16 x+1/2,y+1/2,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Pr1 Pr 2 a 0.00000 0.00000 0.00000 1.00000
O1 O 2 b 0.50000 0.00000 0.00000 1.00000
Li1 Li 4 h 0.65560 0.00000 0.50000 1.00000
O2 O 4 h 0.85570 0.00000 0.50000 1.00000
```

Li_2PrO_3: A2B3C_oC12_65_h_bh_a - POSCAR

```
A2B3C_oC12_65_h_bh_a & a,b/a,c/a,x3,x4 --params=9.5198,0.467551839324,
    ↪ 0.36581650875,0.6556,0.8557 & Cmmm D_{2h}^{19} #65 (abh^2) &
    ↪ oC12 & None & Li2O3Pr & Li2O3Pr & Y. Hinatsu and Y. Doi, J.
    ↪ Alloys Compd. 418, 155-160 (2006)
   1.00000000000000
   4.75990000000000  -2.22550000000000   0.00000000000000
   4.75990000000000   2.22550000000000   0.00000000000000
   0.00000000000000   0.00000000000000   3.48250000000000
     Li    O    Pr
      2    3     1
Direct
   0.65560000000000   0.65560000000000   0.50000000000000   Li   (4h)
  -0.65560000000000  -0.65560000000000   0.50000000000000   Li   (4h)
   0.50000000000000   0.50000000000000   0.00000000000000   O    (2b)
   0.85570000000000   0.85570000000000   0.50000000000000   O    (4h)
  -0.85570000000000  -0.85570000000000   0.50000000000000   O    (4h)
   0.00000000000000   0.00000000000000   0.00000000000000   Pr   (2a)
```

$Mg(NH_3)_2Cl_2$ ($E1_3$): A2B8CD2_oC26_65_h_r_a_i - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Cl2H6MgN2'
_chemical_formula_sum 'C12 H8 Mg N2'

loop_
_publ_author_name
 'A. Leineweber'
 'M. W. Friedriszik'
 'H. Jacobs'
_journal_name_full_name
;
 Journal of Solid State Chemistry
;
_journal_volume 147
_journal_year 1999
_journal_page_first 229
_journal_page_last 234
_publ_Section_title
;
 Preparation and Crystal Structures of Mg(NH$_{3}$)$_{2}$Cl$_{2}$, Mg(
     ↪ NH$_{3}$)$_{2}$Br$_{2}$, and Mg(NH$_{3}$)$_{2}$I$_{2}$
;

_aflow_title 'Mg(NH$_{3}$)$_{2}$Cl$_{2}$ ($E1_{3}$) Structure'
_aflow_proto 'A2B8CD2_oC26_65_h_r_a_i'
_aflow_params 'a,b/a,c/a,x_{2},y_{3},x_{4},y_{4},z_{4}'
_aflow_params_values '8.18099,1.00314142909,0.458990904524,0.2133,0.2595
    ↪ ,0.045,0.312,0.158'
_aflow_Strukturbericht '$E1_{3}$'
_aflow_Pearson 'oC26'

_symmetry_space_group_name_H-M "C 2/m 2/m 2/m"
_symmetry_Int_Tables_number 65

_cell_length_a      8.18099
_cell_length_b      8.20669
_cell_length_c      3.75500
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 -x,-y,-z
6 -x,y,z
7 x,-y,z
8 x,y,-z
9 x+1/2,y+1/2,z
10 x+1/2,-y+1/2,-z
11 -x+1/2,y+1/2,-z
12 -x+1/2,-y+1/2,z
13 -x+1/2,-y+1/2,-z
14 -x+1/2,y+1/2,z
15 x+1/2,-y+1/2,z
16 x+1/2,y+1/2,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Mg1 Mg 2 a 0.00000 0.00000 0.00000 1.00000
Cl1 Cl 4 h 0.21330 0.00000 0.50000 1.00000
N1 N 4 i 0.00000 0.25950 0.00000 1.00000
H1 H 16 r 0.04500 0.31200 0.15800 0.75000
```

$Mg(NH_3)_2Cl_2$ ($E1_3$): A2B8CD2_oC26_65_h_r_a_i - POSCAR

```
A2B8CD2_oC26_65_h_r_a_i & a,b/a,c/a,x2,y3,x4,y4,z4 --params=8.18099,
    ↪ 1.00314142909,0.458990904524,0.2133,0.2595,0.045,0.312,0.158 &
    ↪ Cmmm D_{2h}^{19} #65 (ahir) & oC26 & $E1_{3}$ & Cl2H6MgN2 &
    ↪ Cl2H6MgN2 & A. Leineweber and M. W. Friedriszik and H. Jacobs,
    ↪ J. Solid State Chem. 147, 229-234 (1999)
   1.00000000000000
   4.09049500000000  -4.10334500000000   0.00000000000000
   4.09049500000000   4.10334500000000   0.00000000000000
   0.00000000000000   0.00000000000000   3.75500000000000
     Cl    H    Mg    N
      2    8     1    2
Direct
   0.21330000000000   0.21330000000000   0.50000000000000   Cl   (4h)
  -0.21330000000000  -0.21330000000000   0.50000000000000   Cl   (4h)
  -0.26700000000000   0.35700000000000   0.15800000000000   H   (16r)
   0.26700000000000  -0.35700000000000   0.15800000000000   H   (16r)
  -0.35700000000000   0.26700000000000  -0.15800000000000   H   (16r)
   0.35700000000000  -0.26700000000000  -0.15800000000000   H   (16r)
   0.26700000000000  -0.35700000000000  -0.15800000000000   H   (16r)
  -0.26700000000000   0.35700000000000  -0.15800000000000   H   (16r)
   0.35700000000000  -0.26700000000000   0.15800000000000   H   (16r)
  -0.35700000000000   0.26700000000000   0.15800000000000   H   (16r)
   0.00000000000000   0.00000000000000   0.00000000000000   Mg   (2a)
  -0.25950000000000   0.25950000000000   0.00000000000000   N   (4i)
   0.25950000000000  -0.25950000000000   0.00000000000000   N   (4i)
```

Nb_3O_7F: A3B8_oC22_65_ag_bd2gh - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'FNb3O7'
_chemical_formula_sum 'Nb3 O8'

loop_
_publ_author_name
```

```
 'S. Andersson'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 18
_journal_year 1964
_journal_page_first 2339
_journal_page_last 2344
_publ_Section_title
;
 The Crystal Structure of Nb$_{3}$O$_{7}$F
;

_aflow_title 'Nb$_{3}$O$_{7}$F Structure'
_aflow_proto 'A3B8_oC22_65_ag_bd2gh'
_aflow_params 'a,b/a,c/a,x_{4},x_{5},x_{6},x_{7}'
_aflow_params_values '20.66999,0.185437922321,0.189985578126,0.1836,
      ↪ 0.094,0.71,0.189'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oC22'

_symmetry_space_group_name_H-M "C 2/m 2/m 2/m"
_symmetry_Int_Tables_number 65

_cell_length_a     20.66999
_cell_length_b     3.83300
_cell_length_c     3.92700
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 -x,-y,-z
6 -x,y,z
7 x,-y,z
8 x,y,-z
9 x+1/2,y+1/2,z
10 x+1/2,-y+1/2,-z
11 -x+1/2,y+1/2,-z
12 -x+1/2,-y+1/2,z
13 -x+1/2,-y+1/2,-z
14 -x+1/2,y+1/2,z
15 x+1/2,-y+1/2,z
16 x+1/2,y+1/2,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Nb1 Nb 2 a 0.00000 0.00000 0.00000 1.00000
O1 O 2 b 0.50000 0.00000 0.00000 1.00000
O2 O 2 d 0.00000 0.00000 0.50000 1.00000
Nb2 Nb 4 g 0.18360 0.00000 0.00000 1.00000
O3 O 4 g 0.09400 0.00000 0.00000 1.00000
O4 O 4 g 0.71000 0.00000 0.00000 1.00000
O5 O 4 h 0.18900 0.00000 0.50000 1.00000
```

Nb_3O_7F: A3B8_oC22_65_ag_bd2gh - POSCAR

```
A3B8_oC22_65_ag_bd2gh & a,b/a,c/a,x4,x5,x6,x7 --params=20.66999,
      ↪ 0.185437922321,0.189985578126,0.1836,0.094,0.71,0.189 & Cmmm D_
      ↪ {2h}^{19} #65 (abdg^3h) & oC22 & None & FNb3O7 & FNb3O7 & S.
      ↪ Andersson, Acta Chem. Scand. 18, 2339-2344 (1964)
   1.00000000000000
  10.33499500000000   -1.91650000000000    0.00000000000000
  10.33499500000000    1.91650000000000    0.00000000000000
   0.00000000000000    0.00000000000000    3.92700000000000
     Nb      O
      3      8
Direct
   0.00000000000000    0.00000000000000    0.00000000000000    Nb    (2a)
   0.18360000000000    0.18360000000000    0.00000000000000    Nb    (4g)
  -0.18360000000000   -0.18360000000000    0.00000000000000    Nb    (4g)
   0.50000000000000    0.50000000000000    0.00000000000000    O     (2b)
   0.00000000000000    0.00000000000000    0.50000000000000    O     (2d)
   0.09400000000000    0.09400000000000    0.00000000000000    O     (4g)
  -0.09400000000000   -0.09400000000000    0.00000000000000    O     (4g)
   0.71000000000000    0.71000000000000    0.00000000000000    O     (4g)
  -0.71000000000000   -0.71000000000000    0.00000000000000    O     (4g)
   0.18900000000000    0.18900000000000    0.50000000000000    O     (4h)
  -0.18900000000000   -0.18900000000000    0.50000000000000    O     (4h)
```

$NH_4H_2PO_2$ ($F5_7$): A2BC2D_oC24_67_m_a_n_g - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'H2(NH4)O2P'
_chemical_formula_sum 'H2 (NH4) O2 P'

loop_
_publ_author_name
 'W. H. Zachariasen'
 'R. C. L. Mooney'
_journal_name_full_name
;
 Journal of Chemical Physics
;
_journal_volume 2
_journal_year 1934
_journal_page_first 34
_journal_page_last 37
_publ_Section_title
;
 The Structure of the Hypophosphite Group as Determined from the Crystal
      ↪ Lattice of Ammonium Hypophosphite
;

# Found in Strukturbericht Band III 1933-1935, 1937

_aflow_title 'NH$_{4}$H$_{2}$PO$_{2}$ ($F5_{7}$) Structure'
_aflow_proto 'A2BC2D_oC24_67_m_a_n_g'
_aflow_params 'a,b/a,c/a,z_{2},y_{3},z_{3},x_{4},z_{4}'
_aflow_params_values '11.47,0.659982563208,0.346992153444,0.458,0.608,
      ↪ 0.806,0.386,0.348'
_aflow_Strukturbericht '$F5_{7}$'
_aflow_Pearson 'oC24'

_symmetry_space_group_name_H-M "C 2/m 2/m 2/a"
_symmetry_Int_Tables_number 67

_cell_length_a     11.47000
_cell_length_b     7.57000
_cell_length_c     3.98000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x+1/2,y,-z
4 -x+1/2,-y,z
5 -x,-y,-z
6 -x,y,z
7 x+1/2,-y,z
8 x+1/2,y,-z
9 x+1/2,y+1/2,z
10 x+1/2,-y+1/2,-z
11 -x,y+1/2,-z
12 -x,-y+1/2,z
13 -x+1/2,-y+1/2,-z
14 -x+1/2,y+1/2,z
15 x,-y+1/2,z
16 x,y+1/2,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
NH41 NH4 4 a 0.25000 0.00000 0.00000 1.00000
P1 P 4 g 0.00000 0.25000 0.45800 1.00000
H1 H 8 m 0.00000 0.60800 0.80600 1.00000
O1 O 8 n 0.38600 0.25000 0.34800 1.00000
```

$NH_4H_2PO_2$ ($F5_7$): A2BC2D_oC24_67_m_a_n_g - POSCAR

```
A2BC2D_oC24_67_m_a_n_g & a,b/a,c/a,z2,y3,z3,x4,z4 --params=11.47,
      ↪ 0.659982563208,0.346992153444,0.458,0.608,0.806,0.386,0.348 &
      ↪ Cmma D_{2h}^{21} #67 (agmn) & oC24 & $F5_{7}$ & H2(NH4)O2P & H2
      ↪ (NH4)O2P & W. H. Zachariasen and R. C. L. Mooney, J. Chem.
      ↪ Phys. 2, 34-37 (1934)
   1.00000000000000
   5.73500000000000   -3.78500000000000    0.00000000000000
   5.73500000000000    3.78500000000000    0.00000000000000
   0.00000000000000    0.00000000000000    3.98000000000000
     H    NH4      O      P
     4      2      4      2
Direct
  -0.60800000000000    0.60800000000000    0.80600000000000    H     (8m)
   1.10800000000000   -0.10800000000000    0.80600000000000    H     (8m)
  -0.10800000000000    1.10800000000000   -0.80600000000000    H     (8m)
   0.60800000000000   -0.60800000000000   -0.80600000000000    H     (8m)
   0.25000000000000    0.25000000000000    0.00000000000000    NH4   (4a)
   0.75000000000000    0.75000000000000    0.00000000000000    NH4   (4a)
   1.13600000000000    0.63600000000000    0.34800000000000    O     (8n)
   0.36400000000000   -0.13600000000000    0.34800000000000    O     (8n)
  -0.13600000000000    0.36400000000000   -0.34800000000000    O     (8n)
   0.63600000000000    1.13600000000000   -0.34800000000000    O     (8n)
   0.75000000000000    0.25000000000000    0.45800000000000    P     (4g)
   0.25000000000000    0.75000000000000   -0.45800000000000    P     (4g)
```

Thenardite [Na_2SO_4 (V), $H1_7$]: A2B4C_oF56_70_g_h_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Thenardite'
_chemical_formula_sum 'Na2 O4 S'
```

```
loop_
_publ_author_name
 'A. G. Nord'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 27
_journal_year 1973
_journal_page_first 814
_journal_page_last 822
_publ_Section_title
;
 Refinement of the Crystal Structure of Thenardite, Na$_{2}$SO$_{4}$(V)
;

_aflow_title 'Thenardite [Na$_{2}$SO$_{4}$ (V), $H1_{7}$] Structure'
_aflow_proto 'A2B4C_oF56_70_g_h_a'
_aflow_params 'a,b/a,c/a,z_{2},x_{3},y_{3},z_{3}'
_aflow_params_values '5.8596,2.09987029831,1.67537033245,0.4414,-0.0203,
      ↪ 0.0572,0.2137'
_aflow_Strukturbericht '$H1_{7}$'
_aflow_Pearson 'oF56'

_symmetry_space_group_name_H-M "F 2/d 2/d 2/d (origin choice 2)"
_symmetry_Int_Tables_number 70

_cell_length_a     5.85960
_cell_length_b     12.30440
_cell_length_c     9.81700
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y+3/4,-z+3/4
3 -x+3/4,y,-z+3/4
4 -x+3/4,-y+3/4,z
5 -x,-y,-z
6 -x,y+1/4,z+1/4
7 x+1/4,-y,z+1/4
8 x+1/4,y+1/4,-z
9 x,y+1/2,z+1/2
10 x,-y+1/4,-z+1/4
11 -x+3/4,y+1/2,-z+1/4
12 -x+3/4,-y+1/4,z+1/2
13 -x,-y+1/2,-z+1/2
14 -x,y+3/4,z+3/4
15 x+1/4,-y+1/2,z+3/4
16 x+1/4,y+3/4,-z+1/2
17 x+1/2,y,z+1/2
18 x+1/2,-y+3/4,-z+1/4
19 -x+1/4,y,-z+1/4
20 -x+1/4,-y+3/4,z+1/2
21 -x+1/2,-y,-z+1/2
22 -x+1/2,y+1/4,z+3/4
23 x+3/4,-y,z+3/4
24 x+3/4,y+1/4,-z+1/2
25 x+1/2,y+1/2,z
26 x+1/2,-y+1/4,-z+3/4
27 -x+1/4,y+1/2,-z+3/4
28 -x+1/4,-y+1/4,z
29 -x+1/2,-y+1/2,-z
30 -x+1/2,y+3/4,z+1/4
31 x+3/4,-y+1/2,z+1/4
32 x+3/4,y+3/4,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
S1 S 8 a 0.12500 0.12500 0.12500 1.00000
Na1 Na 16 g 0.12500 0.12500 0.44140 1.00000
O1 O 32 h -0.02030 0.05720 0.21370 1.00000
```

Thenardite [Na_2SO_4 (V), $H1_7$]: A2B4C_oF56_70_g_h_a - POSCAR

```
A2B4C_oF56_70_g_h_a & a,b/a,c/a,z2,x3,y3,z3 --params=5.8596,
      ↪ 2.09987029831,1.67537033245,0.4414,-0.0203,0.0572,0.2137 & Fddd
      ↪ D_{2h}^{24} #70 (agh) & oF56 & $H1_{7}$ & Na2O4S & Thenardite
      ↪ & A. G. Nord, Acta Chem. Scand. 27, 814-822 (1973)
   1.00000000000000
   0.00000000000000    6.15220000000000    4.90850000000000
   2.92980000000000    0.00000000000000    4.90850000000000
   2.92980000000000    6.15220000000000    0.00000000000000
    Na      O      S
     4      8      2
Direct
   0.44140000000000    0.44140000000000   -0.19140000000000   Na  (16g)
  -0.19140000000000   -0.19140000000000    0.44140000000000   Na  (16g)
  -0.44140000000000   -0.44140000000000    1.19140000000000   Na  (16g)
   1.19140000000000    1.19140000000000   -0.44140000000000   Na  (16g)
   0.29120000000000    0.13620000000000   -0.17680000000000   O   (32h)
   0.13620000000000    0.29120000000000    0.24940000000000   O   (32h)
  -0.17680000000000    0.24940000000000    0.29120000000000   O   (32h)
   0.24940000000000   -0.17680000000000    0.13620000000000   O   (32h)
  -0.29120000000000   -0.13620000000000    0.17680000000000   O   (32h)
  -0.13620000000000   -0.29120000000000    0.75060000000000   O   (32h)
   0.17680000000000    0.75060000000000   -0.29120000000000   O   (32h)
   0.75060000000000    0.17680000000000   -0.13620000000000   O   (32h)
   0.12500000000000    0.12500000000000    0.12500000000000   S   (8a)
   0.87500000000000    0.87500000000000    0.87500000000000   S   (8a)
```

Mg_2Cu (C_b): AB2_oF48_70_g_fg - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'CuMg2'
_chemical_formula_sum 'Cu Mg2'

loop_
_publ_author_name
 'F. Gingl'
 'P. Selvam'
 'K. Yvon'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 49
_journal_year 1993
_journal_page_first 201
_journal_page_last 203
_publ_Section_title
;
 Structure refinement of Mg$_{2}$Cu and a comparison of the Mg$_{2}$Cu,
      ↪ Mg$_{2}$Ni and Al$_{2}$Cu structure types
;

# Found in Crystal Structure of CuMg$_{2}$, {Crystallography online.com
      ↪ }, 2018

_aflow_title 'Mg$_{2}$Cu ($C_{b}$) Structure'
_aflow_proto 'AB2_oF48_70_g_fg'
_aflow_params 'a,b/a,c/a,y_{1},z_{2},z_{3}'
_aflow_params_values '5.275,1.71450236967,3.47450236967,0.4586,0.49819,
      ↪ 0.0415'
_aflow_Strukturbericht '$C_{b}$'
_aflow_Pearson 'oF48'

_symmetry_space_group_name_H-M "F 2/d 2/d 2/d (origin choice 2)"
_symmetry_Int_Tables_number 70

_cell_length_a     5.27500
_cell_length_b     9.04400
_cell_length_c     18.32800
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y+3/4,-z+3/4
3 -x+3/4,y,-z+3/4
4 -x+3/4,-y+3/4,z
5 -x,-y,-z
6 -x,y+1/4,z+1/4
7 x+1/4,-y,z+1/4
8 x+1/4,y+1/4,-z
9 x,y+1/2,z+1/2
10 x,-y+1/4,-z+1/4
11 -x+3/4,y+1/2,-z+1/4
12 -x+3/4,-y+1/4,z+1/2
13 -x,-y+1/2,-z+1/2
14 -x,y+3/4,z+3/4
15 x+1/4,-y+1/2,z+3/4
16 x+1/4,y+3/4,-z+1/2
17 x+1/2,y,z+1/2
18 x+1/2,-y+3/4,-z+1/4
19 -x+1/4,y,-z+1/4
20 -x+1/4,-y+3/4,z+1/2
21 -x+1/2,-y,-z+1/2
22 -x+1/2,y+1/4,z+3/4
23 x+3/4,-y,z+3/4
24 x+3/4,y+1/4,-z+1/2
25 x+1/2,y+1/2,z
26 x+1/2,-y+1/4,-z+3/4
27 -x+1/4,y+1/2,-z+3/4
28 -x+1/4,-y+1/4,z
29 -x+1/2,-y+1/2,-z
30 -x+1/2,y+3/4,z+1/4
31 x+3/4,-y+1/2,z+1/4
32 x+3/4,y+3/4,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Mg1 Mg 16 f 0.12500 0.45860 0.12500 1.00000
Cu1 Cu 16 g 0.12500 0.12500 0.49819 1.00000
Mg2 Mg 16 g 0.12500 0.12500 0.04150 1.00000
```

Mg_2Cu (C_b): AB2_oF48_70_g_fg - POSCAR

```
AB2_oF48_70_g_fg & a,b/a,c/a,y1,z2,z3 --params=5.275,1.71450236967,
    ↪ 3.47450236967,0.4586,0.49819,0.0415 & Fddd D_{2h}^{24} #70 (fg^
    ↪ 2) & oF48 & $C_{b}$ & CuMg2 & CuMg2 & F. Gingl and P. Selvam
    ↪ and K. Yvon, Acta Crystallogr. Sect. B Struct. Sci. 49, 201-203
    ↪ (1993)
   1.00000000000000
   0.00000000000000   4.52200000000000   9.16400000000000
   2.63750000000000   0.00000000000000   9.16400000000000
   2.63750000000000   4.52200000000000   0.00000000000000
    Cu    Mg
     4     8
Direct
   0.49819000000000   0.49819000000000  -0.24819000000000   Cu   (16g)
  -0.24819000000000  -0.24819000000000   0.49819000000000   Cu   (16g)
  -0.49819000000000  -0.49819000000000   1.24819000000000   Cu   (16g)
   1.24819000000000   1.24819000000000  -0.49819000000000   Cu   (16g)
   0.45860000000000  -0.20860000000000   0.45860000000000   Mg   (16f)
  -0.20860000000000   0.45860000000000  -0.20860000000000   Mg   (16f)
  -0.45860000000000   1.20860000000000  -0.45860000000000   Mg   (16f)
   1.20860000000000  -0.45860000000000   1.20860000000000   Mg   (16f)
   0.04150000000000   0.04150000000000   0.20850000000000   Mg   (16g)
   0.20850000000000   0.20850000000000   0.04150000000000   Mg   (16g)
  -0.04150000000000  -0.04150000000000   0.79150000000000   Mg   (16g)
   0.79150000000000   0.79150000000000  -0.04150000000000   Mg   (16g)
```

High-Temperature Cryolite (Na_3AlF_6): AB6C3_oI20_71_a_in_cj - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Cryolite'
_chemical_formula_sum 'Al F6 Na3'

loop_
_publ_author_name
 'H. Yang'
 'S. Ghose'
 'D. M. Hatch'
_journal_name_full_name
;
 Physics and Chemistry of Minerals
;
_journal_volume 19
_journal_year 1993
_journal_page_first 528
_journal_page_last 544
_publ_Section_title
;
 Ferroelastic phase transition in cryolite, Na$_{3}$AlF$_{6}$, a mixed
    ↪ fluoride perovskite: High temperature single crystal X-ray
    ↪ diffraction study and symmetry analysis of the transition
    ↪ mechanism
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'High-Temperature Cryolite (Na$_{3}$AlF$_{6}$) Structure'
_aflow_proto 'AB6C3_oI20_71_a_in_cj'
_aflow_params 'a,b/a,c/a,z_{3},z_{4},x_{5},y_{5}'
_aflow_params_values '5.6333,0.99893490494,1.41336339268,0.2192,0.2485,
    ↪ 0.2335,0.2166'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oI20'

_symmetry_space_group_name_H-M "I 2/m 2/m 2/m"
_symmetry_Int_Tables_number 71

_cell_length_a    5.63330
_cell_length_b    5.62730
_cell_length_c    7.96190
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 -x,-y,-z
6 -x,y,z
7 x,-y,z
8 x,y,-z
9 x+1/2,y+1/2,z+1/2
10 x+1/2,-y+1/2,-z+1/2
11 -x+1/2,y+1/2,-z+1/2
12 -x+1/2,-y+1/2,z+1/2
13 -x+1/2,-y+1/2,-z+1/2
14 -x+1/2,y+1/2,z+1/2
15 x+1/2,-y+1/2,z+1/2
16 x+1/2,y+1/2,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Al1 Al 2 a 0.00000 0.00000 0.00000 1.00000
Na1 Na 2 c 0.50000 0.50000 0.00000 1.00000
F1 F 4 i 0.00000 0.00000 0.21920 1.00000
Na2 Na 4 j 0.50000 0.00000 0.24850 1.00000
F2 F 8 n 0.23350 0.21660 0.00000 1.00000
```

High-Temperature Cryolite (Na_3AlF_6): AB6C3_oI20_71_a_in_cj - POSCAR

```
AB6C3_oI20_71_a_in_cj & a,b/a,c/a,z3,z4,x5,y5 --params=5.6333,
    ↪ 0.99893490494,1.41336339268,0.2192,0.2485,0.2335,0.2166 & Immm
    ↪ D_{2h}^{25} #71 (acijn) & oI20 & None & AlF6Na3 & Cryolite & H.
    ↪ Yang and S. Ghose and D. M. Hatch, Phys. Chem. Miner. 19,
    ↪ 528-544 (1993)
   1.00000000000000
  -2.81665000000000   2.81365000000000   3.98095000000000
   2.81665000000000  -2.81365000000000   3.98095000000000
   2.81665000000000   2.81365000000000  -3.98095000000000
    Al    F    Na
     1    6     3
Direct
   0.00000000000000   0.00000000000000   0.00000000000000   Al   (2a)
   0.21920000000000   0.21920000000000   0.00000000000000   F    (4i)
  -0.21920000000000  -0.21920000000000   0.00000000000000   F    (4i)
   0.21660000000000   0.23350000000000   0.45010000000000   F    (8n)
  -0.21660000000000  -0.23350000000000  -0.45010000000000   F    (8n)
   0.21660000000000  -0.23350000000000  -0.01690000000000   F    (8n)
  -0.21660000000000   0.23350000000000   0.01690000000000   F    (8n)
   0.50000000000000   0.50000000000000   0.00000000000000   Na   (2c)
   0.24850000000000   0.74850000000000   0.50000000000000   Na   (4j)
  -0.24850000000000   0.25150000000000   0.50000000000000   Na   (4j)
```

$CsFeS_2$ (100 K): ABC2_oI16_71_g_i_eh - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'CsFeS2'
_chemical_formula_sum 'Cs Fe S2'

loop_
_publ_author_name
 'Y. Ito'
 'M. Nishi'
 'C. F. Majkrzak'
 'L. Passell'
_journal_name_full_name
;
 Journal of the Physical Society of Japan
;
_journal_volume 54
_journal_year 1985
_journal_page_first 348
_journal_page_last 357
_publ_Section_title
;
 Low Temperature Powder Neutron Diffraction Studies of CsFeS$_{2}$
;

# Found in CsFeS2 (100K) Crystal Structure, {PAULING FILE in: Inorganic
    ↪ Solid Phases, SpringerMaterials (online database), Springer,
    ↪ Heidelberg (ed.) SpringerMaterials }, 2016

_aflow_title 'CsFeS$_{2}$ (100~K) Structure'
_aflow_proto 'ABC2_oI16_71_g_i_eh'
_aflow_params 'a,b/a,c/a,x_{1},y_{2},y_{3},z_{4}'
_aflow_params_values '7.09,1.67136812412,0.764456981664,0.256,0.33,0.148
    ↪ ,0.249'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oI16'

_symmetry_space_group_name_H-M "I 2/m 2/m 2/m"
_symmetry_Int_Tables_number 71

_cell_length_a    7.09000
_cell_length_b    11.85000
_cell_length_c    5.42000
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 -x,-y,-z
6 -x,y,z
7 x,-y,z
8 x,y,-z
9 x+1/2,y+1/2,z+1/2
10 x+1/2,-y+1/2,-z+1/2
11 -x+1/2,y+1/2,-z+1/2
12 -x+1/2,-y+1/2,z+1/2
13 -x+1/2,-y+1/2,-z+1/2
14 -x+1/2,y+1/2,z+1/2
15 x+1/2,-y+1/2,z+1/2
16 x+1/2,y+1/2,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
```

```
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
S1 S 4 e 0.25600 0.00000 0.00000 1.00000
Cs1 Cs 4 g 0.00000 0.33000 0.00000 1.00000
S2 S 4 h 0.00000 0.14800 0.50000 1.00000
Fe1 Fe 4 i 0.00000 0.00000 0.24900 1.00000
```

$CsFeS_2$ (100 K): ABC2_oI16_71_g_i_eh - POSCAR

```
ABC2_oI16_71_g_i_eh & a,b/a,c/a,x1,y2,y3,z4 --params=7.09,1.67136812412,
    ↪ 0.764456981664,0.256,0.33,0.148,0.249 & Immm D_{2h}^{25} #71 (
    ↪ eghi) & oI16 & None & CsFeS2 & CsFeS2 & Y. Ito et al., J. Phys.
    ↪ Soc. Jpn. 54, 348-357 (1985)
   1.00000000000000
  -3.54500000000000    5.92500000000000    2.71000000000000
   3.54500000000000   -5.92500000000000    2.71000000000000
   3.54500000000000    5.92500000000000   -2.71000000000000
     Cs     Fe      S
      2      2      4
Direct
   0.33000000000000    0.00000000000000    0.33000000000000   Cs   (4g)
  -0.33000000000000    0.00000000000000   -0.33000000000000   Cs   (4g)
   0.24900000000000    0.24900000000000    0.00000000000000   Fe   (4i)
  -0.24900000000000   -0.24900000000000    0.00000000000000   Fe   (4i)
   0.00000000000000    0.25600000000000    0.25600000000000   S    (4e)
   0.00000000000000   -0.25600000000000   -0.25600000000000   S    (4e)
   0.64800000000000    0.50000000000000    0.14800000000000   S    (4h)
   0.35200000000000    0.50000000000000   -0.14800000000000   S    (4h)
```

CsO: AB_oI8_71_g_i - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'CsO'
_chemical_formula_sum 'Cs O'

loop_
_publ_author_name
 'M. E. Rengade'
_journal_name_full_name
;
 Comptes Rendus de l'Acad{\'e}mie des Sciences
;
_journal_volume 148
_journal_year 1909
_journal_page_first 1199
_journal_page_last 1202
_publ_Section_title
;
 Sur les Sous-Oxydes de Caesium
;

# Found in Binary Alloy Phase Diagrams, {Cd-Ce to Hf-Rb}, 1990

_aflow_title 'CsO Structure'
_aflow_proto 'AB_oI8_71_g_i'
_aflow_params 'a,b/a,c/a,y_{1},z_{2}'
_aflow_params_values '4.322,1.73924109209,1.48773715872,0.25,0.38'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oI8'

_symmetry_space_group_name_H-M "I 2/m 2/m 2/m"
_symmetry_Int_Tables_number 71

_cell_length_a     4.32200
_cell_length_b     7.51700
_cell_length_c     6.43000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 -x,-y,-z
6 -x,y,z
7 x,-y,z
8 x,y,-z
9 x+1/2,y+1/2,z+1/2
10 x+1/2,-y+1/2,-z+1/2
11 -x+1/2,y+1/2,-z+1/2
12 -x+1/2,-y+1/2,z+1/2
13 -x+1/2,-y+1/2,-z+1/2
14 -x+1/2,y+1/2,z+1/2
15 x+1/2,-y+1/2,z+1/2
16 x+1/2,y+1/2,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cs1 Cs 4 g 0.00000 0.25000 0.00000 1.00000
O1 O 4 i 0.00000 0.00000 0.38000 1.00000
```

CsO: AB_oI8_71_g_i - POSCAR

```
AB_oI8_71_g_i & a,b/a,c/a,y1,z2 --params=4.322,1.73924109209,
    ↪ 1.48773715872,0.25,0.38 & Immm D_{2h}^{25} #71 (gi) & oI8 &
    ↪ None & CsO & CsO & M. E. Rengade, {C. R. Acad. Sci. C 148,
    ↪ 1199-1202 (1909)
   1.00000000000000
  -2.16100000000000    3.75850000000000    3.21500000000000
   2.16100000000000   -3.75850000000000    3.21500000000000
   2.16100000000000    3.75850000000000   -3.21500000000000
     Cs      O
      2      2
Direct
   0.25000000000000    0.00000000000000    0.25000000000000   Cs   (4g)
  -0.25000000000000    0.00000000000000   -0.25000000000000   Cs   (4g)
   0.38000000000000    0.38000000000000    0.00000000000000   O    (4i)
  -0.38000000000000   -0.38000000000000    0.00000000000000   O    (4i)
```

Ga_2Mg_5 ($D8_g$): A2B5_oI28_72_j_bfj - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Ga2Mg5'
_chemical_formula_sum 'Ga2 Mg5'

loop_
_publ_author_name
 'K. Schubert'
 'K. Frank'
 'R. Gohle'
 'A. Maldonado'
 'H. G. Meissner'
 'A. Raman'
 'W. Rossteutscher'
_journal_name_full_name
;
 Naturwissenschaften
;
_journal_volume 50
_journal_year 1963
_journal_page_first 41
_journal_page_last 41
_publ_Section_title
;
 Einige Strukturdaten metallischer Phasen (8)
;

# Found in Crystal Structure of Ga$_{2}$Mg$_{5}$, 2018 Found in Crystal
    ↪ Structure of Ga$_{2}$Mg$_{5}$, {Crystallography online.com},

_aflow_title 'Ga$_{2}$Mg$_{5}$ ($D8_{g}$) Structure'
_aflow_proto 'A2B5_oI28_72_j_bfj'
_aflow_params 'a,b/a,c/a,x_{2},x_{3},y_{3},x_{4},y_{4}'
_aflow_params_values '13.7,0.511678832117,0.439416058394,0.26,0.378,
    ↪ 0.255,0.27,0.345'
_aflow_Strukturbericht '$D8_{g}$'
_aflow_Pearson 'oI28'

_symmetry_space_group_name_H-M "I 2/b 2/a 2/m"
_symmetry_Int_Tables_number 72

_cell_length_a     13.70000
_cell_length_b     7.01000
_cell_length_c     6.02000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z+1/2
3 -x,y,-z+1/2
4 -x,-y,z
5 -x,-y,-z
6 -x,y,z+1/2
7 x,-y,z+1/2
8 x,y,-z
9 x+1/2,y+1/2,z+1/2
10 x+1/2,-y+1/2,-z
11 -x+1/2,y+1/2,-z
12 -x+1/2,-y+1/2,z+1/2
13 -x+1/2,-y+1/2,-z+1/2
14 -x+1/2,y+1/2,z
15 x+1/2,-y+1/2,z
16 x+1/2,y+1/2,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Mg1 Mg 4 b 0.50000 0.00000 0.25000 1.00000
Mg2 Mg 8 f 0.26000 0.00000 0.25000 1.00000
```

```
Ga1  Ga  8  j  0.37800  0.25500  0.00000  1.00000
Mg3  Mg  8  j  0.27000  0.34500  0.00000  1.00000
```

Ga_2Mg_5 ($D8_g$): A2B5_oI28_72_j_bfj - POSCAR

```
A2B5_oI28_72_j_bfj & a,b/a,c/a,x2,x3,y3,x4,y4 --params=13.7,
      ↪ 0.511678832117,0.439416058394,0.26,0.378,0.255,0.27,0.345 &
      ↪ Ibam D_{2h}^{26} #72 (bfj^2) & oI28 & $D8_{g}$ & Ga2Mg5 &
      ↪ Ga2Mg5 & K. Schubert et al., Naturwissenschaften 50, 41(1963)
   1.00000000000000
  -6.85000000000000    3.50500000000000    3.01000000000000
   6.85000000000000   -3.50500000000000    3.01000000000000
   6.85000000000000    3.50500000000000   -3.01000000000000
     Ga      Mg
      4      10
Direct
   0.25500000000000    0.37800000000000    0.63300000000000   Ga    (8j)
  -0.25500000000000   -0.37800000000000   -0.63300000000000   Ga    (8j)
   0.75500000000000    0.12200000000000   -0.12300000000000   Ga    (8j)
   0.24500000000000    0.87800000000000    0.12300000000000   Ga    (8j)
   0.25000000000000    0.75000000000000    0.50000000000000   Mg    (4b)
   0.75000000000000    0.25000000000000    0.50000000000000   Mg    (4b)
   0.25000000000000    0.51000000000000    0.26000000000000   Mg    (8f)
   0.25000000000000   -0.01000000000000   -0.26000000000000   Mg    (8f)
   0.75000000000000    0.49000000000000   -0.26000000000000   Mg    (8f)
   0.75000000000000    1.01000000000000    0.26000000000000   Mg    (8f)
   0.34500000000000    0.27000000000000    0.61500000000000   Mg    (8j)
  -0.34500000000000   -0.27000000000000   -0.61500000000000   Mg    (8j)
   0.84500000000000    0.23000000000000    0.07500000000000   Mg    (8j)
   0.15500000000000    0.77000000000000   -0.07500000000000   Mg    (8j)
```

$Zn(NH_3)_2Cl_2$ ($E1_2$): A2B6C2D_oI44_74_h_ij_i_e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Cl2H6N2Zn'
_chemical_formula_sum 'C12 H6 N2 Zn'

loop_
_publ_author_name
 'T. Iv\v{s}i\\'{c}'
 'D. W. Bi'
 'A. Magrez'
_journal_name_full_name
;
 Acta Crystallographica Section E: Crystallographic Communications
;
_journal_volume 75
_journal_year 2019
_journal_page_first 1386
_journal_page_last 1388
_publ_Section_title
;
 New refinement of the crystal structure of Zn(NH$_{3}$)$_{2}$Cl$_{2}$
      ↪ at 100K
;

_aflow_title 'Zn(NH$_{3}$)$_{2}$Cl$_{2}$ ($E1_{2}$) Structure'
_aflow_proto 'A2B6C2D_oI44_74_h_ij_i_e'
_aflow_params 'a,b/a,c/a,z_{1},y_{2},z_{2},x_{3},z_{3},x_{4},z_{4},x_{5
      ↪ },y_{5},z_{5}'
_aflow_params_values '7.7077,1.04085524865,1.09664361612,0.88832,0.47954
      ↪ ,0.73085,0.206,0.533,0.28237,0.47949,0.274,0.3281,0.416'
_aflow_Strukturbericht '$E1_{2}$'
_aflow_Pearson 'oI44'

_symmetry_space_group_name_H-M "I 21/m 21/m 21/a"
_symmetry_Int_Tables_number 74

_cell_length_a     7.70770
_cell_length_b     8.02260
_cell_length_c     8.45260
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y+1/2,-z
4 -x,-y+1/2,z
5 -x,-y,-z
6 -x,y,z
7 x,-y+1/2,z
8 x,y+1/2,-z
9 x+1/2,y+1/2,z+1/2
10 x+1/2,-y+1/2,-z+1/2
11 -x+1/2,y,-z+1/2
12 -x+1/2,-y,z+1/2
13 -x+1/2,-y+1/2,-z+1/2
14 -x+1/2,y+1/2,z+1/2
15 x+1/2,-y,z+1/2
16 x+1/2,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Zn1  Zn  4  e  0.00000  0.25000  0.88832  1.00000
Cl1  Cl  8  h  0.00000  0.47954  0.73085  1.00000
H1  H  8  i  0.20600  0.25000  0.53300  1.00000
N1  N  8  i  0.28237  0.25000  0.47949  1.00000
H2  H  16  j  0.27400  0.32810  0.41600  1.00000
```

$Zn(NH_3)_2Cl_2$ ($E1_2$): A2B6C2D_oI44_74_h_ij_i_e - POSCAR

```
A2B6C2D_oI44_74_h_ij_i_e & a,b/a,c/a,z1,y2,z2,x3,z3,x4,z4,x5,y5,z5 --
      ↪ params=7.7077,1.04085524865,1.09664361612,0.88832,0.47954,
      ↪ 0.73085,0.206,0.533,0.28237,0.47949,0.274,0.3281,0.416 & Imma
      ↪ D_{2h}^{28} #74 (ehi^2j) & oI44 & $E1_{2}$ & Cl2H6N2Zn &
      ↪ Cl2H6N2Zn & T. Iv\v{s}i\'{c} and D. W. Bi and A. Magrez, Acta
      ↪ Crystallogr. E 75, 1386-1388 (2019)
   1.00000000000000
  -3.85385000000000    4.01130000000000    4.22630000000000
   3.85385000000000   -4.01130000000000    4.22630000000000
   3.85385000000000    4.01130000000000   -4.22630000000000
     Cl      H      N      Zn
      4     12      4       2
Direct
   1.21039000000000    0.73085000000000    0.47954000000000   Cl    (8h)
   0.75131000000000    0.73085000000000    0.02046000000000   Cl    (8h)
   0.24869000000000   -0.73085000000000    0.97954000000000   Cl    (8h)
  -1.21039000000000   -0.73085000000000   -0.47954000000000   Cl    (8h)
   0.78300000000000    0.73900000000000    0.45600000000000   H     (8i)
   0.78300000000000    0.32700000000000    0.04400000000000   H     (8i)
   0.21700000000000   -0.73900000000000    0.54400000000000   H     (8i)
   0.21700000000000   -0.32700000000000    0.95600000000000   H     (8i)
   0.74410000000000    0.69000000000000    0.60210000000000   H    (16j)
   0.58790000000000    0.14200000000000   -0.10210000000000   H    (16j)
   0.41210000000000   -0.69000000000000    0.55410000000000   H    (16j)
  -0.74410000000000   -0.14200000000000   -0.05410000000000   H    (16j)
  -0.74410000000000   -0.69000000000000   -0.60210000000000   H    (16j)
   0.41210000000000   -0.14200000000000    1.10210000000000   H    (16j)
   0.58790000000000    0.69000000000000    0.44590000000000   H    (16j)
   0.74410000000000    0.14200000000000    0.05410000000000   H    (16j)
   0.72949000000000    0.76186000000000    0.53237000000000   N     (8i)
   0.72949000000000    0.19712000000000   -0.03237000000000   N     (8i)
   0.27051000000000   -0.76186000000000    0.46763000000000   N     (8i)
   0.27051000000000   -0.19712000000000    1.03237000000000   N     (8i)
   1.13832000000000    0.88832000000000    0.25000000000000   Zn    (4e)
  -0.13832000000000   -0.88832000000000    0.75000000000000   Zn    (4e)
```

$CeCu_2$: AB2_oI12_74_e_h - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'CeCu2'
_chemical_formula_sum 'Ce Cu2'

loop_
_publ_author_name
 'A. C. Larson'
 'D. T. Cromer'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 14
_journal_year 1961
_journal_page_first 73
_journal_page_last 74
_publ_Section_title
;
 The crystal structure of CeCu$_{2}$
;

_aflow_title 'CeCu$_{2}$ Structure'
_aflow_proto 'AB2_oI12_74_e_h'
_aflow_params 'a,b/a,c/a,z_{1},y_{2},z_{2}'
_aflow_params_values '4.43,1.5914221219,1.68171557562,0.5377,0.051,
      ↪ 0.1648'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oI12'

_symmetry_space_group_name_H-M "I 21/m 21/m 21/a"
_symmetry_Int_Tables_number 74

_cell_length_a     4.43000
_cell_length_b     7.05000
_cell_length_c     7.45000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y+1/2,-z
4 -x,-y+1/2,z
5 -x,-y,-z
6 -x,y,z
7 x,-y+1/2,z
8 x,y+1/2,-z
9 x+1/2,y+1/2,z+1/2
10 x+1/2,-y+1/2,-z+1/2
11 -x+1/2,y,-z+1/2
12 -x+1/2,-y,z+1/2
```

```
13 -x+1/2,-y+1/2,-z+1/2
14 -x+1/2,y+1/2,z+1/2
15 x+1/2,-y,z+1/2
16 x+1/2,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ce1 Ce 4 e 0.00000 0.25000 0.53770 1.00000
Cu1 Cu 8 h 0.00000 0.05100 0.16480 1.00000
```

$CeCu_2$: AB2_oI12_74_e_h - POSCAR

```
AB2_oI12_74_e_h & a,b/a,c/a,z1,y2,z2 --params=4.43,1.5914221219,
    ↪ 1.68171557562,0.5377,0.051,0.1648 & Imma D_{2h}^{28} #74 (eh) &
    ↪  oI12 & None & CeCu2 & CeCu2 & A. C. Larson and D. T. Cromer,
    ↪ Acta Cryst. 14, 73-74 (1961)
   1.00000000000000
  -2.21500000000000    3.52500000000000    3.72500000000000
   2.21500000000000   -3.52500000000000    3.72500000000000
   2.21500000000000    3.52500000000000   -3.72500000000000
    Ce     Cu
     2      4
Direct
   0.78770000000000    0.53770000000000    0.25000000000000   Ce   (4e)
   0.21230000000000   -0.53770000000000    0.75000000000000   Ce   (4e)
   0.21580000000000    0.16480000000000    0.05100000000000   Cu   (8h)
   0.61380000000000    0.16480000000000    0.44900000000000   Cu   (8h)
   0.38620000000000   -0.16480000000000    0.55100000000000   Cu   (8h)
  -0.21580000000000   -0.16480000000000   -0.05100000000000   Cu   (8h)
```

$LiCuVO_4$: ABC4D_oI28_74_a_d_hi_e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'CuLiO4V'
_chemical_formula_sum 'Cu Li O4 V'

loop_
_publ_author_name
 'M. A. Lafontaine'
 'M. Leblanc'
 'G. Ferey'
_journal_name_full_name
;
 Acta Crystallographica Section C: Structural Chemistry
;
_journal_volume 45
_journal_year 1989
_journal_page_first 1205
_journal_page_last 1206
_publ_Section_title
;
 New refinement of the room-temperature structure of LiCuVO${4}$
;

# Found in Structure, stoichiometry and magnetic properties of the
    ↪ low-dimensional structure phase LiCuVO${4}$, 2004

_aflow_title 'LiCuVO${4}$ Structure'
_aflow_proto 'ABC4D_oI28_74_a_d_hi_e'
_aflow_params 'a,b/a,c/a,z_{3},y_{4},z_{4},x_{5},z_{5}'
_aflow_params_values '5.662,1.0259625574,1.54680324974,0.386,0.0164,
    ↪ 0.2748,0.2352,-0.0007'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'oI28'

_symmetry_space_group_name_H-M "I 21/m 21/m 21/a"
_symmetry_Int_Tables_number 74

_cell_length_a      5.66200
_cell_length_b      5.80900
_cell_length_c      8.75800
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y+1/2,-z
4 -x,-y+1/2,z
5 -x,-y,-z
6 -x,y,z
7 x,-y+1/2,z
8 x,y+1/2,-z
9 x+1/2,y+1/2,z+1/2
10 x+1/2,-y+1/2,-z+1/2
11 -x+1/2,y,-z+1/2
12 -x+1/2,-y,z+1/2
13 -x+1/2,-y+1/2,-z+1/2
14 -x+1/2,y+1/2,z+1/2
15 x+1/2,-y,z+1/2
16 x+1/2,y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cu1 Cu 4 a 0.00000 0.00000 0.00000 1.00000
Li1 Li 4 d 0.25000 0.25000 0.75000 1.00000
V1 V 4 e 0.00000 0.25000 0.38600 1.00000
O1 O 8 h 0.00000 0.01640 0.27480 1.00000
O2 O 8 i 0.23520 0.25000 -0.00070 1.00000
```

$LiCuVO_4$: ABC4D_oI28_74_a_d_hi_e - POSCAR

```
ABC4D_oI28_74_a_d_hi_e & a,b/a,c/a,z3,y4,z4,x5,z5 --params=5.662,
    ↪ 1.0259625574,1.54680324974,0.386,0.0164,0.2748,0.2352,-0.0007 &
    ↪  Imma D_{2h}^{28} #74 (adehi) & oI28 & None & CuLiO4V & CuLiO4V
    ↪  & M. A. Lafontaine and M. Leblanc and G. Ferey, Acta
    ↪ Crystallogr. C 45, 1205-1206 (1989)
   1.00000000000000
  -2.83100000000000    2.90450000000000    4.37900000000000
   2.83100000000000   -2.90450000000000    4.37900000000000
   2.83100000000000    2.90450000000000   -4.37900000000000
    Cu     Li     O      V
     2      2      8      2
Direct
   0.00000000000000    0.00000000000000    0.00000000000000   Cu   (4a)
   0.50000000000000    0.00000000000000    0.50000000000000   Cu   (4a)
   0.00000000000000    0.00000000000000    0.50000000000000   Li   (4d)
   0.00000000000000    0.50000000000000    0.00000000000000   Li   (4d)
   0.29120000000000    0.27480000000000    0.01640000000000   O    (8h)
   0.75840000000000    0.27480000000000    0.48360000000000   O    (8h)
   0.24160000000000   -0.27480000000000    0.51640000000000   O    (8h)
  -0.29120000000000   -0.27480000000000   -0.01640000000000   O    (8h)
   0.24930000000000    0.23450000000000    0.48520000000000   O    (8i)
   0.24930000000000   -0.23590000000000    0.01480000000000   O    (8i)
   0.75070000000000   -0.23450000000000    0.51480000000000   O    (8i)
   0.75070000000000    0.23590000000000    0.98520000000000   O    (8i)
   0.63600000000000    0.38600000000000    0.25000000000000   V    (4e)
   0.36400000000000   -0.38600000000000    0.75000000000000   V    (4e)
```

Gwihabaite [NH_4NO_3 (V)]: A4B2C3_tP72_77_8d_ab2c2d_6d - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Gwihabaite'
_chemical_formula_sum 'H4 N2 O3'

loop_
_publ_author_name
 'J. L. Amor\\'{o}s'
 'F. Arrese'
 'M. Canut'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 117
_journal_year 1962
_journal_page_first 92
_journal_page_last 107
_publ_Section_title
;
 The crystal structure of the low-temperature phase of NH$_{4}$NO$_{3}$
    ↪ (V) at -- 150$^{\circ}$C
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'Gwihabaite [NH$_{4}$NO$_{3}$ (V)] Structure'
_aflow_proto 'A4B2C3_tP72_77_8d_ab2c2d_6d'
_aflow_params 'a,c/a,z_{1},z_{2},z_{3},z_{4},x_{5},y_{5},z_{5},x_{6},y_{
    ↪ 6},z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},
    ↪ x_{10},y_{10},z_{10},x_{11},y_{11},z_{11},x_{12},y_{12},z_{12},
    ↪ x_{13},y_{13},z_{13},x_{14},y_{14},z_{14},x_{15},y_{15},z_{15},
    ↪ x_{16},y_{16},z_{16},x_{17},y_{17},z_{17},x_{18},y_{18},z_{18},
    ↪ x_{19},y_{19},z_{19},x_{20},y_{20},z_{20}'
_aflow_params_values '7.98,1.22556390977,0.25,0.25,0.75,0.25,0.076,0.076
    ↪ ,0.19,0.076,-0.076,0.31,0.576,0.076,0.19,0.424,0.076,0.31,0.424
    ↪ ,0.424,0.31,0.424,0.576,0.19,-0.076,0.424,0.31,-0.076,0.576,
    ↪ 0.19,0.25,0.25,0.518,0.25,0.25,0.018,0.12,0.27,0.456,0.33,0.23,
    ↪ 0.456,0.25,0.25,0.642,0.12,0.23,-0.044,0.38,0.27,-0.044,0.25,
    ↪ 0.25,0.142'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tP72'

_symmetry_space_group_name_H-M "P 42"
_symmetry_Int_Tables_number 77

_cell_length_a      7.98000
_cell_length_b      7.98000
_cell_length_c      9.78000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
```

```
2 -x,-y,z
3 -y,x,z+1/2
4 y,-x,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
N1 N 2 a 0.00000 0.00000 0.25000 1.00000
N2 N 2 b 0.50000 0.50000 0.25000 1.00000
N3 N 2 c 0.00000 0.50000 0.75000 1.00000
N4 N 2 c 0.00000 0.50000 0.25000 1.00000
H1 H 4 d 0.07600 0.07600 0.19000 1.00000
H2 H 4 d 0.07600 -0.07600 0.31000 1.00000
H3 H 4 d 0.57600 0.07600 0.19000 1.00000
H4 H 4 d 0.42400 0.07600 0.31000 1.00000
H5 H 4 d 0.42400 0.42400 0.31000 1.00000
H6 H 4 d 0.42400 0.57600 0.19000 1.00000
H7 H 4 d -0.07600 0.42400 0.31000 1.00000
H8 H 4 d -0.07600 0.57600 0.19000 1.00000
N5 N 4 d 0.25000 0.25000 0.51800 1.00000
N6 N 4 d 0.25000 0.25000 0.01800 1.00000
O1 O 4 d 0.12000 0.27000 0.45600 1.00000
O2 O 4 d 0.33000 0.23000 0.45600 1.00000
O3 O 4 d 0.25000 0.25000 0.64200 1.00000
O4 O 4 d 0.12000 0.23000 -0.04400 1.00000
O5 O 4 d 0.38000 0.27000 -0.04400 1.00000
O6 O 4 d 0.25000 0.25000 0.14200 1.00000
```

Gwihabaite [NH_4NO_3 (V)]: A4B2C3_tP72_77_8d_ab2c2d_6d - POSCAR

```
A4B2C3_tP72_77_8d_ab2c2d_6d & a,c/a,z1,z2,z3,z4,x5,y5,z5,x6,y6,z6,x7,y7,
    ↪ z7,x8,y8,z8,x9,y9,z9,x10,y10,z10,x11,y11,z11,x12,y12,z12,x13,
    ↪ y13,z13,x14,y14,z14,x15,y15,z15,x16,y16,z16,x17,y17,z17,x18,y18
    ↪ ,z18,x19,y19,z19,x20,y20,z20 --params=7.98,1.22556390977,0.25,
    ↪ 0.25,0.75,0.25,0.076,0.076,0.19,0.076,-0.076,0.31,0.576,0.076,
    ↪ 0.19,0.424,0.076,0.31,0.424,0.424,0.31,0.424,0.576,0.19,-0.076,
    ↪ 0.424,0.31,-0.076,0.576,0.19,0.25,0.25,0.518,0.25,0.25,0.018,
    ↪ 0.12,0.27,0.456,0.33,0.23,0.456,0.25,0.25,0.642,0.12,0.23,-
    ↪ 0.044,0.38,0.27,-0.044,0.25,0.25,0.142 & P4_{2} C_{4}^{3} #77 (
    ↪ abc^2d^16) & tP72 & None & H4N2O3 & Gwihabaite & J. L. Amor\'{o
    ↪ }s and F. Arrese and M. Canut, Zeitschrift f{"u}r
    ↪ Kristallographie - Crystalline Materials 117, 92-107 (1962)
   1.00000000000000
   7.98000000000000    0.00000000000000    0.00000000000000
   0.00000000000000    7.98000000000000    0.00000000000000
   0.00000000000000    0.00000000000000    9.78000000000000
     H     N     O
    32    16    24
Direct
   0.07600000000000    0.07600000000000    0.19000000000000  H   (4d)
  -0.07600000000000   -0.07600000000000    0.19000000000000  H   (4d)
  -0.07600000000000    0.07600000000000    0.69000000000000  H   (4d)
   0.07600000000000   -0.07600000000000    0.69000000000000  H   (4d)
   0.07600000000000   -0.07600000000000    0.31000000000000  H   (4d)
  -0.07600000000000    0.07600000000000    0.31000000000000  H   (4d)
   0.07600000000000    0.07600000000000    0.81000000000000  H   (4d)
  -0.07600000000000   -0.07600000000000    0.81000000000000  H   (4d)
   0.57600000000000    0.07600000000000    0.19000000000000  H   (4d)
  -0.57600000000000   -0.07600000000000    0.19000000000000  H   (4d)
  -0.07600000000000    0.57600000000000    0.69000000000000  H   (4d)
   0.07600000000000   -0.57600000000000    0.69000000000000  H   (4d)
   0.42400000000000    0.07600000000000    0.31000000000000  H   (4d)
  -0.42400000000000   -0.07600000000000    0.31000000000000  H   (4d)
  -0.07600000000000    0.42400000000000    0.81000000000000  H   (4d)
   0.07600000000000   -0.42400000000000    0.81000000000000  H   (4d)
   0.42400000000000    0.42400000000000    0.31000000000000  H   (4d)
  -0.42400000000000   -0.42400000000000    0.31000000000000  H   (4d)
  -0.42400000000000    0.42400000000000    0.81000000000000  H   (4d)
   0.42400000000000   -0.42400000000000    0.81000000000000  H   (4d)
   0.42400000000000    0.57600000000000    0.19000000000000  H   (4d)
  -0.42400000000000   -0.57600000000000    0.19000000000000  H   (4d)
  -0.57600000000000    0.42400000000000    0.69000000000000  H   (4d)
   0.57600000000000   -0.42400000000000    0.69000000000000  H   (4d)
  -0.07600000000000    0.42400000000000    0.31000000000000  H   (4d)
   0.07600000000000   -0.42400000000000    0.31000000000000  H   (4d)
  -0.42400000000000   -0.07600000000000    0.81000000000000  H   (4d)
   0.42400000000000    0.07600000000000    0.81000000000000  H   (4d)
  -0.07600000000000    0.57600000000000    0.19000000000000  H   (4d)
   0.07600000000000   -0.57600000000000    0.19000000000000  H   (4d)
  -0.57600000000000   -0.07600000000000    0.69000000000000  H   (4d)
   0.57600000000000    0.07600000000000    0.69000000000000  H   (4d)
   0.00000000000000    0.00000000000000    0.25000000000000  N   (2a)
   0.00000000000000    0.00000000000000    0.75000000000000  N   (2a)
   0.50000000000000    0.50000000000000    0.25000000000000  N   (2b)
   0.50000000000000    0.50000000000000    0.75000000000000  N   (2b)
   0.00000000000000    0.50000000000000    0.75000000000000  N   (2c)
   0.50000000000000    0.00000000000000    1.25000000000000  N   (2c)
   0.00000000000000    0.50000000000000    0.25000000000000  N   (2c)
   0.50000000000000    0.00000000000000    0.75000000000000  N   (2c)
   0.25000000000000    0.25000000000000    0.51800000000000  N   (4d)
  -0.25000000000000   -0.25000000000000    0.51800000000000  N   (4d)
  -0.25000000000000    0.25000000000000    1.01800000000000  N   (4d)
   0.25000000000000   -0.25000000000000    1.01800000000000  N   (4d)
   0.25000000000000    0.25000000000000    0.01800000000000  N   (4d)
  -0.25000000000000   -0.25000000000000    0.01800000000000  N   (4d)
  -0.25000000000000    0.25000000000000    0.51800000000000  N   (4d)
   0.25000000000000   -0.25000000000000    0.51800000000000  N   (4d)
   0.12000000000000    0.27000000000000    0.45600000000000  O   (4d)
  -0.12000000000000   -0.27000000000000    0.45600000000000  O   (4d)
  -0.27000000000000    0.12000000000000    0.95600000000000  O   (4d)
   0.27000000000000   -0.12000000000000    0.95600000000000  O   (4d)
   0.33000000000000    0.23000000000000    0.45600000000000  O   (4d)
  -0.33000000000000   -0.23000000000000    0.45600000000000  O   (4d)
  -0.23000000000000    0.33000000000000    0.95600000000000  O   (4d)
   0.23000000000000   -0.33000000000000    0.95600000000000  O   (4d)
   0.25000000000000    0.25000000000000    0.64200000000000  O   (4d)
  -0.25000000000000   -0.25000000000000    0.64200000000000  O   (4d)
  -0.25000000000000    0.25000000000000    1.14200000000000  O   (4d)
   0.25000000000000   -0.25000000000000    1.14200000000000  O   (4d)
   0.12000000000000    0.23000000000000   -0.04400000000000  O   (4d)
  -0.12000000000000   -0.23000000000000   -0.04400000000000  O   (4d)
  -0.23000000000000    0.12000000000000    0.45600000000000  O   (4d)
   0.23000000000000   -0.12000000000000    0.45600000000000  O   (4d)
   0.38000000000000    0.27000000000000   -0.04400000000000  O   (4d)
  -0.38000000000000   -0.27000000000000   -0.04400000000000  O   (4d)
  -0.27000000000000    0.38000000000000    0.45600000000000  O   (4d)
   0.27000000000000   -0.38000000000000    0.45600000000000  O   (4d)
   0.25000000000000    0.25000000000000    0.14200000000000  O   (4d)
  -0.25000000000000   -0.25000000000000    0.14200000000000  O   (4d)
  -0.25000000000000    0.25000000000000    0.64200000000000  O   (4d)
   0.25000000000000   -0.25000000000000    0.64200000000000  O   (4d)
```

Kesterite [$Cu_2(Zn,Fe)SnS_4$]: A2BCD4_tI16_82_ac_b_d_g - CIF

```
# AFLOW.org Repositories
# CuFeSSn/A2BCD4_tI16_82_ac_b_d_g-001.ABCD params=-1,2.00313248572,
    ↪ 0.7434,0.256,0.6278 SG=82 [ANRL doi: 10.1016/
    ↪ j.commatsci.2017.01.017 (part 1), doi: 10.1016/
    ↪ j.commatsci.2018.10.043 (part 2)]
data_CuFeSSn
_pd_phase_name A2BCD4_tI16_82_ac_b_d_g-001.ABCD

_chemical_name_mineral 'Kesterite'
_chemical_formula_sum 'Cu2 Fe S Sn4'

loop_
_publ_author_name
 'S. R. Hall'
 'J. T. Szyma{\\'n}ski'
 'J. M. Stewart'
_journal_name_full_name
;
 Canadian Mineralogist
;
_journal_volume 16
_journal_year 1978
_journal_page_first 131
_journal_page_last 137
_publ_Section_title
;
 Kesterite, Cu$_{2}$(Zn,Fe)SnS$_{4}$, and stannite, Cu$_{2}$(Fe,Zn)SnS$_
    ↪ {4}$, structurally similar but distinct minerals
;

_aflow_title 'Kesterite [Cu$_{2}$(Zn,Fe)SnS$_{4}$] Structure'
_aflow_proto 'A2BCD4_tI16_82_ac_b_d_g'
_aflow_params 'a,c/a,x_{5},y_{5},z_{5}'
_aflow_params_values '5.428437725,2.00313248571,0.7434,0.256,0.6278'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tI16'

_cell_length_a  5.4284377250
_cell_length_b  5.4284377250
_cell_length_c  10.8738799536
_cell_angle_alpha  90.0000000000
_cell_angle_beta  90.0000000000
_cell_angle_gamma  90.0000000000
_symmetry_space_group_name_H-M  'I-4'
_symmetry_Int_Tables_Number  82
loop_
_symmetry_equiv_pos_site_id
_symmetry_equiv_pos_as_xyz
  1  x,y,z
  2  -x,-y,z
  3  y,-x,-z
  4  -y,x,-z
  5  x+1/2,y+1/2,z+1/2
  6  -x+1/2,-y+1/2,z+1/2
  7  y+1/2,-x+1/2,-z+1/2
  8  -y+1/2,x+1/2,-z+1/2
loop_
 _atom_site_label
 _atom_site_occupancy
 _atom_site_fract_x
 _atom_site_fract_y
 _atom_site_fract_z
 _atom_site_thermal_displace_type
 _atom_site_B_iso_or_equiv
 _atom_site_type_symbol
 _atom_site_symmetry_multiplicity
 _atom_site_Wyckoff_label
Cu1 1.0000000000 -0.0000000000 0.0000000000 0.0000000000 Biso 1.0 Cu 2 a
Fe1 1.0000000000 0.0000000000 0.0000000000 0.5000000000 Biso 1.0 Fe 2 b
Cu2 1.0000000000 -0.0000000000 0.5000000000 0.2500000000 Biso 1.0 Cu 2 c
S1 1.0000000000 0.0000000000 0.5000000000 0.7500000000 Biso 1.0 S 2 d
Sn1 1.0000000000 0.7434000000 0.2560000000 0.6278000000 Biso 1.0 Sn 8 g
```

Kesterite [$Cu_2(Zn,Fe)SnS_4$]: A2BCD4_tI16_82_ac_b_d_g - POSCAR

```
A2BCD4_tI16_82_ac_b_d_g & a,c/a,x5,y5,z5 --params=5.428437725,
    ↪ 2.00313248571,0.7434,0.256,0.6278 & I-4 S_{4}^{2} #82 (abcdg) &
    ↪  tI16 & None & Cu2(Fe,Zn)SnS4 & Kesterite & S. R. Hall and J.
    ↪ T. Szyma{\'n}ski and J. M. Stewart, Can. Mineral. 16, 131-137 (
    ↪ 1978)
```

```
   1.00000000000000
  -2.71421886250000    2.71421886250000    5.43693997680000
   2.71421886250000   -2.71421886250000    5.43693997680000
   2.71421886250000    2.71421886250000   -5.43693997680000
     Cu     Fe     S     Sn
      2      1      1      4
Direct
   0.00000000000000    0.00000000000000    0.00000000000000   Cu   (2a)
   0.75000000000000    0.25000000000000    0.50000000000000   Cu   (2c)
   0.50000000000000    0.50000000000000    0.00000000000000   Fe   (2b)
   0.25000000000000    0.75000000000000    0.50000000000000    S   (2d)
   0.88380000000000    1.37120000000000    0.99940000000000   Sn   (8g)
   0.37180000000000   -0.11560000000000   -0.99940000000000   Sn   (8g)
  -1.37120000000000   -0.37180000000000   -0.48740000000000   Sn   (8g)
   0.11560000000000   -0.88380000000000    0.48740000000000   Sn   (8g)
```

Bromocarnallite ($KMg(H_2O)_6(Cl,Br)_3$, $E2_6$): A3B6CD_tP44_85_bcg_3g_ac_e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Bromocarnallite'
_chemical_formula_sum 'Br3 (H2O)6 K Mg'

loop_
_publ_author_name
 'K. R. Andre{\ss}'
 'O. Saffe'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 101
_journal_year 1939
_journal_page_first 451
_journal_page_last 469
_publ_Section_title
;
 R\"{o}ntgenographische Untersuchung der Mischkristallreihe
       ↪ Karnallit-Bromkarnallit
;

# Found in Strukturbericht Band VII 1939, 1943

_aflow_title 'Bromocarnallite (KMg(H$_{2}$O)$_{6}$(Cl,Br)$_{3}$, $E2_{6}
      ↪ $) Structure'
_aflow_proto 'A3B6CD_tP44_85_bcg_3g_ac_e'
_aflow_params 'a,c/a,z_{3},z_{4},x_{6},y_{6},z_{6},x_{7},y_{7},z_{7},x_{
      ↪ 8},y_{8},z_{8},x_{9},y_{9},z_{9}'
_aflow_params_values '13.51,0.501110288675,0.58,0.08,0.25,0.0,0.008,0.55
      ↪ ,0.05,0.78,0.54,0.85,0.58,0.34,-0.03,0.58'
_aflow_Strukturbericht '$E2_{6}$'
_aflow_Pearson 'tP44'

_symmetry_space_group_name_H-M "P 4/n (origin choice 2)"
_symmetry_Int_Tables_number 85

_cell_length_a     13.51000
_cell_length_b     13.51000
_cell_length_c     6.77000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x+1/2,-y+1/2,z
3 -y+1/2,x,z
4 y,-x+1/2,z
5 -x,-y,-z
6 x+1/2,y+1/2,-z
7 y+1/2,-x,-z
8 -y,x+1/2,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
K1 K 2 a 0.25000 0.75000 0.00000 1.00000
Br1 Br 2 b 0.25000 0.75000 0.50000 1.00000
Br2 Br 2 c 0.25000 0.25000 0.58000 1.00000
K2 K 2 c 0.25000 0.25000 0.08000 1.00000
Mg1 Mg 4 e 0.00000 0.00000 0.50000 1.00000
Br3 Br 8 g 0.25000 0.00000 0.00800 1.00000
H2O1 H2O 8 g 0.55000 0.05000 0.78000 1.00000
H2O2 H2O 8 g 0.54000 0.85000 0.58000 1.00000
H2O3 H2O 8 g 0.34000 -0.03000 0.58000 1.00000
```

Bromocarnallite ($KMg(H_2O)_6(Cl,Br)_3$, $E2_6$): A3B6CD_tP44_85_bcg_3g_ac_e - POSCAR

```
A3B6CD_tP44_85_bcg_3g_ac_e & a,c/a,z3,z4,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,
      ↪ y9,z9 --params=13.51,0.501110288675,0.58,0.08,0.25,0.0,0.008,
      ↪ 0.55,0.05,0.78,0.54,0.85,0.58,0.34,-0.03,0.58 & P4/n C_{4h}^{3}
      ↪  #85 (abc^2eg^4) & tP44 & $E2_{6}$ & (Br,Cl)3(H2O)6KMg &
      ↪ Bromocarnallite & K. R. Andre{\ss} and O. Saffe, Zeitschrift f
      ↪ {"u}r Kristallographie - Crystalline Materials 101, 451-469 (
      ↪ 1939)
   1.00000000000000
  13.51000000000000    0.00000000000000    0.00000000000000
   0.00000000000000   13.51000000000000    0.00000000000000
   0.00000000000000    0.00000000000000    6.77000000000000
     Br    H2O     K    Mg
     12     24     4     4
Direct
   0.25000000000000    0.75000000000000    0.50000000000000   Br   (2b)
   0.75000000000000    0.25000000000000    0.50000000000000   Br   (2b)
   0.25000000000000    0.25000000000000    0.58000000000000   Br   (2c)
   0.75000000000000    0.75000000000000   -0.58000000000000   Br   (2c)
   0.25000000000000    0.00000000000000    0.00800000000000   Br   (8g)
   0.25000000000000    0.50000000000000    0.00800000000000   Br   (8g)
   0.50000000000000    0.25000000000000    0.00800000000000   Br   (8g)
   0.00000000000000    0.25000000000000    0.00800000000000   Br   (8g)
  -0.25000000000000    0.00000000000000   -0.00800000000000   Br   (8g)
   0.75000000000000    0.50000000000000   -0.00800000000000   Br   (8g)
   0.50000000000000   -0.25000000000000   -0.00800000000000   Br   (8g)
   0.00000000000000    0.75000000000000   -0.00800000000000   Br   (8g)
   0.55000000000000    0.05000000000000    0.78000000000000   H2O  (8g)
  -0.05000000000000    0.45000000000000    0.78000000000000   H2O  (8g)
   0.45000000000000    0.55000000000000    0.78000000000000   H2O  (8g)
   0.05000000000000   -0.05000000000000    0.78000000000000   H2O  (8g)
  -0.55000000000000   -0.05000000000000   -0.78000000000000   H2O  (8g)
   1.05000000000000    0.55000000000000   -0.78000000000000   H2O  (8g)
   0.55000000000000   -0.55000000000000   -0.78000000000000   H2O  (8g)
  -0.05000000000000    1.05000000000000   -0.78000000000000   H2O  (8g)
   0.54000000000000    0.85000000000000    0.58000000000000   H2O  (8g)
  -0.04000000000000   -0.35000000000000    0.58000000000000   H2O  (8g)
  -0.35000000000000    0.54000000000000    0.58000000000000   H2O  (8g)
   0.85000000000000   -0.04000000000000    0.58000000000000   H2O  (8g)
  -0.54000000000000   -0.85000000000000   -0.58000000000000   H2O  (8g)
   1.04000000000000    1.35000000000000   -0.58000000000000   H2O  (8g)
   1.35000000000000   -0.54000000000000   -0.58000000000000   H2O  (8g)
  -0.85000000000000    1.04000000000000   -0.58000000000000   H2O  (8g)
   0.34000000000000   -0.03000000000000    0.58000000000000   H2O  (8g)
   0.16000000000000    0.53000000000000    0.58000000000000   H2O  (8g)
   0.53000000000000    0.34000000000000    0.58000000000000   H2O  (8g)
  -0.03000000000000    0.16000000000000    0.58000000000000   H2O  (8g)
  -0.34000000000000    0.03000000000000   -0.58000000000000   H2O  (8g)
   0.84000000000000    0.47000000000000   -0.58000000000000   H2O  (8g)
   0.47000000000000   -0.34000000000000   -0.58000000000000   H2O  (8g)
   0.03000000000000    0.84000000000000   -0.58000000000000   H2O  (8g)
   0.25000000000000    0.75000000000000    0.00000000000000    K   (2a)
   0.75000000000000    0.25000000000000    0.00000000000000    K   (2a)
   0.25000000000000    0.25000000000000    0.08000000000000    K   (2c)
   0.75000000000000    0.75000000000000   -0.08000000000000    K   (2c)
   0.00000000000000    0.00000000000000    0.50000000000000   Mg   (4e)
   0.50000000000000    0.50000000000000    0.50000000000000   Mg   (4e)
   0.50000000000000    0.00000000000000    0.50000000000000   Mg   (4e)
   0.00000000000000    0.50000000000000    0.50000000000000   Mg   (4e)
```

$MoPO_5$: AB5C_tP14_85_c_cg_b - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'MoO5P'
_chemical_formula_sum 'Mo O5 P'

loop_
_publ_author_name
 'P. Kierkegaard'
 'M. Westerlund'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 18
_journal_year 1964
_journal_page_first 2217
_journal_page_last 2225
_publ_Section_title
;
 The Crystal Structure of MoOPO$_{4}$
;

_aflow_title 'MoPO$_{5}$ Structure'
_aflow_proto 'AB5C_tP14_85_c_cg_b'
_aflow_params 'a,c/a,z_{2},z_{3},x_{4},y_{4},z_{4}'
_aflow_params_values '6.1768,0.695052454345,0.1975,0.8102,0.3125,-0.0554
      ↪ ,0.2994'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tP14'

_symmetry_space_group_name_H-M "P 4/n (origin choice 2)"
_symmetry_Int_Tables_number 85

_cell_length_a     6.17680
_cell_length_b     6.17680
_cell_length_c     4.29320
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x+1/2,-y+1/2,z
3 -y+1/2,x,z
4 y,-x+1/2,z
5 -x,-y,-z
6 x+1/2,y+1/2,-z
```

```
7 y+1/2,-x,-z
8 -y,x+1/2,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
P1 P 2 b 0.25000 0.75000 0.50000 1.00000
Mo1 Mo 2 c 0.25000 0.25000 0.19750 1.00000
O1 O 2 c 0.25000 0.25000 0.81020 1.00000
O2 O 8 g 0.31250 -0.05540 0.29940 1.00000
```

$MoPO_5$: AB5C_tP14_85_c_cg_b - POSCAR

```
AB5C_tP14_85_c_cg_b & a,c/a,z2,z3,x4,y4,z4 --params=6.1768,
    ↪ 0.695052454345,0.1975,0.8102,0.3125,-0.0554,0.2994 & P4/n C_{4h
    ↪ }^{3} #85 (bc^2g) & tP14 & None & MoO5P & MoO5P & P.
    ↪ Kierkegaard and M. Westerlund, Acta Chem. Scand. 18, 2217-2225
    ↪ (1964)
   1.00000000000000
   6.17680000000000   0.00000000000000   0.00000000000000
   0.00000000000000   6.17680000000000   0.00000000000000
   0.00000000000000   0.00000000000000   4.29320000000000
    Mo    O    P
     2   10    2
Direct
   0.25000000000000   0.25000000000000   0.19750000000000  Mo   (2c)
   0.75000000000000   0.75000000000000  -0.19750000000000  Mo   (2c)
   0.25000000000000   0.25000000000000   0.81020000000000  O    (2c)
   0.75000000000000   0.75000000000000  -0.81020000000000  O    (2c)
   0.31250000000000  -0.05540000000000   0.29940000000000  O    (8g)
   0.18750000000000   0.55540000000000   0.29940000000000  O    (8g)
   0.55540000000000   0.31250000000000   0.29940000000000  O    (8g)
  -0.05540000000000   0.18750000000000   0.29940000000000  O    (8g)
  -0.31250000000000   0.05540000000000  -0.29940000000000  O    (8g)
   0.81250000000000   0.44460000000000  -0.29940000000000  O    (8g)
   0.44460000000000  -0.31250000000000  -0.29940000000000  O    (8g)
   0.05540000000000   0.81250000000000  -0.29940000000000  O    (8g)
   0.25000000000000   0.75000000000000   0.50000000000000  P    (2b)
   0.75000000000000   0.25000000000000   0.50000000000000  P    (2b)
```

$PNCl_2$ ($E1_4$): A2BC_tP32_86_2g_g_g - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Cl2NP'
_chemical_formula_sum 'Cl2 N P'

loop_
_publ_author_name
 'J. A. A. Ketelaar'
 'T. A. {de Vries}'
_journal_name_full_name
;
 Recueil des Travaux Chimiques des Pays-Bas
;
_journal_volume 58
_journal_year 1939
_journal_page_first 1081
_journal_page_last 1099
_publ_Section_title
;
 The crystal structure of tetra phosphonitrile chloride, P$_{4}$N$_{4}
    ↪ $Cl$_{8}$
;

_aflow_title 'PNCl$_{2}$ ($E1_{4}$) Structure'
_aflow_proto 'A2BC_tP32_86_2g_g_g'
_aflow_params 'a,c/a,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3},y_{3},z_{
    ↪ 3},x_{4},y_{4},z_{4}'
_aflow_params_values '10.82,0.549907578558,0.551,0.852,0.86,0.594,0.198,
    ↪ 0.69,0.625,0.645,0.14,0.574,0.791,0.17'
_aflow_Strukturbericht '$E1_{4}$'
_aflow_Pearson 'tP32'

_symmetry_space_group_name_H-M "P 42/n (origin choice 2)"
_symmetry_Int_Tables_number 86

_cell_length_a    10.82000
_cell_length_b    10.82000
_cell_length_c    5.95000
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x+1/2,-y+1/2,z
3 -y,x+1/2,z+1/2
4 y+1/2,-x,z+1/2
5 -x,-y,-z
6 x+1/2,y+1/2,-z
7 y,-x+1/2,-z+1/2
8 -y+1/2,x,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cl1 Cl 8 g 0.55100 0.85200 0.86000 1.00000
Cl2 Cl 8 g 0.59400 0.19800 0.69000 1.00000
N1 N 8 g 0.62500 0.64500 0.14000 1.00000
P1 P 8 g 0.57400 0.79100 0.17000 1.00000
```

$PNCl_2$ ($E1_4$): A2BC_tP32_86_2g_g_g - POSCAR

```
A2BC_tP32_86_2g_g_g & a,c/a,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,z4 --params
    ↪ =10.82,0.549907578558,0.551,0.852,0.86,0.594,0.198,0.69,0.625,
    ↪ 0.645,0.14,0.574,0.791,0.17 & P4_{2}/n C_{4h}^{4} #86 (g^4) &
    ↪ tP32 & $E1_{4}$ & Cl2NP & Cl2NP & J. A. A. Ketelaar and T. A. {
    ↪ de Vries}, Rec. Trav. Chim. Pays-Bas 58, 1081-1099 (1939)
   1.00000000000000
  10.82000000000000   0.00000000000000   0.00000000000000
   0.00000000000000  10.82000000000000   0.00000000000000
   0.00000000000000   0.00000000000000   5.95000000000000
    Cl    N    P
    16    8    8
Direct
   0.55100000000000   0.85200000000000   0.86000000000000  Cl   (8g)
  -0.05100000000000  -0.35200000000000   0.86000000000000  Cl   (8g)
  -0.85200000000000   1.05100000000000   1.36000000000000  Cl   (8g)
   1.35200000000000  -0.55100000000000   1.36000000000000  Cl   (8g)
  -0.55100000000000  -0.85200000000000  -0.86000000000000  Cl   (8g)
   1.05100000000000   1.35200000000000  -0.86000000000000  Cl   (8g)
   0.85200000000000  -0.05100000000000  -0.36000000000000  Cl   (8g)
  -0.35200000000000   0.55100000000000  -0.36000000000000  Cl   (8g)
   0.59400000000000   0.19800000000000   0.69000000000000  Cl   (8g)
  -0.09400000000000   0.30200000000000   0.69000000000000  Cl   (8g)
  -0.19800000000000   1.09400000000000   1.19000000000000  Cl   (8g)
   0.69800000000000  -0.59400000000000   1.19000000000000  Cl   (8g)
  -0.59400000000000  -0.19800000000000  -0.69000000000000  Cl   (8g)
   1.09400000000000   0.69800000000000  -0.69000000000000  Cl   (8g)
   0.19800000000000  -0.09400000000000  -0.19000000000000  Cl   (8g)
   0.30200000000000   0.59400000000000  -0.19000000000000  Cl   (8g)
   0.62500000000000   0.64500000000000   0.14000000000000  N    (8g)
  -0.12500000000000  -0.14500000000000   0.14000000000000  N    (8g)
  -0.64500000000000   1.12500000000000   0.64000000000000  N    (8g)
   1.14500000000000  -0.62500000000000   0.64000000000000  N    (8g)
  -0.62500000000000  -0.64500000000000  -0.14000000000000  N    (8g)
   1.12500000000000   1.14500000000000  -0.14000000000000  N    (8g)
   0.64500000000000  -0.12500000000000   0.36000000000000  N    (8g)
  -0.14500000000000   0.62500000000000   0.36000000000000  N    (8g)
   0.57400000000000   0.79100000000000   0.17000000000000  P    (8g)
  -0.07400000000000  -0.29100000000000   0.17000000000000  P    (8g)
  -0.79100000000000   1.07400000000000   0.67000000000000  P    (8g)
   1.29100000000000  -0.57400000000000   0.67000000000000  P    (8g)
  -0.57400000000000  -0.79100000000000  -0.17000000000000  P    (8g)
   1.07400000000000   1.29100000000000  -0.17000000000000  P    (8g)
   0.79100000000000  -0.07400000000000   0.33000000000000  P    (8g)
  -0.29100000000000   0.57400000000000   0.33000000000000  P    (8g)
```

$Nd_4Re_2O_{11}$: A4B11C2_tP68_86_2g_ab5g_g - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Nd4O11Re2'
_chemical_formula_sum 'Nd4 O11 Re2'

loop_
_publ_author_name
 'K.-A. Wilhelmi'
 'E. Lagervall'
 'O. Muller'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 24
_journal_year 1970
_journal_page_first 3406
_journal_page_last 3408
_publ_Section_title
;
 On the Crystal Structure of Nd$_{4}$Re$_{2}$O$_{11}$
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'Nd$_{4}$Re$_{2}$O$_{11}$ Structure'
_aflow_proto 'A4B11C2_tP68_86_2g_ab5g_g'
_aflow_params 'a,c/a,x_{3},y_{3},z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{
    ↪ 5},x_{6},y_{6},z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},
    ↪ y_{9},z_{9},x_{10},y_{10},z_{10}'
_aflow_params_values '12.676,0.441858630483,0.184,0.1182,0.4979,0.1107,
    ↪ 0.8064,0.603,0.0301,0.1938,0.7009,0.0013,0.4086,0.6937,0.0415,
    ↪ 0.6291,0.7905,0.1484,0.8032,-0.0057,0.4559,0.34,0.1868,0.0247,-
    ↪ 0.087,0.0707'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tP68'

_symmetry_space_group_name_H-M "P 42/n (origin choice 2)"
_symmetry_Int_Tables_number 86

_cell_length_a    12.67600
```

```
_cell_length_b      12.67600
_cell_length_c      5.60100
_cell_angle_alpha   90.00000
_cell_angle_beta    90.00000
_cell_angle_gamma   90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x+1/2,-y+1/2,z
3 -y,x+1/2,z+1/2
4 y+1/2,-x,z+1/2
5 -x,-y,-z
6 x+1/2,y+1/2,-z
7 y,-x+1/2,-z+1/2
8 -y+1/2,x,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 2 a 0.25000 0.25000 0.25000 1.00000
O2 O 2 b 0.25000 0.25000 0.75000 1.00000
Nd1 Nd 8 g 0.18400 0.11820 0.49790 1.00000
Nd2 Nd 8 g 0.11070 0.80640 0.60300 1.00000
O3 O 8 g 0.03010 0.19380 0.70090 1.00000
O4 O 8 g 0.00130 0.40860 0.69370 1.00000
O5 O 8 g 0.04150 0.62910 0.79050 1.00000
O6 O 8 g 0.14840 0.80320 -0.00570 1.00000
O7 O 8 g 0.45590 0.34000 0.18680 1.00000
Re1 Re 8 g 0.02470 -0.08700 0.07070 1.00000
```

$Nd_4Re_2O_{11}$: A4B11C2_tP68_86_2g_ab5g_g - POSCAR

```
A4B11C2_tP68_86_2g_ab5g_g & a,c/a,x3,y3,z3,x4,y4,z4,x5,y5,z5,x6,y6,z6,x7
    ↪ ,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10 --params=12.676,
    ↪ 0.441858630483,0.184,0.1182,0.4979,0.1107,0.8064,0.603,0.0301,
    ↪ 0.1938,0.7009,0.0013,0.4086,0.6937,0.0415,0.6291,0.7905,0.1484,
    ↪ 0.8032,-0.0057,0.4559,0.34,0.1868,0.0247,-0.087,0.0707 & P4_{2
    ↪ }/n C_{4h}^{4} #86 (abg^8) & tP68 & None & Nd4O11Re2 &
    ↪ Nd4O11Re2 & K.-A. Wilhelmi and E. Lagervall and O. Muller, Acta
    ↪  Chem. Scand. 24, 3406-3408 (1970)
   1.00000000000000
  12.67600000000000    0.00000000000000    0.00000000000000
   0.00000000000000   12.67600000000000    0.00000000000000
   0.00000000000000    0.00000000000000    5.60100000000000
    Nd     O     Re
    16    44     8
Direct
   0.18400000000000    0.11820000000000    0.49790000000000  Nd  (8g)
   0.31600000000000    0.38180000000000    0.49790000000000  Nd  (8g)
  -0.11820000000000    0.68400000000000    0.99790000000000  Nd  (8g)
   0.61820000000000   -0.18400000000000    0.99790000000000  Nd  (8g)
  -0.18400000000000   -0.11820000000000   -0.49790000000000  Nd  (8g)
   0.68400000000000    0.61820000000000   -0.49790000000000  Nd  (8g)
   0.11820000000000    0.31600000000000    0.00210000000000  Nd  (8g)
   0.38180000000000    0.18400000000000    0.00210000000000  Nd  (8g)
   0.11070000000000    0.80640000000000    0.60300000000000  Nd  (8g)
   0.38930000000000   -0.30640000000000    0.60300000000000  Nd  (8g)
  -0.80640000000000    0.61070000000000    1.10300000000000  Nd  (8g)
   1.30640000000000   -0.11070000000000    1.10300000000000  Nd  (8g)
  -0.11070000000000   -0.80640000000000   -0.60300000000000  Nd  (8g)
   0.61070000000000    1.30640000000000   -0.60300000000000  Nd  (8g)
   0.80640000000000    0.38930000000000   -0.10300000000000  Nd  (8g)
  -0.30640000000000    0.11070000000000   -0.10300000000000  Nd  (8g)
   0.25000000000000    0.25000000000000    0.25000000000000  O   (2a)
   0.75000000000000    0.75000000000000    0.75000000000000  O   (2a)
   0.25000000000000    0.25000000000000    0.75000000000000  O   (2b)
   0.75000000000000    0.75000000000000    0.25000000000000  O   (2b)
   0.03010000000000    0.19380000000000    0.70090000000000  O   (8g)
   0.46990000000000    0.30620000000000    0.70090000000000  O   (8g)
  -0.19380000000000    0.53010000000000    1.20090000000000  O   (8g)
   0.69380000000000   -0.03010000000000    1.20090000000000  O   (8g)
  -0.03010000000000   -0.19380000000000   -0.70090000000000  O   (8g)
   0.53010000000000    0.69380000000000   -0.70090000000000  O   (8g)
   0.19380000000000    0.46990000000000   -0.20090000000000  O   (8g)
   0.30620000000000    0.03010000000000   -0.20090000000000  O   (8g)
   0.00130000000000    0.40860000000000    0.69370000000000  O   (8g)
   0.49870000000000    0.09140000000000    0.69370000000000  O   (8g)
  -0.40860000000000    0.50130000000000    1.19370000000000  O   (8g)
   0.90860000000000   -0.00130000000000    1.19370000000000  O   (8g)
  -0.00130000000000   -0.40860000000000   -0.69370000000000  O   (8g)
   0.50130000000000    0.90860000000000   -0.69370000000000  O   (8g)
   0.40860000000000    0.49870000000000   -0.19370000000000  O   (8g)
   0.09140000000000    0.00130000000000   -0.19370000000000  O   (8g)
   0.04150000000000    0.62910000000000    0.79050000000000  O   (8g)
   0.45850000000000   -0.12910000000000    0.79050000000000  O   (8g)
  -0.62910000000000    0.54150000000000    1.29050000000000  O   (8g)
   1.12910000000000   -0.04150000000000    1.29050000000000  O   (8g)
  -0.04150000000000   -0.62910000000000   -0.79050000000000  O   (8g)
   0.54150000000000    1.12910000000000   -0.79050000000000  O   (8g)
   0.62910000000000    0.45850000000000   -0.29050000000000  O   (8g)
  -0.12910000000000    0.04150000000000   -0.29050000000000  O   (8g)
   0.14840000000000    0.80320000000000   -0.00570000000000  O   (8g)
   0.35160000000000   -0.30320000000000   -0.00570000000000  O   (8g)
  -0.80320000000000    0.64840000000000    0.49430000000000  O   (8g)
   1.30320000000000   -0.14840000000000    0.49430000000000  O   (8g)
  -0.14840000000000   -0.80320000000000    0.00570000000000  O   (8g)
   0.64840000000000    1.30320000000000    0.00570000000000  O   (8g)
   0.80320000000000    0.35160000000000    0.50570000000000  O   (8g)
  -0.30320000000000    0.14840000000000    0.50570000000000  O   (8g)
   0.45590000000000    0.34000000000000    0.18680000000000  O   (8g)
   0.04410000000000    0.16000000000000    0.18680000000000  O   (8g)
  -0.34000000000000    0.95590000000000    0.68680000000000  O   (8g)
   0.84000000000000   -0.45590000000000    0.68680000000000  O   (8g)
  -0.45590000000000   -0.34000000000000   -0.18680000000000  O   (8g)
   0.95590000000000    0.84000000000000   -0.18680000000000  O   (8g)
   0.34000000000000    0.04410000000000    0.31320000000000  O   (8g)
   0.16000000000000    0.45590000000000    0.31320000000000  O   (8g)
   0.02470000000000   -0.08700000000000    0.07070000000000  Re  (8g)
   0.47530000000000    0.58700000000000    0.07070000000000  Re  (8g)
   0.08700000000000    0.52470000000000    0.57070000000000  Re  (8g)
   0.41300000000000   -0.02470000000000    0.57070000000000  Re  (8g)
  -0.02470000000000    0.08700000000000   -0.07070000000000  Re  (8g)
   0.52470000000000    0.41300000000000   -0.07070000000000  Re  (8g)
  -0.08700000000000    0.47530000000000    0.42930000000000  Re  (8g)
   0.58700000000000    0.02470000000000    0.42930000000000  Re  (8g)
```

$NaSb(OH)_6$ ($J1_{11}$): AB6C_tP32_86_d_3g_c - CIF

```
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_audit_creation_method FINDSYM

_chemical_name_mineral 'Na(OH)6Sb'
_chemical_formula_sum 'Na O6 Sb'

loop_
_publ_author_name
 'T. Asai'
_journal_name_full_name
;
 Bulletin of the Chemical Society of Japan
;
_journal_volume 48
_journal_year 1975
_journal_page_first 2677
_journal_page_last 2679
_publ_Section_title
;
 Refinement of the Crystal Structure of Sodium Hexahydroxoantimonate(V),
     ↪ NaSb(OH)$_{6}$
;

# Found in The Fascination of Crystals and Symmetry, {NaSb(OH)$_{6}$},
    ↪ 2014

_aflow_title 'NaSb(OH)$_{6}$ ($J1_{11}$) Structure'
_aflow_proto 'AB6C_tP32_86_d_3g_c'
_aflow_params 'a,c/a,x_{3},y_{3},z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{
    ↪ 5}'
_aflow_params_values '8.029,0.983185950928,0.0558,0.2226,-0.0982,0.2792,
    ↪ 0.582,-0.0828,0.0895,0.0685,0.2219'
_aflow_Strukturbericht '$J1_{11}$'
_aflow_Pearson 'tP32'

_symmetry_space_group_name_H-M "P 42/n (origin choice 2)"
_symmetry_Int_Tables_number 86

_cell_length_a      8.02900
_cell_length_b      8.02900
_cell_length_c      7.89400
_cell_angle_alpha   90.00000
_cell_angle_beta    90.00000
_cell_angle_gamma   90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x+1/2,-y+1/2,z
3 -y,x+1/2,z+1/2
4 y+1/2,-x,z+1/2
5 -x,-y,-z
6 x+1/2,y+1/2,-z
7 y,-x+1/2,-z+1/2
8 -y+1/2,x,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Sb1 Sb 4 c 0.00000 0.00000 0.00000 1.00000
Na1 Na 4 d 0.00000 0.00000 0.50000 1.00000
O1 O 8 g 0.05580 0.22260 -0.09820 1.00000
O2 O 8 g 0.27920 0.58200 -0.08280 1.00000
O3 O 8 g 0.08950 0.06850 0.22190 1.00000
```

$NaSb(OH)_6$ ($J1_{11}$): AB6C_tP32_86_d_3g_c - POSCAR

```
AB6C_tP32_86_d_3g_c & a,c/a,x3,y3,z3,x4,y4,z4,x5,y5,z5 --params=8.029,
    ↪ 0.983185950928,0.0558,0.2226,-0.0982,0.2792,0.582,-0.0828,
    ↪ 0.0895,0.0685,0.2219 & P4_{2}/n C_{4h}^{4} #86 (cdg^3) & tP32 &
    ↪  $J1_{11}$ & Na(OH)6Sb & Na(OH)6Sb & T. Asai, Bull. Chem. Soc.
    ↪ Jpn. 48, 2677-2679 (1975)
   1.00000000000000
   8.02900000000000    0.00000000000000    0.00000000000000
   0.00000000000000    8.02900000000000    0.00000000000000
   0.00000000000000    0.00000000000000    7.89400000000000
```

```
    Na     O     Sb
     4    24     4
Direct
     0.00000000000000     0.00000000000000     0.50000000000000   Na   (4d)
     0.50000000000000     0.50000000000000     0.50000000000000   Na   (4d)
     0.00000000000000     0.50000000000000     0.00000000000000   Na   (4d)
     0.50000000000000     0.00000000000000     0.00000000000000   Na   (4d)
     0.05580000000000     0.22260000000000    -0.09820000000000   O    (8g)
     0.44420000000000     0.27740000000000    -0.09820000000000   O    (8g)
    -0.22260000000000     0.55580000000000     0.40180000000000   O    (8g)
     0.72260000000000    -0.05580000000000     0.40180000000000   O    (8g)
    -0.05580000000000    -0.22260000000000     0.09820000000000   O    (8g)
     0.55580000000000     0.72260000000000     0.09820000000000   O    (8g)
     0.22260000000000     0.44420000000000     0.59820000000000   O    (8g)
     0.27740000000000     0.05580000000000     0.59820000000000   O    (8g)
     0.27920000000000     0.58200000000000    -0.08280000000000   O    (8g)
     0.22080000000000    -0.08200000000000    -0.08280000000000   O    (8g)
    -0.58200000000000     0.77920000000000     0.41720000000000   O    (8g)
     1.08200000000000    -0.27920000000000     0.41720000000000   O    (8g)
    -0.27920000000000    -0.58200000000000     0.08280000000000   O    (8g)
     0.77920000000000     1.08200000000000     0.08280000000000   O    (8g)
     0.58200000000000     0.22080000000000     0.58280000000000   O    (8g)
    -0.08200000000000     0.27920000000000     0.58280000000000   O    (8g)
     0.08950000000000     0.06850000000000     0.22190000000000   O    (8g)
     0.41050000000000     0.43150000000000     0.22190000000000   O    (8g)
    -0.06850000000000     0.58950000000000     0.72190000000000   O    (8g)
     0.56850000000000    -0.08950000000000     0.72190000000000   O    (8g)
    -0.08950000000000    -0.06850000000000    -0.22190000000000   O    (8g)
     0.58950000000000     0.56850000000000    -0.22190000000000   O    (8g)
     0.06850000000000     0.41050000000000     0.27810000000000   O    (8g)
     0.43150000000000     0.08950000000000     0.27810000000000   O    (8g)
     0.00000000000000     0.00000000000000     0.00000000000000   Sb   (4c)
     0.50000000000000     0.50000000000000     0.00000000000000   Sb   (4c)
     0.00000000000000     0.50000000000000     0.50000000000000   Sb   (4c)
     0.50000000000000     0.00000000000000     0.50000000000000   Sb   (4c)
```

β-$LiIO_3$: ABC3_tP40_86_g_g_3g - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'ILiO3'
_chemical_formula_sum 'I Li O3'

loop_
_publ_author_name
 'H. Schulz'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 29
_journal_year 1973
_journal_page_first 2285
_journal_page_last 2289
_publ_Section_title
;
 The structure of $\beta$-LiIO$_{3}$
;

_aflow_title '$\beta$-LiIO$_{3}$ Structure'
_aflow_proto 'ABC3_tP40_86_g_g_3g'
_aflow_params 'a,c/a,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3},y_{3},z_{
      ↪ 3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{5}'
_aflow_params_values '9.7329,0.63255555898,0.0343,0.7576,0.115,0.426,
      ↪ 0.236,0.426,0.836,0.121,-0.05,0.094,0.209,0.094,0.847,0.552,
      ↪ 0.172'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tP40'

_symmetry_space_group_name_H-M "P 42/n (origin choice 2)"
_symmetry_Int_Tables_number 86

_cell_length_a     9.73290
_cell_length_b     9.73290
_cell_length_c     6.15660
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x+1/2,-y+1/2,z
3 -y,x+1/2,z+1/2
4 y+1/2,-x,z+1/2
5 -x,-y,-z
6 x+1/2,y+1/2,-z
7 y,-x+1/2,-z+1/2
8 -y+1/2,x,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
I1 I 8 g 0.03430 0.75760 0.11500 1.00000
Li1 Li 8 g 0.42600 0.23600 0.42600 1.00000
O1 O 8 g 0.83600 0.12100 -0.05000 1.00000
O2 O 8 g 0.09400 0.20900 0.09400 1.00000
O3 O 8 g 0.84700 0.55200 0.17200 1.00000
```

β-$LiIO_3$: ABC3_tP40_86_g_g_3g - POSCAR

```
ABC3_tP40_86_g_g_3g & a,c/a,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,z4,x5,y5,z5
      ↪ --params=9.7329,0.63255555898,0.0343,0.7576,0.115,0.426,0.236,
      ↪ 0.426,0.836,0.121,-0.05,0.094,0.209,0.094,0.847,0.552,0.172 &
      ↪ P4_{2}/n C_{4h}^{4} #86 (g^5) & tP40 & None & ILiO3 & ILiO3 &
      ↪ H. Schulz, Acta Crystallogr. Sect. B Struct. Sci. 29, 2285-2289
      ↪ (1973)
   1.00000000000000
   9.73290000000000     0.00000000000000     0.00000000000000
   0.00000000000000     9.73290000000000     0.00000000000000
   0.00000000000000     0.00000000000000     6.15660000000000
     I    Li    O
     8     8    24
Direct
     0.03430000000000     0.75760000000000     0.11500000000000   I    (8g)
     0.46570000000000    -0.25760000000000     0.11500000000000   I    (8g)
    -0.75760000000000     0.53430000000000     0.61500000000000   I    (8g)
     1.25760000000000    -0.03430000000000     0.61500000000000   I    (8g)
    -0.03430000000000    -0.75760000000000    -0.11500000000000   I    (8g)
     0.53430000000000     1.25760000000000    -0.11500000000000   I    (8g)
     0.75760000000000     0.46570000000000     0.38500000000000   I    (8g)
    -0.25760000000000     0.03430000000000     0.38500000000000   I    (8g)
     0.42600000000000     0.23600000000000     0.42600000000000   Li   (8g)
     0.07400000000000     0.26400000000000     0.42600000000000   Li   (8g)
    -0.23600000000000     0.92600000000000     0.92600000000000   Li   (8g)
     0.73600000000000    -0.42600000000000     0.92600000000000   Li   (8g)
    -0.42600000000000    -0.23600000000000    -0.42600000000000   Li   (8g)
     0.92600000000000     0.73600000000000    -0.42600000000000   Li   (8g)
     0.23600000000000     0.07400000000000     0.07400000000000   Li   (8g)
     0.26400000000000     0.42600000000000     0.07400000000000   Li   (8g)
     0.83600000000000     0.12100000000000    -0.05000000000000   O    (8g)
    -0.33600000000000     0.37900000000000    -0.05000000000000   O    (8g)
    -0.12100000000000     1.33600000000000     0.45000000000000   O    (8g)
     0.62100000000000    -0.83600000000000     0.45000000000000   O    (8g)
    -0.83600000000000    -0.12100000000000     0.05000000000000   O    (8g)
     1.33600000000000     0.62100000000000     0.05000000000000   O    (8g)
     0.12100000000000    -0.33600000000000     0.55000000000000   O    (8g)
     0.37900000000000     0.83600000000000     0.55000000000000   O    (8g)
     0.09400000000000     0.20900000000000     0.09400000000000   O    (8g)
     0.40600000000000     0.29100000000000     0.09400000000000   O    (8g)
    -0.20900000000000     0.59400000000000     0.59400000000000   O    (8g)
     0.70900000000000    -0.09400000000000     0.59400000000000   O    (8g)
    -0.09400000000000    -0.20900000000000    -0.09400000000000   O    (8g)
     0.59400000000000     0.70900000000000    -0.09400000000000   O    (8g)
     0.20900000000000     0.40600000000000     0.40600000000000   O    (8g)
     0.29100000000000     0.09400000000000     0.40600000000000   O    (8g)
     0.84700000000000     0.55200000000000     0.17200000000000   O    (8g)
    -0.34700000000000    -0.05200000000000     0.17200000000000   O    (8g)
    -0.55200000000000     1.34700000000000     0.67200000000000   O    (8g)
     1.05200000000000    -0.84700000000000     0.67200000000000   O    (8g)
    -0.84700000000000    -0.55200000000000    -0.17200000000000   O    (8g)
     1.34700000000000     1.05200000000000    -0.17200000000000   O    (8g)
     0.55200000000000    -0.34700000000000     0.32800000000000   O    (8g)
    -0.05200000000000     0.84700000000000     0.32800000000000   O    (8g)
```

Marialite Scapolite [$Na_4Cl(AlSi_3)_3O_{24}$, $S6_4$]: AB4C24D12_tI82_87_a_h_2h2i_hi - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Marialite scapolite'
_chemical_formula_sum 'Cl Na4 O24 Si12'

loop_
_publ_author_name
 'J. J. Papike'
 'T. Zoltai'
_journal_name_full_name
;
 American Mineralogist
;
_journal_volume 50
_journal_year 1965
_journal_page_first 641
_journal_page_last 655
_publ_Section_title
;
 The crystal structure of a marialite scapolite
;

_aflow_title 'Marialite Scapolite [Na$_{4}$Cl(AlSi$_{3}$)$_{3}$O$_{24}$,
      ↪ $S6_{4}$] Structure'
_aflow_proto 'AB4C24D12_tI82_87_a_h_2h2i_hi'
_aflow_params 'a,c/a,x_{2},y_{2},x_{3},y_{3},x_{4},y_{4},x_{5},y_{5},x_{
      ↪ 6},y_{6},z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},z_{8}'
_aflow_params_values '12.06,0.627860696517,0.134,0.2113,0.4587,0.3483,
      ↪ 0.3066,0.1206,0.3388,0.4104,0.0517,0.35,0.2148,0.2293,0.1289,
      ↪ 0.3281,0.3374,0.0851,0.206'
_aflow_Strukturbericht '$S6_{4}$'
_aflow_Pearson 'tI82'

_symmetry_space_group_name_H-M "I 4/m"
_symmetry_Int_Tables_number 87

_cell_length_a     12.06000
_cell_length_b     12.06000
_cell_length_c     7.57200
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000
```

```
loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,z
3 -y,x,z
4 y,-x,z
5 -x,-y,-z
6 x,y,-z
7 y,-x,-z
8 -y,x,-z
9 x+1/2,y+1/2,z+1/2
10 -x+1/2,-y+1/2,z+1/2
11 -y+1/2,x+1/2,z+1/2
12 y+1/2,-x+1/2,z+1/2
13 -x+1/2,-y+1/2,-z+1/2
14 x+1/2,y+1/2,-z+1/2
15 y+1/2,-x+1/2,-z+1/2
16 -y+1/2,x+1/2,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cl1 Cl 2 a 0.00000 0.00000 0.00000 1.00000
Na1 Na 8 h 0.13400 0.21130 0.00000 1.00000
O1 O 8 h 0.45870 0.34830 0.00000 1.00000
O2 O 8 h 0.30660 0.12060 0.00000 1.00000
Si1 Si 8 h 0.33880 0.41040 0.00000 1.00000
O3 O 16 i 0.05170 0.35000 0.21480 1.00000
O4 O 16 i 0.22930 0.12890 0.32810 1.00000
Si2 Si 16 i 0.33740 0.08510 0.20600 1.00000
```

Marialite Scapolite [$Na_4Cl(AlSi_3)_3O_{24}$, $S6_4$]: AB4C24D12_tI82_87_a_h_2h2i_hi - POSCAR

```
AB4C24D12_tI82_87_a_h_2h2i_hi & a,c/a,x2,y2,x3,y3,x4,y4,x5,y5,x6,y6,z6,
    ↪ x7,y7,z7,x8,y8,z8 --params=12.06,0.627860696517,0.134,0.2113,
    ↪ 0.4587,0.3483,0.3066,0.1206,0.3388,0.4104,0.0517,0.35,0.2148,
    ↪ 0.2293,0.1289,0.3281,0.3374,0.0851,0.206 & I4/m C_{4h}^{5} #87
    ↪ (ah^4i^3) & tI82 & $S6_{4}$ & ClNa4O24(Al3Si9) & Marialite
    ↪ scapolite & J. J. Papike and T. Zoltai, Am. Mineral. 50,
    ↪ 641-655 (1965)
   1.00000000000000
  -6.03000000000000    6.03000000000000    3.78600000000000
   6.03000000000000   -6.03000000000000    3.78600000000000
   6.03000000000000    6.03000000000000   -3.78600000000000
     Cl     Na      O     Si
      1      4     24     12
Direct
   0.00000000000000    0.00000000000000    0.00000000000000    Cl    (2a)
   0.21130000000000    0.13400000000000    0.34530000000000    Na    (8h)
  -0.21130000000000   -0.13400000000000   -0.34530000000000    Na    (8h)
   0.13400000000000   -0.21130000000000   -0.07730000000000    Na    (8h)
  -0.13400000000000    0.21130000000000    0.07730000000000    Na    (8h)
   0.34830000000000    0.45870000000000    0.80700000000000    O     (8h)
  -0.34830000000000   -0.45870000000000   -0.80700000000000    O     (8h)
   0.45870000000000   -0.34830000000000    0.11040000000000    O     (8h)
  -0.45870000000000    0.34830000000000   -0.11040000000000    O     (8h)
   0.12060000000000    0.30660000000000    0.42720000000000    O     (8h)
  -0.12060000000000   -0.30660000000000   -0.42720000000000    O     (8h)
   0.30660000000000   -0.12060000000000    0.18600000000000    O     (8h)
  -0.30660000000000    0.12060000000000   -0.18600000000000    O     (8h)
   0.56480000000000    0.26650000000000    0.40170000000000    O    (16i)
  -0.13520000000000    0.16310000000000   -0.40170000000000    O    (16i)
   0.26650000000000   -0.13520000000000   -0.29830000000000    O    (16i)
   0.16310000000000    0.56480000000000    0.29830000000000    O    (16i)
  -0.56480000000000   -0.26650000000000   -0.40170000000000    O    (16i)
   0.13520000000000   -0.16310000000000    0.40170000000000    O    (16i)
  -0.26650000000000    0.13520000000000    0.29830000000000    O    (16i)
  -0.16310000000000   -0.56480000000000   -0.29830000000000    O    (16i)
   0.45700000000000    0.55740000000000    0.35820000000000    O    (16i)
   0.19920000000000    0.09880000000000   -0.35820000000000    O    (16i)
   0.55740000000000    0.19920000000000    0.10040000000000    O    (16i)
   0.09880000000000    0.45700000000000   -0.10040000000000    O    (16i)
  -0.45700000000000   -0.55740000000000   -0.35820000000000    O    (16i)
  -0.19920000000000   -0.09880000000000    0.35820000000000    O    (16i)
  -0.55740000000000   -0.19920000000000   -0.10040000000000    O    (16i)
  -0.09880000000000   -0.45700000000000    0.10040000000000    O    (16i)
   0.41040000000000    0.33880000000000    0.74920000000000    Si    (8h)
  -0.41040000000000   -0.33880000000000   -0.74920000000000    Si    (8h)
   0.33880000000000   -0.41040000000000   -0.07160000000000    Si    (8h)
  -0.33880000000000    0.41040000000000    0.07160000000000    Si    (8h)
   0.29110000000000    0.54340000000000    0.42250000000000    Si    (16i)
   0.12090000000000   -0.13140000000000   -0.42250000000000    Si    (16i)
   0.54340000000000    0.12090000000000    0.25230000000000    Si    (16i)
  -0.13140000000000    0.29110000000000   -0.25230000000000    Si    (16i)
  -0.29110000000000   -0.54340000000000   -0.42250000000000    Si    (16i)
  -0.12090000000000    0.13140000000000    0.42250000000000    Si    (16i)
  -0.54340000000000   -0.12090000000000   -0.25230000000000    Si    (16i)
   0.13140000000000   -0.29110000000000    0.25230000000000    Si    (16i)
```

Sr_2NiWO_6: AB6C2D_tI20_87_a_eh_d_b - CIF

```
# CIF file
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_audit_creation_method FINDSYM

_chemical_name_mineral 'NiO6Sr2W'
_chemical_formula_sum 'Ni O6 Sr2 W'

loop_
_publ_author_name
 'D. Iwanaga'
 'Y. Inaguma'
 'M. Itoh'
_journal_name_full_name
;
 Materials Research Bulletin
;
_journal_volume 35
_journal_year 2000
_journal_page_first 449
_journal_page_last 457
_publ_Section_title
;
 Structure and Magnetic Properties of Sr$_{2}$Ni$A$O$_{6}$ ($A$ = W, Te)
;

_aflow_title 'Sr$_{2}$NiWO$_{6}$ Structure'
_aflow_proto 'AB6C2D_tI20_87_a_eh_d_b'
_aflow_params 'a,c/a,z_{4},x_{5},y_{5}'
_aflow_params_values '5.5571,1.42396213853,0.255,0.289,0.227'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tI20'

_symmetry_space_group_name_H-M "I 4/m"
_symmetry_Int_Tables_number 87

_cell_length_a      5.55710
_cell_length_b      5.55710
_cell_length_c      7.91310
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,z
3 -y,x,z
4 y,-x,z
5 -x,-y,-z
6 x,y,-z
7 y,-x,-z
8 -y,x,-z
9 x+1/2,y+1/2,z+1/2
10 -x+1/2,-y+1/2,z+1/2
11 -y+1/2,x+1/2,z+1/2
12 y+1/2,-x+1/2,z+1/2
13 -x+1/2,-y+1/2,-z+1/2
14 x+1/2,y+1/2,-z+1/2
15 y+1/2,-x+1/2,-z+1/2
16 -y+1/2,x+1/2,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ni1 Ni 2 a 0.00000 0.00000 0.00000 1.00000
W1 W 2 b 0.00000 0.00000 0.50000 1.00000
Sr1 Sr 4 d 0.00000 0.50000 0.25000 1.00000
O1 O 4 e 0.00000 0.00000 0.25500 1.00000
O2 O 8 h 0.28900 0.22700 0.00000 1.00000
```

Sr_2NiWO_6: AB6C2D_tI20_87_a_eh_d_b - POSCAR

```
AB6C2D_tI20_87_a_eh_d_b & a,c/a,z4,x5,y5 --params=5.5571,1.42396213853,
    ↪ 0.255,0.289,0.227 & I4/m C_{4h}^{5} #87 (abdeh) & tI20 & None &
    ↪ NiO6Sr2W & NiO6Sr2W & D. Iwanaga and Y. Inaguma and M. Itoh,
    ↪ Mater. Res. Bull. 35, 449-457 (2000)
   1.00000000000000
  -2.77855000000000    2.77855000000000    3.95655000000000
   2.77855000000000   -2.77855000000000    3.95655000000000
   2.77855000000000    2.77855000000000   -3.95655000000000
     Ni      O     Sr      W
      1      6      2      1
Direct
   0.00000000000000    0.00000000000000    0.00000000000000    Ni    (2a)
   0.25500000000000    0.25500000000000    0.00000000000000    O     (4e)
  -0.25500000000000   -0.25500000000000    0.00000000000000    O     (4e)
   0.22700000000000    0.28900000000000    0.51600000000000    O     (8h)
  -0.22700000000000   -0.28900000000000   -0.51600000000000    O     (8h)
   0.28900000000000   -0.22700000000000    0.06200000000000    O     (8h)
  -0.28900000000000    0.22700000000000   -0.06200000000000    O     (8h)
   0.75000000000000    0.25000000000000    0.50000000000000    Sr    (4d)
   0.25000000000000    0.75000000000000    0.50000000000000    Sr    (4d)
   0.50000000000000    0.50000000000000    0.00000000000000    W     (2b)
```

$Na_4Ge_9O_{20}$: A9B4C20_tI132_88_a2f_f_5f - CIF

```
# CIF file
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_audit_creation_method FINDSYM

_chemical_name_mineral 'Ge9Na4O20'
_chemical_formula_sum 'Ge9 Na4 O20'

loop_
```

```
_publ_author_name
 'N. Ingri'
 'G. Lundgren'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 17
_journal_year 1963
_journal_page_first 617
_journal_page_last 633
_publ_Section_title
;
 The Crystal Structure of Na$_{4}$Ge$_{9}$O$_{20}$
;

_aflow_title 'Na$_{4}$Ge$_{9}$O$_{20}$ Structure'
_aflow_proto 'A9B4C20_tI132_88_a2f_f_5f'
_aflow_params 'a,c/a,x_{2},y_{2},z_{2},x_{3},y_{3},z_{3},x_{4},y_{4},z_{
      ↪ 4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{6},x_{7},y_{7},z_{7},x_{8},
      ↪ y_{8},z_{8},x_{9},y_{9},z_{9}'
_aflow_params_values '14.97999,0.492924227586,0.1366,0.046,0.7009,0.0956
      ↪ ,0.218,0.4912,0.0878,0.0544,0.1702,0.0928,0.2127,0.2448,0.0824,
      ↪ 0.2227,0.7476,0.1889,0.0739,-0.0926,0.0253,0.0446,0.7555,0.1554
      ↪ ,0.1088,0.507'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tI132'

_symmetry_space_group_name_H-M "I 41/a (origin choice 2)"
_symmetry_Int_Tables_number 88

_cell_length_a     14.97999
_cell_length_b     14.97999
_cell_length_c     7.38400
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y+1/2,z
3 -y+3/4,x+1/4,z+1/4
4 y+1/4,-x+1/4,z+1/4
5 -x,-y,-z
6 x,y+1/2,-z
7 y+1/4,-x+3/4,-z+3/4
8 -y+3/4,x+3/4,-z+3/4
9 x+1/2,y+1/2,z+1/2
10 -x+1/2,-y,z+1/2
11 -y+1/4,x+3/4,z+3/4
12 y+3/4,-x+3/4,z+3/4
13 -x+1/2,-y+1/2,-z+1/2
14 x+1/2,y,-z+1/2
15 y+3/4,-x+1/4,-z+1/4
16 -y+1/4,x+1/4,-z+1/4

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ge1 Ge 4 a 0.00000 0.25000 0.12500 1.00000
Ge2 Ge 16 f 0.13660 0.04600 0.70090 1.00000
Ge3 Ge 16 f 0.09560 0.21800 0.49120 1.00000
Na1 Na 16 f 0.08780 0.05440 0.17020 1.00000
O1 O 16 f 0.09280 0.21270 0.24480 1.00000
O2 O 16 f 0.08240 0.22270 0.74760 1.00000
O3 O 16 f 0.18890 0.07390 -0.09260 1.00000
O4 O 16 f 0.02530 0.04460 0.75550 1.00000
O5 O 16 f 0.15540 0.10880 0.50700 1.00000
```

$Na_4Ge_9O_{20}$: A9B4C20_tI132_88_a2f_f_5f - POSCAR

```
A9B4C20_tI132_88_a2f_f_5f & a,c/a,x2,y2,z2,x3,y3,z3,x4,y4,z4,x5,y5,z5,x6
      ↪ ,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9 --params=14.97999,
      ↪ 0.492924227586,0.1366,0.046,0.7009,0.0956,0.218,0.4912,0.0878,
      ↪ 0.0544,0.1702,0.0928,0.2127,0.2448,0.0824,0.2227,0.7476,0.1889,
      ↪ 0.0739,-0.0926,0.0253,0.0446,0.7555,0.1554,0.1088,0.507 & I4_{1
      ↪ }/a C_{4h}^{6} #88 (af^8) & tI132 & None & Ge9Na4O20 &
      ↪ Ge9Na4O20 & N. Ingri and G. Lundgren, Acta Chem. Scand. 17,
      ↪ 617-633 (1963)
   1.00000000000000
  -7.48999500000000    7.48999500000000    3.69200000000000
   7.48999500000000   -7.48999500000000    3.69200000000000
   7.48999500000000    7.48999500000000   -3.69200000000000
     Ge     Na      O
     18      8     40
Direct
   0.37500000000000    0.12500000000000    0.25000000000000   Ge   (4a)
   0.62500000000000    0.87500000000000    0.75000000000000   Ge   (4a)
   0.74690000000000    0.83750000000000    0.18260000000000   Ge   (16f)
   1.15490000000000    0.56430000000000    0.31740000000000   Ge   (16f)
   1.33750000000000    0.65490000000000    0.09060000000000   Ge   (16f)
   1.06430000000000    1.24690000000000    0.40940000000000   Ge   (16f)
  -0.74690000000000   -0.83750000000000   -0.18260000000000   Ge   (16f)
  -0.15490000000000   -0.56430000000000    0.68260000000000   Ge   (16f)
  -0.33750000000000   -0.65490000000000   -0.09060000000000   Ge   (16f)
  -0.06430000000000   -0.24690000000000    0.59060000000000   Ge   (16f)
   0.70920000000000    0.58680000000000    0.31360000000000   Ge   (16f)
   0.77320000000000    0.39560000000000    0.18640000000000   Ge   (16f)
   1.08680000000000    0.27320000000000   -0.12240000000000   Ge   (16f)
   0.89560000000000    1.20920000000000    0.62240000000000   Ge   (16f)
  -0.70920000000000   -0.58680000000000   -0.31360000000000   Ge   (16f)
   0.22680000000000   -0.39560000000000    0.81360000000000   Ge   (16f)
  -0.08680000000000   -0.27320000000000    0.12240000000000   Ge   (16f)
   0.10440000000000   -0.20920000000000    0.37760000000000   Ge   (16f)
   0.22460000000000    0.25800000000000    0.14220000000000   Na   (16f)
   0.61580000000000    0.08240000000000    0.35780000000000   Na   (16f)
   0.75800000000000    0.11580000000000    0.03340000000000   Na   (16f)
   0.58240000000000    0.72460000000000    0.46660000000000   Na   (16f)
  -0.22460000000000   -0.25800000000000   -0.14220000000000   Na   (16f)
   0.38420000000000   -0.08240000000000    0.64220000000000   Na   (16f)
   0.24200000000000   -0.11580000000000   -0.03340000000000   Na   (16f)
   0.41760000000000    0.27540000000000    0.53340000000000   Na   (16f)
   0.45750000000000    0.33760000000000    0.30550000000000   O    (16f)
   0.53210000000000    0.15200000000000    0.19450000000000   O    (16f)
   0.83760000000000    0.03210000000000   -0.11990000000000   O    (16f)
   0.65200000000000    0.95750000000000    0.61990000000000   O    (16f)
  -0.45750000000000   -0.33760000000000   -0.30550000000000   O    (16f)
   0.46790000000000   -0.15200000000000    0.80550000000000   O    (16f)
   0.16240000000000   -0.03210000000000    0.11990000000000   O    (16f)
   0.34800000000000    0.04250000000000    0.38010000000000   O    (16f)
   0.97030000000000    0.83000000000000    0.30510000000000   O    (16f)
   1.02490000000000    0.66520000000000    0.19490000000000   O    (16f)
   1.33000000000000    0.52490000000000   -0.14030000000000   O    (16f)
   1.16520000000000    1.47030000000000    0.64030000000000   O    (16f)
  -0.97030000000000   -0.83000000000000   -0.30510000000000   O    (16f)
  -0.02490000000000   -0.66520000000000    0.80510000000000   O    (16f)
  -0.33000000000000   -0.52490000000000    0.14030000000000   O    (16f)
  -0.16520000000000   -0.47030000000000    0.35970000000000   O    (16f)
  -0.01870000000000    0.09630000000000    0.26280000000000   O    (16f)
   0.33350000000000   -0.28150000000000    0.23720000000000   O    (16f)
   0.59630000000000   -0.16650000000000    0.11500000000000   O    (16f)
   0.21850000000000    0.48130000000000    0.38500000000000   O    (16f)
   0.01870000000000   -0.09630000000000   -0.26280000000000   O    (16f)
   0.66650000000000    0.28150000000000    0.76280000000000   O    (16f)
   0.40370000000000    0.16650000000000   -0.11500000000000   O    (16f)
   0.78150000000000    0.51870000000000    0.61500000000000   O    (16f)
   0.80010000000000    0.78080000000000    0.06990000000000   O    (16f)
   1.21090000000000    0.73020000000000    0.43010000000000   O    (16f)
   1.28080000000000    0.71090000000000   -0.01930000000000   O    (16f)
   1.23020000000000    1.30010000000000    0.51930000000000   O    (16f)
  -0.80010000000000   -0.78080000000000   -0.06990000000000   O    (16f)
  -0.21090000000000   -0.73020000000000    0.56990000000000   O    (16f)
  -0.28080000000000   -0.71090000000000    0.01930000000000   O    (16f)
  -0.23020000000000   -0.30010000000000    0.48070000000000   O    (16f)
   0.61580000000000    0.66240000000000    0.26420000000000   O    (16f)
   0.89820000000000    0.35160000000000    0.23580000000000   O    (16f)
   1.16240000000000    0.39820000000000    0.04660000000000   O    (16f)
   0.85160000000000    1.11580000000000    0.45340000000000   O    (16f)
  -0.61580000000000   -0.66240000000000   -0.26420000000000   O    (16f)
   0.10180000000000   -0.35160000000000    0.76420000000000   O    (16f)
  -0.16240000000000   -0.39820000000000   -0.04660000000000   O    (16f)
   0.14840000000000   -0.11580000000000    0.54660000000000   O    (16f)
```

Copper (I) Azide (CuN_3): AB3_tI32_88_d_cf - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Copper (i) azide'
_chemical_formula_sum 'Cu N3'

loop_
_publ_author_name
 'H. Wilsdorf'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 1
_journal_year 1948
_journal_page_first 115
_journal_page_last 118
_publ_Section_title
;
 Die Kristallstruktur des einwertigen Kupferazids, CuN$_{3}$
;

# Found in {\em Ab initio} study of electronic structure and optical
      ↪ properties of heavy-metal azides: TlN$_{3}$, AgN$_{3}$, and
      ↪ CuN$_{3}$, 2008

_aflow_title 'Copper (I) Azide (CuN$_{3}$) Structure'
_aflow_proto 'AB3_tI32_88_d_cf'
_aflow_params 'a,c/a,x_{3},y_{3},z_{3}'
_aflow_params_values '8.65001,0.646242027466,0.173,0.173,0.375'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tI32'

_symmetry_space_group_name_H-M "I 41/a (origin choice 2)"
_symmetry_Int_Tables_number 88

_cell_length_a     8.65001
_cell_length_b     8.65001
_cell_length_c     5.59000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
```

```
1 x,y,z
2 -x,-y+1/2,z
3 -y+3/4,x+1/4,z+1/4
4 y+1/4,-x+1/4,z+1/4
5 -x,-y,-z
6 x,y+1/2,-z
7 y+1/4,-x+3/4,-z+3/4
8 -y+3/4,x+3/4,-z+3/4
9 x+1/2,y+1/2,z+1/2
10 -x+1/2,-y,z+1/2
11 -y+1/4,x+3/4,z+3/4
12 y+3/4,-x+3/4,z+3/4
13 -x+1/2,-y+1/2,-z+1/2
14 x+1/2,y,-z+1/2
15 y+3/4,-x+1/4,-z+1/4
16 -y+1/4,x+1/4,-z+1/4

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
N1 N 8 c 0.00000 0.00000 0.00000 1.00000
Cu1 Cu 8 d 0.00000 0.00000 0.50000 1.00000
N2 N 16 f 0.17300 0.17300 0.37500 1.00000
```

Copper (I) Azide (CuN_3): AB3_tI32_88_d_cf - POSCAR

```
AB3_tI32_88_d_cf & a,c/a,x3,y3,z3 --params=8.65001,0.646242027466,0.173,
    ↪ 0.173,0.375 & I4_{1}/a C_{4h}^{6} #88 (cdf) & tI32 & None &
    ↪ CuN3 & Copper (i) azide & H. Wilsdorf, Acta Cryst. 1, 115-118 (
    ↪ 1948)
   1.00000000000000
  -4.32500500000000    4.32500500000000    2.79500000000000
   4.32500500000000   -4.32500500000000    2.79500000000000
   4.32500500000000    4.32500500000000   -2.79500000000000
    Cu     N
     4    12
Direct
   0.50000000000000    0.50000000000000    0.00000000000000   Cu    (8d)
   0.00000000000000    0.50000000000000    0.50000000000000   Cu    (8d)
   0.00000000000000    0.50000000000000    0.00000000000000   Cu    (8d)
   0.00000000000000    0.00000000000000    0.50000000000000   Cu    (8d)
   0.00000000000000    0.00000000000000    0.00000000000000   N     (8c)
   0.50000000000000    0.00000000000000    0.50000000000000   N     (8c)
   0.50000000000000    0.00000000000000    0.00000000000000   N     (8c)
   0.50000000000000    0.50000000000000    0.50000000000000   N     (8c)
   0.54800000000000    0.54800000000000    0.34600000000000   N    (16f)
   0.70200000000000    0.20200000000000    0.15400000000000   N    (16f)
   1.04800000000000    0.20200000000000    0.00000000000000   N    (16f)
   0.70200000000000    1.04800000000000    0.50000000000000   N    (16f)
  -0.54800000000000   -0.54800000000000   -0.34600000000000   N    (16f)
   0.29800000000000   -0.20200000000000    0.84600000000000   N    (16f)
  -0.04800000000000   -0.20200000000000    0.00000000000000   N    (16f)
   0.29800000000000   -0.04800000000000    0.50000000000000   N    (16f)
```

Scheelite ($CaWO_4$, $H0_4$): AB4C_tI24_88_b_f_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Scheelite'
_chemical_formula_sum 'Ca O4 W'

loop_
_publ_author_name
 'R. M. Hazen'
 'L. W. Finger'
 'J. W. E. Mariathasan'
_journal_name_full_name
;
 Journal of Physics and Chemistry of Solids
;
_journal_volume 46
_journal_year 1985
_journal_page_first 253
_journal_page_last 263
_publ_Section_title
;
 High-pressure crystal chemistry of scheelite-type tungstates and
    ↪ molybdates
;

# Found in Electronic band structures of the scheelite materials CaMoO$_
    ↪ {4}$, CaWO$_{4}$, PbMoO$_{4}$, and PbWO$_{4}$, 1988

_aflow_title 'Scheelite (CaWO$_{4}$, $H0_{4}$) Structure'
_aflow_proto 'AB4C_tI24_88_b_f_a'
_aflow_params 'a,c/a,x_{3},y_{3},z_{3}'
_aflow_params_values '5.2429,2.16935283908,0.1488,0.0038,0.2159'
_aflow_Strukturbericht '$H0_{4}$'
_aflow_Pearson 'tI24'

_symmetry_space_group_name_H-M "I 41/a (origin choice 2)"
_symmetry_Int_Tables_number 88

_cell_length_a    5.24290
_cell_length_b    5.24290
_cell_length_c    11.37370
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y+1/2,z
3 -y+3/4,x+1/4,z+1/4
4 y+1/4,-x+1/4,z+1/4
5 -x,-y,-z
6 x,y+1/2,-z
7 y+1/4,-x+3/4,-z+3/4
8 -y+3/4,x+3/4,-z+3/4
9 x+1/2,y+1/2,z+1/2
10 -x+1/2,-y,z+1/2
11 -y+1/4,x+3/4,z+3/4
12 y+3/4,-x+3/4,z+3/4
13 -x+1/2,-y+1/2,-z+1/2
14 x+1/2,y,-z+1/2
15 y+3/4,-x+1/4,-z+1/4
16 -y+1/4,x+1/4,-z+1/4

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
W1 W 4 a 0.00000 0.25000 0.12500 1.00000
Ca1 Ca 4 b 0.00000 0.25000 0.62500 1.00000
O1 O 16 f 0.14880 0.00380 0.21590 1.00000
```

Scheelite ($CaWO_4$, $H0_4$): AB4C_tI24_88_b_f_a - POSCAR

```
AB4C_tI24_88_b_f_a & a,c/a,x3,y3,z3 --params=5.2429,2.16935283908,0.1488
    ↪ ,0.0038,0.2159 & I4_{1}/a C_{4h}^{6} #88 (abf) & tI24 & $H0_{4}
    ↪ $ & CaO4W & Scheelite & R. M. Hazen and L. W. Finger and J. W.
    ↪ E. Mariathasan, J. Phys. Chem. Solids 46, 253-263 (1985)
   1.00000000000000
  -2.62145000000000    2.62145000000000    5.68685000000000
   2.62145000000000   -2.62145000000000    5.68685000000000
   2.62145000000000    2.62145000000000   -5.68685000000000
    Ca     O     W
     2     8     2
Direct
   0.87500000000000    0.62500000000000    0.25000000000000   Ca    (4b)
   0.12500000000000    0.37500000000000    0.75000000000000   Ca    (4b)
   0.21970000000000    0.36470000000000    0.15260000000000   O    (16f)
   0.71210000000000    0.06710000000000    0.34740000000000   O    (16f)
   0.86470000000000    0.21210000000000    0.14500000000000   O    (16f)
   0.56710000000000    0.71970000000000    0.35500000000000   O    (16f)
  -0.21970000000000   -0.36470000000000   -0.15260000000000   O    (16f)
   0.28790000000000   -0.06710000000000    0.65260000000000   O    (16f)
   0.13530000000000   -0.21210000000000   -0.14500000000000   O    (16f)
   0.43290000000000    0.28030000000000    0.64500000000000   O    (16f)
   0.37500000000000    0.12500000000000    0.25000000000000   W    (4a)
   0.62500000000000    0.87500000000000    0.75000000000000   W    (4a)
```

$G7_5$ ($PbCO_3 \cdot PbCl_2$, Phosgenite) (*obsolete*): AB2C3D2_tP16_90_c_f_ce_e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Phosgenite'
_chemical_formula_sum 'C Cl2 O3 Pb2'

loop_
_publ_author_name
 'E. Onorato'
_journal_name_full_name
;
 Periodico di Mineralogia
;
_journal_volume 5
_journal_year 1934
_journal_page_first 1
_journal_page_last 27
_publ_Section_title
;
 La struttura della Fosgenite
;

# Found in Reexamination of the crystal structure of phosgenite, Pb$_{2}
    ↪ $Cl$_{2}$($CO$_{3}$), 1974

_aflow_title '$G7_{5}$ (PbCO$_{3}$ $\cdot$ PbCl$_{2}$, Phosgenite) ({\em
    ↪ {obsolete}}) Structure'
_aflow_proto 'AB2C3D2_tP16_90_c_f_ce_e'
_aflow_params 'a,c/a,z_{1},z_{2},x_{3},x_{4},x_{5}'
_aflow_params_values '8.12,0.541871921182,0.55,0.84,0.11,0.34,0.24'
_aflow_Strukturbericht '$G7_{5}$'
_aflow_Pearson 'tP16'

_symmetry_space_group_name_H-M "P 4 21 2"
_symmetry_Int_Tables_number 90

_cell_length_a    8.12000
_cell_length_b    8.12000
_cell_length_c    4.40000
_cell_angle_alpha 90.00000
```

```
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z
3 -x+1/2,y+1/2,-z
4 -x,-y,z
5 -y,-x,-z
6 -y+1/2,x+1/2,z
7 y+1/2,-x+1/2,z
8 y,x,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
C1 C 2 c 0.00000 0.50000 0.55000 1.00000
O1 O 2 c 0.00000 0.50000 0.84000 1.00000
O2 O 4 e 0.11000 0.11000 0.00000 1.00000
Pb1 Pb 4 e 0.34000 0.34000 0.00000 1.00000
Cl1 Cl 4 f 0.24000 0.24000 0.50000 1.00000
```

$G7_5$ ($PbCO_3 \cdot PbCl_2$, Phosgenite) (*obsolete*): AB2C3D2_tP16_90_c_f_ce_e - POSCAR

```
AB2C3D2_tP16_90_c_f_ce_e & a,c/a,z1,z2,x3,x4,x5 --params=8.12,
      ↪ 0.541871921182,0.55,0.84,0.11,0.34,0.24 & P42_{1}2 D_{4}^{2} #
      ↪ 90 (c^2e^2f) & tP16 & $G7_{5}$ & CCl2O3Pb2 & Phosgenite & E.
      ↪ Onorato, Period. Mineral. 5, 1-27 (1934)
   1.00000000000000
   8.12000000000000   0.00000000000000   0.00000000000000
   0.00000000000000   8.12000000000000   0.00000000000000
   0.00000000000000   0.00000000000000   4.40000000000000
     C    Cl     O    Pb
     2     4     6     4
Direct
   0.00000000000000   0.50000000000000   0.55000000000000    C    (2c)
   0.50000000000000   0.00000000000000  -0.55000000000000    C    (2c)
   0.24000000000000   0.24000000000000   0.50000000000000   Cl    (4f)
  -0.24000000000000  -0.24000000000000   0.50000000000000   Cl    (4f)
   0.26000000000000   0.74000000000000   0.50000000000000   Cl    (4f)
   0.74000000000000   0.26000000000000   0.50000000000000   Cl    (4f)
   0.00000000000000   0.50000000000000   0.84000000000000    O    (2c)
   0.50000000000000   0.00000000000000  -0.84000000000000    O    (2c)
   0.11000000000000   0.11000000000000   0.00000000000000    O    (4e)
  -0.11000000000000  -0.11000000000000   0.00000000000000    O    (4e)
   0.39000000000000   0.61000000000000   0.00000000000000    O    (4e)
   0.61000000000000   0.39000000000000   0.00000000000000    O    (4e)
   0.34000000000000   0.34000000000000   0.00000000000000   Pb    (4e)
  -0.34000000000000  -0.34000000000000   0.00000000000000   Pb    (4e)
   0.16000000000000   0.84000000000000   0.00000000000000   Pb    (4e)
   0.84000000000000   0.16000000000000   0.00000000000000   Pb    (4e)
```

Retgersite (α-$NiSO_4 \cdot 6H_2O$, $H4_5$): A12BC10D_tP96_92_6b_a_5b_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Retgersite'
_chemical_formula_sum 'H12 Ni O10 S'

loop_
_publ_author_name
 'K. Stadnicka'
 'A. M. Glazer'
 'M. Koralewski'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 43
_journal_year 1987
_journal_page_first 319
_journal_page_last 325
_publ_Section_title
;
 Structure, absolute configuration and optical activity of $\
      ↪ alpha$-nickel sulfate hexahydrate
;

# Found in Handbook of Mineralogy, 2004

_aflow_title 'Retgersite ($\alpha$-NiSO$_{4}$$\cdot$6H$_{2}$O, $H4_{5}$)
      ↪ Structure'
_aflow_proto 'A12BC10D_tP96_92_6b_a_5b_a'
_aflow_params 'a,c/a,x_{1},x_{2},x_{3},y_{3},z_{3},x_{4},y_{4},z_{4},x_{
      ↪ 5},y_{5},z_{5},x_{6},y_{6},z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},
      ↪ z_{8},x_{9},y_{9},z_{9},x_{10},y_{10},z_{10},x_{11},y_{11},z_{
      ↪ 11},x_{12},y_{12},z_{12},x_{13},y_{13},z_{13}'
_aflow_params_values '6.783,2.69615214507,0.2106,0.70943,0.1097,0.8667,
      ↪ 0.0394,0.2215,-0.0746,0.0867,0.5702,0.1559,0.0507,0.5371,0.3532
      ↪ ,0.0601,-0.0051,0.4453,0.0744,-0.0101,0.2941,0.1149,0.1727,-
      ↪ 0.047,0.0528,0.4705,0.2449,0.0561,0.0658,0.3599,0.085,0.6209,
      ↪ 0.6203,0.0658,0.9237,0.6731,0.0003'
_aflow_Strukturbericht '$H4_{5}$'
_aflow_Pearson 'tP96'
_symmetry_space_group_name_H-M "P 41 21 2"
_symmetry_Int_Tables_number 92

_cell_length_a      6.78300
_cell_length_b      6.78300
_cell_length_c      18.28800
_cell_angle_alpha   90.00000
_cell_angle_beta    90.00000
_cell_angle_gamma   90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+3/4
3 -x+1/2,y+1/2,-z+1/4
4 -x,-y,z+1/2
5 -y,-x,-z+1/2
6 -y+1/2,x+1/2,z+1/4
7 y+1/2,-x+1/2,z+3/4
8 y,x,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ni1 Ni 4 a 0.21060 0.21060 0.00000 1.00000
S1 S 4 a 0.70943 0.70943 0.00000 1.00000
H1 H 8 b 0.10970 0.86670 0.03940 1.00000
H2 H 8 b 0.22150 -0.07460 0.08670 1.00000
H3 H 8 b 0.57020 0.15590 0.05070 1.00000
H4 H 8 b 0.53710 0.35320 0.06010 1.00000
H5 H 8 b -0.00510 0.44530 0.07440 1.00000
H6 H 8 b -0.01010 0.29410 0.11490 1.00000
O1 O 8 b 0.17270 -0.04700 0.05280 1.00000
O2 O 8 b 0.47050 0.24490 0.05610 1.00000
O3 O 8 b 0.06580 0.35990 0.08500 1.00000
O4 O 8 b 0.62090 0.62030 0.06580 1.00000
O5 O 8 b 0.92370 0.67310 0.00030 1.00000
```

Retgersite (α-$NiSO_4 \cdot 6H_2O$, $H4_5$): A12BC10D_tP96_92_6b_a_5b_a - POSCAR

```
A12BC10D_tP96_92_6b_a_5b_a & a,c/a,x1,x2,x3,y3,z3,x4,y4,z4,x5,y5,z5,x6,
      ↪ y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10,x11,y11,z11,x12,
      ↪ y12,z12,x13,y13,z13 --params=6.783,2.69615214507,0.2106,0.70943
      ↪ ,0.1097,0.8667,0.0394,0.2215,-0.0746,0.0867,0.5702,0.1559,
      ↪ 0.0507,0.5371,0.3532,0.0601,-0.0051,0.4453,0.0744,-0.0101,
      ↪ 0.2941,0.1149,0.1727,-0.047,0.0528,0.4705,0.2449,0.0561,0.0658,
      ↪ 0.3599,0.085,0.6209,0.6203,0.0658,0.9237,0.6731,0.0003 & P4_{1}
      ↪ 2_{1}2 D_{4}^{4} #92 (a^2b^11) & tP96 & $H4_{5}$ & H12NiO10S &
      ↪ Retgersite & K. Stadnicka and A. M. Glazer and M. Koralewski,
      ↪ Acta Crystallogr. Sect. B Struct. Sci. 43, 319-325 (1987)
   1.00000000000000
   6.78300000000000   0.00000000000000   0.00000000000000
   0.00000000000000   6.78300000000000   0.00000000000000
   0.00000000000000   0.00000000000000  18.28800000000000
     H    Ni     O     S
    48     4    40     4
Direct
   0.10970000000000   0.86670000000000   0.03940000000000    H    (8b)
  -0.10970000000000  -0.86670000000000   0.53940000000000    H    (8b)
  -0.36670000000000   0.60970000000000   0.28940000000000    H    (8b)
   1.36670000000000   0.39030000000000   0.78940000000000    H    (8b)
   0.39030000000000   1.36670000000000   0.21060000000000    H    (8b)
   0.60970000000000  -0.36670000000000   0.71060000000000    H    (8b)
   0.86670000000000   0.10970000000000  -0.03940000000000    H    (8b)
  -0.86670000000000  -0.10970000000000   0.46060000000000    H    (8b)
   0.22150000000000  -0.07460000000000   0.08670000000000    H    (8b)
  -0.22150000000000   0.07460000000000   0.58670000000000    H    (8b)
   0.57460000000000   0.72150000000000   0.33670000000000    H    (8b)
   0.42540000000000   0.27850000000000   0.83670000000000    H    (8b)
   0.27850000000000   0.42540000000000   0.16330000000000    H    (8b)
   0.72150000000000   0.57460000000000   0.66330000000000    H    (8b)
  -0.07460000000000   0.22150000000000  -0.08670000000000    H    (8b)
   0.07460000000000  -0.22150000000000   0.41330000000000    H    (8b)
   0.57020000000000   0.15590000000000   0.05070000000000    H    (8b)
  -0.57020000000000  -0.15590000000000   0.55070000000000    H    (8b)
   0.34410000000000   1.07020000000000   0.30070000000000    H    (8b)
   0.65590000000000  -0.07020000000000   0.80070000000000    H    (8b)
  -0.07020000000000   0.65590000000000   0.19930000000000    H    (8b)
   1.07020000000000   0.34410000000000   0.69930000000000    H    (8b)
   0.15590000000000   0.57020000000000  -0.05070000000000    H    (8b)
  -0.15590000000000  -0.57020000000000   0.44930000000000    H    (8b)
   0.53710000000000   0.35320000000000   0.06010000000000    H    (8b)
  -0.53710000000000  -0.35320000000000   0.56010000000000    H    (8b)
   0.14680000000000   1.03710000000000   0.31010000000000    H    (8b)
   0.85320000000000  -0.03710000000000   0.81010000000000    H    (8b)
  -0.03710000000000   0.85320000000000   0.18990000000000    H    (8b)
   1.03710000000000   0.14680000000000   0.68990000000000    H    (8b)
   0.35320000000000   0.53710000000000  -0.06010000000000    H    (8b)
  -0.35320000000000  -0.53710000000000   0.43990000000000    H    (8b)
  -0.00510000000000   0.44530000000000   0.07440000000000    H    (8b)
   0.00510000000000  -0.44530000000000   0.57440000000000    H    (8b)
   0.05470000000000   0.49490000000000   0.32440000000000    H    (8b)
   0.94530000000000   0.50510000000000   0.82440000000000    H    (8b)
   0.50510000000000   0.94530000000000   0.17560000000000    H    (8b)
   0.49490000000000   0.05470000000000   0.67560000000000    H    (8b)
   0.44530000000000  -0.00510000000000  -0.07440000000000    H    (8b)
  -0.44530000000000   0.00510000000000   0.42560000000000    H    (8b)
  -0.01010000000000   0.29410000000000   0.11490000000000    H    (8b)
```

```
   0.01010000000000  −0.29410000000000   0.61490000000000   H   (8b)
   0.20590000000000   0.48990000000000   0.36490000000000   H   (8b)
   0.79410000000000   0.51010000000000   0.86490000000000   H   (8b)
   0.51010000000000   0.79410000000000   0.13510000000000   H   (8b)
   0.48990000000000   0.20590000000000   0.63510000000000   H   (8b)
   0.29410000000000  −0.01010000000000  −0.11490000000000   H   (8b)
  −0.29410000000000   0.01010000000000   0.38510000000000   H   (8b)
   0.21060000000000   0.21060000000000   0.00000000000000   Ni  (4a)
  −0.21060000000000  −0.21060000000000   0.50000000000000   Ni  (4a)
   0.28940000000000   0.71060000000000   0.25000000000000   Ni  (4a)
   0.71060000000000   0.28940000000000   0.75000000000000   Ni  (4a)
   0.17270000000000  −0.04700000000000   0.05280000000000   O   (8b)
  −0.17270000000000   0.04700000000000   0.55280000000000   O   (8b)
   0.54700000000000   0.67270000000000   0.30280000000000   O   (8b)
   0.45300000000000   0.32730000000000   0.80280000000000   O   (8b)
   0.32730000000000   0.45300000000000   0.19720000000000   O   (8b)
   0.67270000000000   0.54700000000000   0.69720000000000   O   (8b)
  −0.04700000000000   0.17270000000000  −0.05280000000000   O   (8b)
   0.04700000000000  −0.17270000000000   0.44720000000000   O   (8b)
   0.47050000000000   0.24490000000000   0.05610000000000   O   (8b)
  −0.47050000000000  −0.24490000000000   0.55610000000000   O   (8b)
   0.25510000000000   0.97050000000000   0.30610000000000   O   (8b)
   0.74490000000000   0.02950000000000   0.80610000000000   O   (8b)
   0.02950000000000   0.74490000000000   0.19390000000000   O   (8b)
   0.97050000000000   0.25510000000000   0.69390000000000   O   (8b)
   0.24490000000000   0.47050000000000  −0.05610000000000   O   (8b)
  −0.24490000000000  −0.47050000000000   0.44390000000000   O   (8b)
   0.06580000000000   0.35990000000000   0.08500000000000   O   (8b)
  −0.06580000000000  −0.35990000000000   0.58500000000000   O   (8b)
   0.14010000000000   0.56580000000000   0.33500000000000   O   (8b)
   0.85990000000000   0.43420000000000   0.83500000000000   O   (8b)
   0.43420000000000   0.85990000000000   0.16500000000000   O   (8b)
   0.56580000000000   0.14010000000000   0.66500000000000   O   (8b)
   0.35990000000000   0.06580000000000  −0.08500000000000   O   (8b)
  −0.35990000000000  −0.06580000000000   0.41500000000000   O   (8b)
   0.62090000000000   0.62030000000000   0.06580000000000   O   (8b)
  −0.62090000000000  −0.62030000000000   0.56580000000000   O   (8b)
  −0.12030000000000   1.12090000000000   0.31580000000000   O   (8b)
   1.12030000000000  −0.12090000000000   0.81580000000000   O   (8b)
  −0.12090000000000   1.12030000000000   0.18420000000000   O   (8b)
   1.12090000000000  −0.12030000000000   0.68420000000000   O   (8b)
   0.62030000000000   0.62090000000000  −0.06580000000000   O   (8b)
  −0.62030000000000  −0.62090000000000   0.43420000000000   O   (8b)
   0.92370000000000   0.67310000000000   0.00030000000000   O   (8b)
  −0.92370000000000  −0.67310000000000   0.50030000000000   O   (8b)
  −0.17310000000000   1.42370000000000   0.25030000000000   O   (8b)
   1.17310000000000  −0.42370000000000   0.75030000000000   O   (8b)
  −0.42370000000000   1.17310000000000   0.24970000000000   O   (8b)
   1.42370000000000  −0.17310000000000   0.74970000000000   O   (8b)
   0.67310000000000   0.92370000000000  −0.00030000000000   O   (8b)
  −0.67310000000000  −0.92370000000000   0.49970000000000   O   (8b)
   0.70943000000000   0.70943000000000   0.00000000000000   S   (4a)
  −0.70943000000000  −0.70943000000000   0.50000000000000   S   (4a)
  −0.20943000000000   1.20943000000000   0.25000000000000   S   (4a)
   1.20943000000000  −0.20943000000000   0.75000000000000   S   (4a)
```

Paratellurite (α-TeO_2): A2B_tP12_92_b_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Paratellurite'
_chemical_formula_sum 'O2 Te'

loop_
_publ_author_name
 'P. A. Thomas'
_journal_name_full_name
;
 Journal of Physics C: Solid State Physics
;
_journal_volume 21
_journal_year 1988
_journal_page_first 4611
_journal_page_last 4627
_publ_Section_title
;
 The crystal structure and absolute optical chirality of paratellurite,
      ↪ $\alpha$-TeO$_{2}$
;

# Found in {\em Ab initio} study of the vibrational properties of
      ↪ crystalline TeO$_{2}$: The $\alpha$, $\beta$, and $\gamma$
      ↪ phases, 2006

_aflow_title 'Paratellurite ($\alpha$-TeO$_{2}$) Structure'
_aflow_proto 'A2B_tP12_92_b_a'
_aflow_params 'a,c/a,x_{1},x_{2},y_{2},z_{2}'
_aflow_params_values '4.8082,1.60392662535,0.0268,0.1386,0.2576,0.1862'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tP12'

_symmetry_space_group_name_H-M "P 41 21 2"
_symmetry_Int_Tables_number 92

_cell_length_a     4.80820
_cell_length_b     4.80820
_cell_length_c     7.71200
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+3/4
3 -x+1/2,y+1/2,-z+1/4
4 -x,-y,z+1/2
5 -y,-x,-z+1/2
6 -y+1/2,x+1/2,z+1/4
7 y+1/2,-x+1/2,z+3/4
8 y,x,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Te1 Te 4 a 0.02680 0.02680 0.00000 1.00000
O1 O 8 b 0.13860 0.25760 0.18620 1.00000
```

Paratellurite (α-TeO_2): A2B_tP12_92_b_a - POSCAR

```
A2B_tP12_92_b_a & a,c/a,x1,x2,y2,z2 --params=4.8082,1.60392662535,0.0268
      ↪ ,0.1386,0.2576,0.1862 & P4_{1}2_{1}2 D_{4}^{4} #92 (ab) & tP12
      ↪ & None & O2Te & Paratellurite & P. A. Thomas, J. Phys. C: Solid
      ↪ State Phys. 21, 4611-4627 (1988)
   1.00000000000000
   4.80820000000000   0.00000000000000   0.00000000000000
   0.00000000000000   4.80820000000000   0.00000000000000
   0.00000000000000   0.00000000000000   7.71200000000000
     O     Te
     8     4
Direct
   0.13860000000000   0.25760000000000   0.18620000000000   O   (8b)
  -0.13860000000000  -0.25760000000000   0.68620000000000   O   (8b)
   0.24240000000000   0.63860000000000   0.43620000000000   O   (8b)
   0.75760000000000   0.36140000000000   0.93620000000000   O   (8b)
   0.36140000000000   0.75760000000000   0.06380000000000   O   (8b)
   0.63860000000000   0.24240000000000   0.56380000000000   O   (8b)
   0.25760000000000   0.13860000000000  -0.18620000000000   O   (8b)
  -0.25760000000000  -0.13860000000000   0.31380000000000   O   (8b)
   0.02680000000000   0.02680000000000   0.00000000000000   Te  (4a)
  -0.02680000000000  -0.02680000000000   0.50000000000000   Te  (4a)
   0.47320000000000   0.52680000000000   0.25000000000000   Te  (4a)
   0.52680000000000   0.47320000000000   0.75000000000000   Te  (4a)
```

Phase III $Cd_2Re_2O_7$: A2B7C2_tI44_98_f_bcde_f - CIF

```
# CIF file
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_audit_creation_method FINDSYM

_chemical_name_mineral 'Cd2O7Re2'
_chemical_formula_sum 'Cd2 O7 Re2'

loop_
_publ_author_name
 'S.-W. Huang'
 'H.-T. Jeng'
 'J.-Y. Lin'
 'W. J. Chang'
 'J. M. Chen'
 'G. H. Lee'
 'H. Berger'
 'H. D. Yang'
 'K. S. Liang'
_journal_name_full_name
;
 Journal of Physics: Condensed Matter
;
_journal_volume 21
_journal_year 2009
_journal_page_first 195602
_journal_page_last 195602
_publ_Section_title
;
 Electronic structure of pyrochlore Cd$_{2}$Re$_{2}$O$_{7}$
;

# Found in The crystal structure of the inversion breaking metal Cd$_{2}
      ↪ $Re$_{2}$O$_{7}$, {arXiv:1911.10141 [cond-mat.str-el]}, 2019

_aflow_title 'Phase III Cd$_{2}$Re$_{2}$O$_{7}$ Structure'
_aflow_proto 'A2B7C2_tI44_98_f_bcde_f'
_aflow_params 'a,c/a,z_{2},x_{3},x_{4},x_{5},x_{6}'
_aflow_params_values '7.2313,1.41443447236,0.81048,0.188,0.803,0.5041,-
      ↪ 0.0033'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tI44'

_symmetry_space_group_name_H-M "I 41 2 2"
_symmetry_Int_Tables_number 98

_cell_length_a     7.23130
_cell_length_b     7.23130
_cell_length_c     10.22820
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
```

```
1 x,y,z
2 x,-y+1/2,-z+1/4
3 -x,y+1/2,-z+1/4
4 -x,-y,z
5 -y,-x,-z
6 -y,x+1/2,z+1/4
7 y,-x+1/2,z+1/4
8 y,x,-z
9 x+1/2,y+1/2,z+1/2
10 x+1/2,-y,-z+3/4
11 -x+1/2,y,-z+3/4
12 -x+1/2,-y+1/2,z+1/2
13 -y+1/2,-x+1/2,-z+1/2
14 -y+1/2,x,z+3/4
15 y+1/2,-x,z+3/4
16 y+1/2,x+1/2,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 4 b 0.00000 0.00000 0.50000 1.00000
O2 O 8 c 0.00000 0.00000 0.81048 1.00000
O3 O 8 d 0.18800 0.18800 0.00000 1.00000
O4 O 8 e 0.80300 0.19700 0.00000 1.00000
Cd1 Cd 8 f 0.50410 0.25000 0.12500 1.00000
Re1 Re 8 f -0.00330 0.25000 0.12500 1.00000
```

Phase III $Cd_2Re_2O_7$: A2B7C2_tI44_98_f_bcde_f - POSCAR

```
A2B7C2_tI44_98_f_bcde_f & a,c/a,z2,x3,x4,x5,x6 --params=7.2313,
      ↪ 1.41443447236,0.81048,0.188,0.803,0.5041,-0.0033 & I4_{1}22 D_{
      ↪ 4}^{10} #98 (bcdef^2) & tI44 & None & Cd2O7Re2 & Cd2O7Re2 &
      ↪ S.-W. Huang et al., J. Phys.: Condens. Matter 21, 195602(2009)
   1.00000000000000
  -3.61565000000000    3.61565000000000    5.11410000000000
   3.61565000000000   -3.61565000000000    5.11410000000000
   3.61565000000000    3.61565000000000   -5.11410000000000
    Cd      O     Re
     4     14      4
Direct
   0.37500000000000    0.62910000000000    0.75410000000000   Cd    (8f)
   0.87500000000000   -0.37910000000000    0.24590000000000   Cd    (8f)
   1.37910000000000    0.12500000000000    0.75410000000000   Cd    (8f)
   0.37090000000000    0.62500000000000    0.24590000000000   Cd    (8f)
   0.50000000000000    0.50000000000000    0.00000000000000   O     (4b)
   0.25000000000000    0.75000000000000    0.50000000000000   O     (4b)
   0.81048000000000    0.81048000000000    0.00000000000000   O     (8c)
   1.56048000000000    1.06048000000000    0.50000000000000   O     (8c)
  -0.06048000000000   -0.56048000000000    0.50000000000000   O     (8c)
  -0.81048000000000   -0.81048000000000    0.00000000000000   O     (8c)
   0.18800000000000    0.18800000000000    0.37600000000000   O     (8d)
  -0.18800000000000   -0.18800000000000   -0.37600000000000   O     (8d)
   0.93800000000000    0.06200000000000    0.50000000000000   O     (8d)
   0.56200000000000    0.43800000000000    0.50000000000000   O     (8d)
   0.80300000000000   -0.80300000000000    0.00000000000000   O     (8e)
  -0.80300000000000    0.80300000000000    0.00000000000000   O     (8e)
  -0.05300000000000   -0.55300000000000   -1.10600000000000   O     (8e)
   1.55300000000000    1.05300000000000    2.10600000000000   O     (8e)
   0.37500000000000    0.12170000000000    0.24670000000000   Re    (8f)
   0.87500000000000    0.12830000000000    0.75330000000000   Re    (8f)
   0.87170000000000    0.12500000000000    0.24670000000000   Re    (8f)
   0.87830000000000    0.62500000000000    0.75330000000000   Re    (8f)
```

$F5_4$ (NH_4ClO_2) (*obsolete*): ABC2_tP8_100_b_a_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Cl(NH4)O2'
_chemical_formula_sum 'Cl (NH4) O2'

loop_
_publ_author_name
 'G. R. Levi'
 'A. Scherillo'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 76
_journal_year 1931
_journal_page_first 431
_journal_page_last 452
_publ_Section_title
;
 Ricerche cristallografiche sui sali dell\'acido cloroso
;

# Found in Strukturbericht Band II 1928-1932, 1937

_aflow_title '$F5_{4}$ (NH$_{4}$ClO$_{2}$) ({\em{obsolete}}) Structure'
_aflow_proto 'ABC2_tP8_100_b_a_c'
_aflow_params 'a,c/a,z_{1},z_{2},x_{3},z_{3}'
_aflow_params_values '6.3,0.587301587302,0.0,0.25,0.65,0.5'
_aflow_Strukturbericht '$F5_{4}$'
_aflow_Pearson 'tP8'

_symmetry_space_group_name_H-M "P 4 b m"
_symmetry_Int_Tables_number 100

_cell_length_a     6.30000
_cell_length_b     6.30000
_cell_length_c     3.70000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,z
3 -y,x,z
4 y,-x,z
5 -x+1/2,y+1/2,z
6 x+1/2,-y+1/2,z
7 y+1/2,x+1/2,z
8 -y+1/2,-x+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
NH41 NH4 2 a 0.00000 0.00000 0.00000 1.00000
Cl1 Cl 2 b 0.50000 0.00000 0.25000 1.00000
O1 O 4 c 0.65000 0.15000 0.50000 1.00000
```

$F5_4$ (NH_4ClO_2) (*obsolete*): ABC2_tP8_100_b_a_c - POSCAR

```
ABC2_tP8_100_b_a_c & a,c/a,z1,z2,x3,z3 --params=6.3,0.587301587302,0.0,
      ↪ 0.25,0.65,0.5 & P4bm C_{4v}^{2} #100 (abc) & tP8 & $F5_{4}$ &
      ↪ Cl(NH4)O2 & Cl(NH4)O2 & G. R. Levi and A. Scherillo,
      ↪ Zeitschrift f{"u}r Kristallographie - Crystalline Materials 76,
      ↪ 431-452 (1931)
   1.00000000000000
   6.30000000000000    0.00000000000000    0.00000000000000
   0.00000000000000    6.30000000000000    0.00000000000000
   0.00000000000000    0.00000000000000    3.70000000000000
    Cl    NH4     O
     2      2     4
Direct
   0.50000000000000    0.00000000000000    0.25000000000000   Cl    (2b)
   0.00000000000000    0.50000000000000    0.25000000000000   Cl    (2b)
   0.00000000000000    0.00000000000000    0.00000000000000   NH4   (2a)
   0.50000000000000    0.50000000000000    0.00000000000000   NH4   (2a)
   0.65000000000000    1.15000000000000    0.50000000000000   O     (4c)
  -0.65000000000000   -0.15000000000000    0.50000000000000   O     (4c)
  -0.15000000000000    0.65000000000000    0.50000000000000   O     (4c)
   1.15000000000000   -0.65000000000000    0.50000000000000   O     (4c)
```

NH_4NO_3 II ($G0_9$): ABC3_tP10_100_b_a_bc - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'N(NH4)O3'
_chemical_formula_sum 'N (NH4) O3'

loop_
_publ_author_name
 'Y. Shinnaka'
_journal_name_full_name
;
 Journal of the Physical Society of Japan
;
_journal_volume 11
_journal_year 1956
_journal_page_first 393
_journal_page_last 396
_publ_Section_title
;
 On the Metastable Transition and the Crystal Structure of Ammonium
      ↪ Nitrate (Tetragonal Modification)
;

# Found in The structure of ammonium nitrate (IV), 1972

_aflow_title 'NH$_{4}$NO$_{3}$ II ($G0_{9}$) Structure'
_aflow_proto 'ABC3_tP10_100_b_a_bc'
_aflow_params 'a,c/a,z_{1},z_{2},z_{3},x_{4},z_{4}'
_aflow_params_values '5.74,0.862369337979,0.0,0.53,0.75,0.62,0.42'
_aflow_Strukturbericht '$G0_{9}$'
_aflow_Pearson 'tP10'

_symmetry_space_group_name_H-M "P 4 b m"
_symmetry_Int_Tables_number 100

_cell_length_a     5.74000
_cell_length_b     5.74000
_cell_length_c     4.95000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
```

```
1 x,y,z
2 -x,-y,z
3 -y,x,z
4 y,-x,z
5 -x+1/2,y+1/2,z
6 x+1/2,-y+1/2,z
7 y+1/2,x+1/2,z
8 -y+1/2,-x+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
NH41 NH4 2 a 0.00000 0.00000 0.00000 1.00000
N1 N 2 b 0.50000 0.00000 0.53000 1.00000
O1 O 2 b 0.50000 0.00000 0.75000 1.00000
O2 O 4 c 0.62000 0.12000 0.42000 1.00000
```

NH_4NO_3 II ($G0_9$): ABC3_tP10_100_b_a_bc - POSCAR

```
ABC3_tP10_100_b_a_bc & a,c/a,z1,z2,z3,x4,z4 --params=5.74,0.862369337979
      ↪ ,0.0,0.53,0.75,0.62,0.42 & P4bm C_{4v}^{2} #100 (ab^2c) & tP10
      ↪ & $G0_{9}$ & N(NH4)O3 & N(NH4)O3 & Y. Shinnaka, J. Phys. Soc.
      ↪ Jpn. 11, 393-396 (1956)
   1.00000000000000
   5.74000000000000    0.00000000000000    0.00000000000000
   0.00000000000000    5.74000000000000    0.00000000000000
   0.00000000000000    0.00000000000000    4.95000000000000
     N    NH4     O
     2      2     6
Direct
   0.50000000000000    0.00000000000000    0.53000000000000     N    (2b)
   0.00000000000000    0.50000000000000    0.53000000000000     N    (2b)
   0.00000000000000    0.00000000000000    0.00000000000000   NH4    (2a)
   0.50000000000000    0.50000000000000    0.00000000000000   NH4    (2a)
   0.50000000000000    0.00000000000000    0.75000000000000     O    (2b)
   0.00000000000000    0.50000000000000    0.75000000000000     O    (2b)
   0.62000000000000    1.12000000000000    0.42000000000000     O    (4c)
  -0.62000000000000   -0.12000000000000    0.42000000000000     O    (4c)
  -0.12000000000000    0.62000000000000    0.42000000000000     O    (4c)
   1.12000000000000   -0.62000000000000    0.42000000000000     O    (4c)
```

VSe_2O_6: A6B2C_tP72_103_abc5d_2d_abc - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'O6Se2V'
_chemical_formula_sum 'O6 Se2 V'

loop_
_publ_author_name
 'G. Meunier'
 'M. Bertaud'
 'J. Galy'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 30
_journal_year 1974
_journal_page_first 2834
_journal_page_last 2839
_publ_Section_title
;
 Cristallochimie du s\\'{e}l\\'{e}nium(+IV). I. VSe$_{2}$O$_{6}$, une
      ↪ structure \'{a} trois cha\^{i}nes parall\'{e}les (VO$_{5}$)$^{
      ↪ 6n-}_{n}$ ind\\'{e}pendantes pont\\'{e}es par des groupements
      ↪ (Se$_{2}$O)$^{6+}$
;

_aflow_title 'VSe$_{2}$O$_{6}$ Structure'
_aflow_proto 'A6B2C_tP72_103_abc5d_2d_abc'
_aflow_params 'a,c/a,z_{1},z_{2},z_{3},z_{4},z_{5},z_{6},x_{7},y_{7},z_{
      ↪ 7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{10},y_{10},z_{10},x_{
      ↪ 11},y_{11},z_{11},x_{12},y_{12},z_{12},x_{13},y_{13},z_{13}'
_aflow_params_values '11.22,0.700534759358,0.377,0.1715,0.223,0.4274,
      ↪ 0.391,0.0972,0.1233,0.12,0.13,0.2546,0.255,0.44,0.0919,0.3511,
      ↪ 0.146,0.3271,0.4759,0.468,0.4059,0.1432,0.15,0.1033,0.2405,0.0,
      ↪ 0.3671,0.2818,0.1031'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tP72'

_symmetry_space_group_name_H-M "P 4 c c"
_symmetry_Int_Tables_number 103

_cell_length_a     11.22000
_cell_length_b     11.22000
_cell_length_c     7.86000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,z
3 -y,x,z
4 y,-x,z
5 -x,y,z+1/2
6 x,-y,z+1/2
7 y,x,z+1/2
8 -y,-x,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 2 a 0.00000 0.00000 0.37700 1.00000
V1 V 2 a 0.00000 0.00000 0.17150 1.00000
O2 O 2 b 0.50000 0.50000 0.22300 1.00000
V2 V 2 b 0.50000 0.50000 0.42740 1.00000
O3 O 4 c 0.00000 0.50000 0.39100 1.00000
V3 V 4 c 0.00000 0.50000 0.09720 1.00000
O4 O 8 d 0.12330 0.12000 0.13000 1.00000
O5 O 8 d 0.25460 0.25500 0.44000 1.00000
O6 O 8 d 0.09190 0.35110 0.14600 1.00000
O7 O 8 d 0.32710 0.47590 0.46800 1.00000
O8 O 8 d 0.40590 0.14320 0.15000 1.00000
Se1 Se 8 d 0.10330 0.24050 0.00000 1.00000
Se2 Se 8 d 0.36710 0.28180 0.10310 1.00000
```

VSe_2O_6: A6B2C_tP72_103_abc5d_2d_abc - POSCAR

```
A6B2C_tP72_103_abc5d_2d_abc & a,c/a,z1,z2,z3,z4,z5,z6,x7,y7,z7,x8,y8,z8,
      ↪ x9,y9,z9,x10,y10,z10,x11,y11,z11,x12,y12,z12,x13,y13,z13 --
      ↪ params=11.22,0.700534759358,0.377,0.1715,0.223,0.4274,0.391,
      ↪ 0.0972,0.1233,0.12,0.13,0.2546,0.255,0.44,0.0919,0.3511,0.146,
      ↪ 0.3271,0.4759,0.468,0.4059,0.1432,0.15,0.1033,0.2405,0.0,0.3671
      ↪ ,0.2818,0.1031 & P4cc C_{4v}^{5} #103 (a^2b^2c^2d^7) & tP72 &
      ↪ None & O6Se2V & O6Se2V & G. Meunier and M. Bertaud and J. Galy,
      ↪  Acta Crystallogr. Sect. B Struct. Sci. 30, 2834-2839 (1974)
  1.00000000000000
 11.22000000000000    0.00000000000000    0.00000000000000
  0.00000000000000   11.22000000000000    0.00000000000000
  0.00000000000000    0.00000000000000    7.86000000000000
    O     Se     V
   48     16     8
Direct
   0.00000000000000    0.00000000000000    0.37700000000000     O    (2a)
   0.00000000000000    0.00000000000000    0.87700000000000     O    (2a)
   0.50000000000000    0.50000000000000    0.22300000000000     O    (2b)
   0.50000000000000    0.50000000000000    0.72300000000000     O    (2b)
   0.00000000000000    0.50000000000000    0.39100000000000     O    (4c)
   0.50000000000000    0.00000000000000    0.39100000000000     O    (4c)
   0.00000000000000    0.50000000000000    0.89100000000000     O    (4c)
   0.50000000000000    0.00000000000000    0.89100000000000     O    (4c)
   0.12330000000000    0.12000000000000    0.13000000000000     O    (8d)
  -0.12330000000000   -0.12000000000000    0.13000000000000     O    (8d)
  -0.12000000000000    0.12330000000000    0.13000000000000     O    (8d)
   0.12000000000000   -0.12330000000000    0.13000000000000     O    (8d)
   0.12330000000000   -0.12000000000000    0.63000000000000     O    (8d)
  -0.12330000000000    0.12000000000000    0.63000000000000     O    (8d)
  -0.12000000000000   -0.12330000000000    0.63000000000000     O    (8d)
   0.12000000000000    0.12330000000000    0.63000000000000     O    (8d)
   0.25460000000000    0.25500000000000    0.44000000000000     O    (8d)
  -0.25460000000000   -0.25500000000000    0.44000000000000     O    (8d)
  -0.25500000000000    0.25460000000000    0.44000000000000     O    (8d)
   0.25500000000000   -0.25460000000000    0.44000000000000     O    (8d)
   0.25460000000000   -0.25500000000000    0.94000000000000     O    (8d)
  -0.25460000000000    0.25500000000000    0.94000000000000     O    (8d)
  -0.25500000000000   -0.25460000000000    0.94000000000000     O    (8d)
   0.25500000000000    0.25460000000000    0.94000000000000     O    (8d)
   0.09190000000000    0.35110000000000    0.14600000000000     O    (8d)
  -0.09190000000000   -0.35110000000000    0.14600000000000     O    (8d)
  -0.35110000000000    0.09190000000000    0.14600000000000     O    (8d)
   0.35110000000000   -0.09190000000000    0.14600000000000     O    (8d)
   0.09190000000000   -0.35110000000000    0.64600000000000     O    (8d)
  -0.09190000000000    0.35110000000000    0.64600000000000     O    (8d)
  -0.35110000000000   -0.09190000000000    0.64600000000000     O    (8d)
   0.35110000000000    0.09190000000000    0.64600000000000     O    (8d)
   0.32710000000000    0.47590000000000    0.46800000000000     O    (8d)
  -0.32710000000000   -0.47590000000000    0.46800000000000     O    (8d)
  -0.47590000000000    0.32710000000000    0.46800000000000     O    (8d)
   0.47590000000000   -0.32710000000000    0.46800000000000     O    (8d)
   0.32710000000000   -0.47590000000000    0.96800000000000     O    (8d)
  -0.32710000000000    0.47590000000000    0.96800000000000     O    (8d)
  -0.47590000000000   -0.32710000000000    0.96800000000000     O    (8d)
   0.47590000000000    0.32710000000000    0.96800000000000     O    (8d)
   0.40590000000000    0.14320000000000    0.15000000000000     O    (8d)
  -0.40590000000000   -0.14320000000000    0.15000000000000     O    (8d)
  -0.14320000000000    0.40590000000000    0.15000000000000     O    (8d)
   0.14320000000000   -0.40590000000000    0.15000000000000     O    (8d)
   0.40590000000000   -0.14320000000000    0.65000000000000     O    (8d)
  -0.40590000000000    0.14320000000000    0.65000000000000     O    (8d)
  -0.14320000000000   -0.40590000000000    0.65000000000000     O    (8d)
   0.14320000000000    0.40590000000000    0.65000000000000     O    (8d)
   0.10330000000000    0.24050000000000    0.00000000000000    Se    (8d)
  -0.10330000000000   -0.24050000000000    0.00000000000000    Se    (8d)
  -0.24050000000000    0.10330000000000    0.00000000000000    Se    (8d)
   0.24050000000000   -0.10330000000000    0.00000000000000    Se    (8d)
   0.10330000000000   -0.24050000000000    0.50000000000000    Se    (8d)
  -0.10330000000000    0.24050000000000    0.50000000000000    Se    (8d)
  -0.24050000000000   -0.10330000000000    0.50000000000000    Se    (8d)
   0.24050000000000    0.10330000000000    0.50000000000000    Se    (8d)
   0.36710000000000    0.28180000000000    0.10310000000000    Se    (8d)
  -0.36710000000000   -0.28180000000000    0.10310000000000    Se    (8d)
  -0.28180000000000    0.36710000000000    0.10310000000000    Se    (8d)
```

```
  0.28180000000000  −0.36710000000000   0.10310000000000   Se   (8d)
  0.36710000000000  −0.28180000000000   0.60310000000000   Se   (8d)
 −0.36710000000000   0.28180000000000   0.60310000000000   Se   (8d)
 −0.28180000000000  −0.36710000000000   0.60310000000000   Se   (8d)
  0.28180000000000   0.36710000000000   0.60310000000000   Se   (8d)
  0.00000000000000   0.00000000000000   0.17150000000000   V    (2a)
  0.00000000000000   0.00000000000000   0.67150000000000   V    (2a)
  0.50000000000000   0.50000000000000   0.42740000000000   V    (2b)
  0.50000000000000   0.50000000000000   0.92740000000000   V    (2b)
  0.00000000000000   0.50000000000000   0.09720000000000   V    (4c)
  0.50000000000000   0.00000000000000   0.09720000000000   V    (4c)
  0.00000000000000   0.50000000000000   0.59720000000000   V    (4c)
  0.50000000000000   0.00000000000000   0.59720000000000   V    (4c)
```

BaNiSn$_3$: ABC3_tI10_107_a_a_ab - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'BaNiSn3'
_chemical_formula_sum 'Ba Ni Sn3'

loop_
_publ_author_name
 'W. D\"{o}rrscheidt'
 'H. Sch\"{a}fer'
_journal_name_full_name
;
 Journal of the Less-Common Metals
;
_journal_volume 58
_journal_year 1978
_journal_page_first 209
_journal_page_last 216
_publ_Section_title
;
 Die Struktur des BaPtSn$_{3}$, BaNiSn$_{3}$ und SrNiSn$_{3}$ und ihre
    ↪ Verwandtschaft zum ThCr$_{2}$Si$_{2}$-Strukturtyp
;

_aflow_title 'BaNiSn$_{3}$ Structure'
_aflow_proto 'ABC3_tI10_107_a_a_ab'
_aflow_params 'a,c/a,z_{1},z_{2},z_{3},z_{4}'
_aflow_params_values '4.82,2.26763692946,0.0,0.6554,0.4241,0.2516'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tI10'

_symmetry_space_group_name_H-M "I 4 m m"
_symmetry_Int_Tables_number 107

_cell_length_a     4.82000
_cell_length_b     4.82000
_cell_length_c     10.93001
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,z
3 -y,x,z
4 y,-x,z
5 -x,y,z
6 x,-y,z
7 y,x,z
8 -y,-x,z
9 x+1/2,y+1/2,z+1/2
10 -x+1/2,-y+1/2,z+1/2
11 -y+1/2,x+1/2,z+1/2
12 y+1/2,-x+1/2,z+1/2
13 -x+1/2,y+1/2,z+1/2
14 x+1/2,-y+1/2,z+1/2
15 y+1/2,x+1/2,z+1/2
16 -y+1/2,-x+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ba1 Ba 2 a 0.00000 0.00000 0.00000 1.00000
Ni1 Ni 2 a 0.00000 0.00000 0.65540 1.00000
Sn1 Sn 2 a 0.00000 0.00000 0.42410 1.00000
Sn2 Sn 4 b 0.00000 0.50000 0.25160 1.00000
```

BaNiSn$_3$: ABC3_tI10_107_a_a_ab - POSCAR

```
ABC3_tI10_107_a_a_ab & a,c/a,z1,z2,z3,z4 --params=4.82,2.26763692946,0.0
    ↪ ,0.6554,0.4241,0.2516 & I4mm C_{4v}^{9} #107 (a^3b) & tI10 &
    ↪ None & BaNiSn3 & BaNiSn3 & W. D\"{o}rrscheidt and H. Sch\"{a}
    ↪ fer, J. Less-Common Met. 58, 209-216 (1978)
   1.00000000000000
  -2.41000000000000   2.41000000000000   5.46500500000000
   2.41000000000000  -2.41000000000000   5.46500500000000
   2.41000000000000   2.41000000000000  -5.46500500000000
    Ba    Ni    Sn
     1     1     3
Direct
```

```
  0.00000000000000   0.00000000000000   0.00000000000000   Ba   (2a)
  0.65540000000000   0.65540000000000   0.00000000000000   Ni   (2a)
  0.42410000000000   0.42410000000000   0.00000000000000   Sn   (2a)
  0.75160000000000   0.25160000000000   0.50000000000000   Sn   (4b)
  0.25160000000000   0.75160000000000   0.50000000000000   Sn   (4b)
```

$E3_1$ (β-Ag$_2$HgI$_4$) (*obsolete*): A2BC4_tP7_111_f_a_n - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Ag2HgI4'
_chemical_formula_sum 'Ag2 Hg I4'

loop_
_publ_author_name
 'J. A. A. Ketelaar'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 80
_journal_year 1931
_journal_page_first 190
_journal_page_last 203
_publ_Section_title
;
 Strukturbestimmung der komplexen Quecksilberverbindungen Ag$_{2}$HgJ$_{
    ↪ 4}$ und Cu$_{2}$HgJ$_{4}$
;

# Found in Single-crystal studies of $\beta$-Ag$_{2}$HgI$_{4}$, 1974

_aflow_title '$E3_{1}$ ($\beta$-Ag$_{2}$HgI$_{4}$) ({\em{obsolete}})
    ↪ Structure'
_aflow_proto 'A2BC4_tP7_111_f_a_n'
_aflow_params 'a,c/a,x_{3},z_{3}'
_aflow_params_values '6.34,1.0,0.27,0.225'
_aflow_Strukturbericht '$E3_{1}$'
_aflow_Pearson 'tP7'

_symmetry_space_group_name_H-M "P -4 2 m"
_symmetry_Int_Tables_number 111

_cell_length_a     6.34000
_cell_length_b     6.34000
_cell_length_c     6.34000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 y,x,z
6 y,-x,-z
7 -y,x,-z
8 -y,-x,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Hg1 Hg 1 a 0.00000 0.00000 0.00000 1.00000
Ag1 Ag 2 f 0.50000 0.00000 0.50000 1.00000
I1 I 4 n 0.27000 0.27000 0.22500 1.00000
```

$E3_1$ (β-Ag$_2$HgI$_4$) (*obsolete*): A2BC4_tP7_111_f_a_n - POSCAR

```
A2BC4_tP7_111_f_a_n & a,c/a,x3,z3 --params=6.34,1.0,0.27,0.225 & P-42m
    ↪ D_{2d}^{1} #111 (afn) & tP7 & $E3_{1}$ & Ag2HgI4 & Ag2HgI4 & J.
    ↪ A. A. Ketelaar, Zeitschrift f{"u}r Kristallographie -
    ↪ Crystalline Materials 80, 190-203 (1931)
   1.00000000000000
   6.34000000000000   0.00000000000000   0.00000000000000
   0.00000000000000   6.34000000000000   0.00000000000000
   0.00000000000000   0.00000000000000   6.34000000000000
    Ag    Hg     I
     2     1     4
Direct
   0.50000000000000   0.00000000000000   0.50000000000000   Ag   (2f)
   0.00000000000000   0.50000000000000   0.50000000000000   Ag   (2f)
   0.00000000000000   0.00000000000000   0.00000000000000   Hg   (1a)
   0.27000000000000   0.27000000000000   0.22500000000000    I   (4n)
  -0.27000000000000  -0.27000000000000   0.22500000000000    I   (4n)
   0.27000000000000  -0.27000000000000  -0.22500000000000    I   (4n)
  -0.27000000000000   0.27000000000000  -0.22500000000000    I   (4n)
```

Ammonium Chlorite (NH_4ClO_2): AB4CD2_tP16_113_c_f_a_e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM
```

```
_chemical_name_mineral 'ClH4NO2'
_chemical_formula_sum 'Cl H4 N O2'

loop_
_publ_author_name
 'A. I. Smolentsev'
 'D. Y. Naumov'
_journal_name_full_name
;
 Acta Crystallographica Section E: Crystallographic Communications
;
_journal_volume 61
_journal_year 2005
_journal_page_first i38
_journal_page_last i40
_publ_Section_title
;
 Ammonium chlorite, NH$_{4}$ClO$_{2}$, at 150~K
;

_aflow_title 'Ammonium Chlorite (NH$_{4}$ClO$_{2}$) Structure'
_aflow_proto 'AB4CD2_tP16_113_c_f_a_e'
_aflow_params 'a,c/a,z_{2},x_{3},z_{3},x_{4},y_{4},z_{4}'
_aflow_params_values '6.3397,0.592457056328,0.37123,0.14562,0.6062,0.111
      ↪ ,0.03,0.125'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tP16'

_symmetry_space_group_name_H-M "P -4 21 m"
_symmetry_Int_Tables_number 113

_cell_length_a    6.33970
_cell_length_b    6.33970
_cell_length_c    3.75600
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z
3 -x+1/2,y+1/2,-z
4 -x,-y,z
5 y+1/2,x+1/2,z
6 y,-x,-z
7 -y,x,-z
8 -y+1/2,-x+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
N1 N 2 a 0.00000 0.00000 0.00000 1.00000
Cl1 Cl 2 c 0.00000 0.50000 0.37123 1.00000
O1 O 4 e 0.14562 0.64562 0.60620 1.00000
H1 H 8 f 0.11100 0.03000 0.12500 1.00000
```

Ammonium Chlorite (NH_4ClO_2): AB4CD2_tP16_113_c_f_a_e - POSCAR

```
AB4CD2_tP16_113_c_f_a_e & a,c/a,z2,x3,z3,x4,y4,z4 --params=6.3397,
      ↪ 0.592457056328,0.37123,0.14562,0.6062,0.111,0.03,0.125 & P-42_{
      ↪ 1}m D_{2d}^{3} #113 (acef) & tP16 & None & ClH4NO2 & ClH4NO2 &
      ↪ A. I. Smolentsev and D. Y. Naumov, Acta Crystallogr. E 61,
      ↪ i38-i40 (2005)
   1.00000000000000
   6.33970000000000    0.00000000000000    0.00000000000000
   0.00000000000000    6.33970000000000    0.00000000000000
   0.00000000000000    0.00000000000000    3.75600000000000
     Cl     H     N     O
      2     8     2     4
Direct
   0.00000000000000    0.50000000000000    0.37123000000000   Cl   (2c)
   0.50000000000000    0.00000000000000   -0.37123000000000   Cl   (2c)
   0.11100000000000    0.03000000000000    0.12500000000000   H    (8f)
  -0.11100000000000   -0.03000000000000    0.12500000000000   H    (8f)
   0.03000000000000   -0.11100000000000   -0.12500000000000   H    (8f)
  -0.03000000000000    0.11100000000000   -0.12500000000000   H    (8f)
   0.38900000000000    0.53000000000000   -0.12500000000000   H    (8f)
   0.61100000000000    0.47000000000000   -0.12500000000000   H    (8f)
   0.47000000000000    0.38900000000000    0.12500000000000   H    (8f)
   0.53000000000000    0.61100000000000    0.12500000000000   H    (8f)
   0.00000000000000    0.00000000000000    0.00000000000000   N    (2a)
   0.50000000000000    0.50000000000000    0.00000000000000   N    (2a)
   0.14562000000000    0.64562000000000    0.60620000000000   O    (4e)
  -0.14562000000000    0.35438000000000    0.60620000000000   O    (4e)
   0.64562000000000   -0.14562000000000   -0.60620000000000   O    (4e)
   0.35438000000000    0.14562000000000   -0.60620000000000   O    (4e)
```

$C_{19}Sc_{15}$: A19B15_tP68_114_bc4e_ac3e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'C19Sc15'
_chemical_formula_sum 'C19 Sc15'

loop_
_publ_author_name
 'H. Jedlicka'
 'H. Nowotny'
 'F. Benesovsky'
_journal_name_full_name
;
 Monatshefte f{\"u}r Chemie - Chemical Monthly
;
_journal_volume 102
_journal_year 1971
_journal_page_first 389
_journal_page_last 403
_publ_Section_title
;
 Zum System Scandium-Kohlenstoff, 2. Mitt.: Kristallstruktur des
      ↪ C-reichen Carbids
;

_aflow_title 'C$_{19}$Sc$_{15}$ Structure'
_aflow_proto 'A19B15_tP68_114_bc4e_ac3e'
_aflow_params 'a,c/a,z_{3},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{6},x_{
      ↪ 7},y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{10},y_{10
      ↪ },z_{10},x_{11},y_{11},z_{11}'
_aflow_params_values '7.5,2.0,0.165,0.3198,0.712,0.118,-0.008,0.384,0.2,
      ↪ 0.1333,0.436,0.208,0.212,-0.194,0.375,0.179,0.4021,0.1993,-0.02
      ↪ ,0.0982,0.2823,0.1593,0.7157,0.0982,0.1423'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tP68'

_symmetry_space_group_name_H-M "P -4 21 c"
_symmetry_Int_Tables_number 114

_cell_length_a    7.50000
_cell_length_b    7.50000
_cell_length_c    15.00000
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x+1/2,y+1/2,-z+1/2
4 -x,-y,z
5 y+1/2,x+1/2,z+1/2
6 y,-x,-z
7 -y,x,-z
8 -y+1/2,-x+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Sc1 Sc 2 a 0.00000 0.00000 0.00000 1.00000
C1 C 2 b 0.00000 0.00000 0.50000 1.00000
C2 C 4 c 0.00000 0.00000 0.16500 1.00000
Sc2 Sc 4 c 0.00000 0.00000 0.31980 1.00000
C3 C 8 e 0.71200 0.11800 -0.00800 1.00000
C4 C 8 e 0.38400 0.20000 0.13330 1.00000
C5 C 8 e 0.43600 0.20800 0.21200 1.00000
C6 C 8 e -0.19400 0.37500 0.17900 1.00000
Sc3 Sc 8 e 0.40210 0.19930 -0.02000 1.00000
Sc4 Sc 8 e 0.09820 0.28230 0.15930 1.00000
Sc5 Sc 8 e 0.71570 0.09820 0.14230 1.00000
```

$C_{19}Sc_{15}$: A19B15_tP68_114_bc4e_ac3e - POSCAR

```
A19B15_tP68_114_bc4e_ac3e & a,c/a,z3,z4,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8,y8
      ↪ ,z8,x9,y9,z9,x10,y10,z10,x11,y11,z11 --params=7.5,2.0,0.165,
      ↪ 0.3198,0.712,0.118,-0.008,0.384,0.2,0.1333,0.436,0.208,0.212,-
      ↪ 0.194,0.375,0.179,0.4021,0.1993,-0.02,0.0982,0.2823,0.1593,
      ↪ 0.7157,0.0982,0.1423 & P-42_{1}c D_{2d}^{4} #114 (abc^2e^7) &
      ↪ tP68 & None & C19Sc15 & C19Sc15 & H. Jedlicka and H. Nowotny
      ↪ and F. Benesovsky, Monatshefte f{"u}r Chemie - Chemical Monthly
      ↪ 102, 389-403 (1971)
   1.00000000000000
   7.50000000000000    0.00000000000000    0.00000000000000
   0.00000000000000    7.50000000000000    0.00000000000000
   0.00000000000000    0.00000000000000   15.00000000000000
     C     Sc
    38     30
Direct
   0.00000000000000    0.00000000000000    0.50000000000000   C    (2b)
   0.50000000000000    0.50000000000000    0.00000000000000   C    (2b)
   0.00000000000000    0.00000000000000    0.16500000000000   C    (4c)
   0.00000000000000    0.00000000000000   -0.16500000000000   C    (4c)
   0.50000000000000    0.50000000000000    0.33500000000000   C    (4c)
   0.50000000000000    0.50000000000000    0.66500000000000   C    (4c)
   0.71200000000000    0.11800000000000   -0.00800000000000   C    (8e)
  -0.71200000000000   -0.11800000000000   -0.00800000000000   C    (8e)
   0.11800000000000   -0.71200000000000    0.00800000000000   C    (8e)
  -0.11800000000000    0.71200000000000    0.00800000000000   C    (8e)
  -0.21200000000000    0.61800000000000    0.50800000000000   C    (8e)
   1.21200000000000    0.38200000000000    0.50800000000000   C    (8e)
   0.38200000000000   -0.21200000000000    0.49200000000000   C    (8e)
   0.61800000000000    1.21200000000000    0.49200000000000   C    (8e)
   0.38400000000000    0.20000000000000    0.13330000000000   C    (8e)
  -0.38400000000000   -0.20000000000000    0.13330000000000   C    (8e)
```

```
   0.20000000000000  −0.38400000000000  −0.13330000000000   C   (8e)
  −0.20000000000000   0.38400000000000  −0.13330000000000   C   (8e)
   0.11600000000000   0.70000000000000   0.36670000000000   C   (8e)
   0.88400000000000   0.30000000000000   0.36670000000000   C   (8e)
   0.30000000000000   0.11600000000000   0.63330000000000   C   (8e)
   0.70000000000000   0.88400000000000   0.63330000000000   C   (8e)
   0.43600000000000   0.20800000000000   0.21200000000000   C   (8e)
  −0.43600000000000  −0.20800000000000   0.21200000000000   C   (8e)
   0.20800000000000  −0.43600000000000  −0.21200000000000   C   (8e)
  −0.20800000000000   0.43600000000000  −0.21200000000000   C   (8e)
   0.06400000000000   0.70800000000000   0.28800000000000   C   (8e)
   0.93600000000000   0.29200000000000   0.28800000000000   C   (8e)
   0.29200000000000   0.06400000000000   0.71200000000000   C   (8e)
   0.70800000000000   0.93600000000000   0.71200000000000   C   (8e)
  −0.19400000000000   0.37500000000000   0.17900000000000   C   (8e)
   0.19400000000000  −0.37500000000000   0.17900000000000   C   (8e)
   0.37500000000000   0.19400000000000  −0.17900000000000   C   (8e)
  −0.37500000000000  −0.19400000000000  −0.17900000000000   C   (8e)
   0.69400000000000   0.87500000000000   0.32100000000000   C   (8e)
   0.30600000000000   0.12500000000000   0.32100000000000   C   (8e)
   0.12500000000000   0.69400000000000   0.67900000000000   C   (8e)
   0.87500000000000   0.30600000000000   0.67900000000000   C   (8e)
   0.00000000000000   0.00000000000000   0.00000000000000   Sc  (2a)
   0.50000000000000   0.50000000000000   0.50000000000000   Sc  (2a)
   0.00000000000000   0.00000000000000   0.31980000000000   Sc  (4c)
   0.00000000000000   0.00000000000000  −0.31980000000000   Sc  (4c)
   0.50000000000000   0.50000000000000   0.18020000000000   Sc  (4c)
   0.50000000000000   0.50000000000000   0.81980000000000   Sc  (4c)
   0.40210000000000   0.19930000000000  −0.02000000000000   Sc  (8e)
  −0.40210000000000  −0.19930000000000  −0.02000000000000   Sc  (8e)
   0.19930000000000  −0.40210000000000   0.02000000000000   Sc  (8e)
  −0.19930000000000   0.40210000000000   0.02000000000000   Sc  (8e)
   0.09790000000000   0.69930000000000   0.52000000000000   Sc  (8e)
   0.90210000000000   0.30070000000000   0.52000000000000   Sc  (8e)
   0.30070000000000   0.09790000000000   0.48000000000000   Sc  (8e)
   0.69930000000000   0.90210000000000   0.48000000000000   Sc  (8e)
   0.09820000000000   0.28230000000000   0.15930000000000   Sc  (8e)
  −0.09820000000000  −0.28230000000000   0.15930000000000   Sc  (8e)
   0.28230000000000  −0.09820000000000  −0.15930000000000   Sc  (8e)
  −0.28230000000000   0.09820000000000  −0.15930000000000   Sc  (8e)
   0.40180000000000   0.78230000000000   0.34070000000000   Sc  (8e)
   0.59820000000000   0.21770000000000   0.34070000000000   Sc  (8e)
   0.21770000000000   0.40180000000000   0.65930000000000   Sc  (8e)
   0.78230000000000   0.59820000000000   0.65930000000000   Sc  (8e)
   0.71570000000000   0.09820000000000   0.14230000000000   Sc  (8e)
  −0.71570000000000  −0.09820000000000   0.14230000000000   Sc  (8e)
   0.09820000000000  −0.71570000000000  −0.14230000000000   Sc  (8e)
  −0.09820000000000   0.71570000000000  −0.14230000000000   Sc  (8e)
  −0.21570000000000   0.59820000000000   0.35770000000000   Sc  (8e)
   1.21570000000000   0.40180000000000   0.35770000000000   Sc  (8e)
   0.40180000000000  −0.21570000000000   0.64230000000000   Sc  (8e)
   0.59820000000000   1.21570000000000   0.64230000000000   Sc  (8e)
```

$Ag_2SO_4{\cdot}4NH_3$ ($H4_{17}$): A2B12C4D4E_tP46_114_d_3e_e_e_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Ag2H12N4O4S'
_chemical_formula_sum 'Ag2 H12 N4 O4 S'

loop_
_publ_author_name
 'U. Zachwieja'
 'H. Jacobs'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 201
_journal_year 1992
_journal_page_first 207
_journal_page_last 212
_publ_Section_title
;
 Redetermination of the crystal structure of diammine silver(I)-sulfate,
      ↪ [Ag(NH$_{3}$)$_{2}$]$_{2}$SO$_{4}$
;

_aflow_title 'Ag$_{2}$SO$_{4}$$\cdot$4NH$_{3}$ ($H4_{17}$) Structure'
_aflow_proto 'A2B12C4D4E_tP46_114_d_3e_e_e_a'
_aflow_params 'a,c/a,z_{2},x_{3},y_{3},z_{3},x_{4},y_{4},z_{4},x_{5},y_{
      ↪ 5},z_{5},x_{6},y_{6},z_{6},x_{7},y_{7},z_{7}'
_aflow_params_values '8.442,0.757995735608,0.49194,0.082,0.21,0.416,
      ↪ 0.183,0.287,0.52,0.063,0.227,0.603,0.1081,0.275,0.5083,0.1301,
      ↪ 0.0595,0.1309'
_aflow_Strukturbericht '$H4_{17}$'
_aflow_Pearson 'tP46'

_symmetry_space_group_name_H-M "P -4 21 c"
_symmetry_Int_Tables_number 114

_cell_length_a    8.44200
_cell_length_b    8.44200
_cell_length_c    6.39900
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x+1/2,y+1/2,-z+1/2
4 -x,-y,z
5 y+1/2,x+1/2,z+1/2
6 y,-x,-z
7 -y,x,-z
8 -y+1/2,-x+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
S1 S 2 a 0.00000 0.00000 0.00000 1.00000
Ag1 Ag 4 d 0.00000 0.50000 0.49194 1.00000
H1 H 8 e 0.08200 0.21000 0.41600 1.00000
H2 H 8 e 0.18300 0.28700 0.52000 1.00000
H3 H 8 e 0.06300 0.22700 0.60300 1.00000
N1 N 8 e 0.10810 0.27500 0.50830 1.00000
O1 O 8 e 0.13010 0.05950 0.13090 1.00000
```

$Ag_2SO_4{\cdot}4NH_3$ ($H4_{17}$): A2B12C4D4E_tP46_114_d_3e_e_e_a - POSCAR

```
A2B12C4D4E_tP46_114_d_3e_e_e_a & a,c/a,z2,x3,y3,z3,x4,y4,z4,x5,y5,z5,x6,
      ↪ y6,z6,x7,y7,z7 --params=8.442,0.757995735608,0.49194,0.082,0.21
      ↪ ,0.416,0.183,0.287,0.52,0.063,0.227,0.603,0.1081,0.275,0.5083,
      ↪ 0.1301,0.0595,0.1309 & P-42_{1}c D_{2d}^{4} #114 (ade^5) & tP46
      ↪  & $H4_{17}$ & Ag2H12N4O4S & Ag2H12N4O4S & U. Zachwieja and H.
      ↪ Jacobs, Zeitschrift f{"u}r Kristallographie - Crystalline
      ↪ Materials 201, 207-212 (1992)
   1.00000000000000
   8.44200000000000   0.00000000000000   0.00000000000000
   0.00000000000000   8.44200000000000   0.00000000000000
   0.00000000000000   0.00000000000000   6.39900000000000
    Ag    H    N    O    S
     4   24    8    8    2
Direct
   0.00000000000000   0.50000000000000   0.49194000000000   Ag   (4d)
   0.50000000000000   0.00000000000000  −0.49194000000000   Ag   (4d)
   0.50000000000000   0.00000000000000   0.00806000000000   Ag   (4d)
   0.00000000000000   0.50000000000000   0.99194000000000   Ag   (4d)
   0.08200000000000   0.21000000000000   0.41600000000000   H    (8e)
  −0.08200000000000  −0.21000000000000   0.41600000000000   H    (8e)
   0.21000000000000  −0.08200000000000  −0.41600000000000   H    (8e)
  −0.21000000000000   0.08200000000000  −0.41600000000000   H    (8e)
   0.41800000000000   0.71000000000000   0.08400000000000   H    (8e)
   0.58200000000000   0.29000000000000   0.08400000000000   H    (8e)
   0.29000000000000   0.41800000000000   0.91600000000000   H    (8e)
   0.71000000000000   0.58200000000000   0.91600000000000   H    (8e)
   0.18300000000000   0.28700000000000   0.52000000000000   H    (8e)
  −0.18300000000000  −0.28700000000000   0.52000000000000   H    (8e)
   0.28700000000000  −0.18300000000000  −0.52000000000000   H    (8e)
  −0.28700000000000   0.18300000000000  −0.52000000000000   H    (8e)
   0.31700000000000   0.78700000000000  −0.02000000000000   H    (8e)
   0.68300000000000   0.21300000000000  −0.02000000000000   H    (8e)
   0.21300000000000   0.31700000000000   1.02000000000000   H    (8e)
   0.78700000000000   0.68300000000000   1.02000000000000   H    (8e)
   0.06300000000000   0.22700000000000   0.60300000000000   H    (8e)
  −0.06300000000000  −0.22700000000000   0.60300000000000   H    (8e)
   0.22700000000000  −0.06300000000000  −0.60300000000000   H    (8e)
  −0.22700000000000   0.06300000000000  −0.60300000000000   H    (8e)
   0.43700000000000   0.72700000000000  −0.10300000000000   H    (8e)
   0.56300000000000   0.27300000000000  −0.10300000000000   H    (8e)
   0.27300000000000   0.43700000000000   1.10300000000000   H    (8e)
   0.72700000000000   0.56300000000000   1.10300000000000   H    (8e)
   0.10810000000000   0.27500000000000   0.50830000000000   N    (8e)
  −0.10810000000000  −0.27500000000000   0.50830000000000   N    (8e)
   0.27500000000000  −0.10810000000000  −0.50830000000000   N    (8e)
  −0.27500000000000   0.10810000000000  −0.50830000000000   N    (8e)
   0.39190000000000   0.77500000000000  −0.00830000000000   N    (8e)
   0.60810000000000   0.22500000000000  −0.00830000000000   N    (8e)
   0.22500000000000   0.39190000000000   1.00830000000000   N    (8e)
   0.77500000000000   0.60810000000000   1.00830000000000   N    (8e)
   0.13010000000000   0.05950000000000   0.13090000000000   O    (8e)
  −0.13010000000000  −0.05950000000000   0.13090000000000   O    (8e)
   0.05950000000000  −0.13010000000000  −0.13090000000000   O    (8e)
  −0.05950000000000   0.13010000000000  −0.13090000000000   O    (8e)
   0.36990000000000   0.55950000000000   0.36910000000000   O    (8e)
   0.63010000000000   0.44050000000000   0.36910000000000   O    (8e)
   0.44050000000000   0.36990000000000   0.63090000000000   O    (8e)
   0.55950000000000   0.63010000000000   0.63090000000000   O    (8e)
   0.00000000000000   0.00000000000000   0.00000000000000   S    (2a)
   0.50000000000000   0.50000000000000   0.50000000000000   S    (2a)
```

$F6_1$ (Chalcopyrite, $CuFeS_2$) (*obsolete*): ABC2_tP4_115_a_c_g - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'CuFeS2'
_chemical_formula_sum 'Cu Fe S2'

loop_
_publ_author_name
 'R. Gro{\ss}'
 'N. Gro{\ss}'
_journal_name_full_name
;
 Neues Jahrbuch fur Mineralogie, Monatshefte
;
_journal_volume 48
```

```
_journal_year 1923
_journal_page_first 113
_journal_page_last 135
_publ_Section_title
;
 ~
;

# Found in Strukturbericht 1913-1928, 1931

_aflow_title '$F6_{1}$ (Chalcopyrite, CuFeS$_{2}$) ({\em{obsolete}})
      ↪ Structure'
_aflow_proto 'ABC2_tP4_115_a_c_g'
_aflow_params 'a,c/a,z_{3}'
_aflow_params_values '3.72645,1.39381985536,0.19'
_aflow_Strukturbericht '$F6_{1}$'
_aflow_Pearson 'tP4'

_symmetry_space_group_name_H-M "P -4 m 2"
_symmetry_Int_Tables_number 115

_cell_length_a     3.72645
_cell_length_b     3.72645
_cell_length_c     5.19400
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,z
3 -y,-x,-z
4 y,x,-z
5 -x,y,z
6 x,-y,z
7 y,-x,-z
8 -y,x,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cu1 Cu 1 a 0.00000 0.00000 0.00000 1.00000
Fe1 Fe 1 c 0.50000 0.50000 0.50000 1.00000
S1 S 2 g 0.00000 0.50000 0.19000 1.00000
```

$F6_1$ (Chalcopyrite, $CuFeS_2$) (*obsolete*): ABC2_tP4_115_a_c_g - POSCAR

```
ABC2_tP4_115_a_c_g & a,c/a,z3 --params=3.72645,1.39381985536,0.19 &
      ↪ P-4m2 D_{2d}^{5} #115 (acg) & tP4 & $F6_{1}$ & CuFeS2 & CuFeS2
      ↪ & R. Gro{\ss} and N. Gro{\ss}, Neues Jahrb. Mineral. Monatsh.
      ↪ 48, 113-135 (1923)
   1.00000000000000
   3.72645000000000    0.00000000000000    0.00000000000000
   0.00000000000000    3.72645000000000    0.00000000000000
   0.00000000000000    0.00000000000000    5.19400000000000
    Cu     Fe     S
     1      1     2
Direct
   0.00000000000000    0.00000000000000    0.00000000000000   Cu   (1a)
   0.50000000000000    0.50000000000000    0.50000000000000   Fe   (1c)
   0.00000000000000    0.50000000000000    0.19000000000000    S   (2g)
   0.50000000000000    0.00000000000000   -0.19000000000000    S   (2g)
```

Phase II $Cd_2Re_2O_7$: A2B7C2_tI44_119_i_bdefgh_i - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Cd2O7Re2'
_chemical_formula_sum 'Cd2 O7 Re2'

loop_
_publ_author_name
 'S.-W. Huang'
 'H.-T. Jeng'
 'J.-Y. Lin'
 'W. J. Chang'
 'J. M. Chen'
 'G. H. Lee'
 'H. Berger'
 'H. D. Yang'
 'K. S. Liang'
_journal_name_full_name
;
 Journal of Physics: Condensed Matter
;
_journal_volume 21
_journal_year 2009
_journal_page_first 195602
_journal_page_last 195602
_publ_Section_title
;
 Electronic structure of pyrochlore Cd$_{2}$Re$_{2}$O$_{7}$
;
```

```
# Found in The crystal structure of the inversion breaking metal Cd$_{2}
      ↪ $Re$_{2}$O$_{7}$, {arXiv:1911.10141 [cond-mat.str-el]}, 2019

_aflow_title 'Phase II Cd$_{2}$Re$_{2}$O$_{7}$ Structure'
_aflow_proto 'A2B7C2_tI44_119_i_bdefgh_i'
_aflow_params 'a,c/a,z_{3},z_{4},x_{5},x_{6},x_{7},z_{7},x_{8},z_{8}'
_aflow_params_values '7.2312,1.41410830844,0.1974,0.0683,0.1889,0.1941,
      ↪ 0.2471,0.3741,0.2471,0.87294'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tI44'

_symmetry_space_group_name_H-M "I -4 m 2"
_symmetry_Int_Tables_number 119

_cell_length_a     7.23120
_cell_length_b     7.23120
_cell_length_c     10.22570
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,z
3 -y,-x,-z
4 y,x,-z
5 -x,y,z
6 x,-y,z
7 y,-x,-z
8 -y,x,-z
9 x+1/2,y+1/2,z+1/2
10 -x+1/2,-y+1/2,z+1/2
11 -y+1/2,-x+1/2,-z+1/2
12 y+1/2,x+1/2,-z+1/2
13 -x+1/2,y+1/2,z+1/2
14 x+1/2,-y+1/2,z+1/2
15 y+1/2,-x+1/2,-z+1/2
16 -y+1/2,x+1/2,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 2 b 0.00000 0.00000 0.50000 1.00000
O2 O 2 d 0.00000 0.50000 0.75000 1.00000
O3 O 4 e 0.00000 0.00000 0.19740 1.00000
O4 O 4 f 0.00000 0.50000 0.06830 1.00000
O5 O 8 g 0.18890 0.18890 0.00000 1.00000
O6 O 8 h 0.19410 0.69410 0.25000 1.00000
Cd1 Cd 8 i 0.24710 0.00000 0.37410 1.00000
Re1 Re 8 i 0.24710 0.00000 0.87294 1.00000
```

Phase II $Cd_2Re_2O_7$: A2B7C2_tI44_119_i_bdefgh_i - POSCAR

```
A2B7C2_tI44_119_i_bdefgh_i & a,c/a,z3,z4,x5,x6,x7,z7,x8,z8 --params=
      ↪ 7.2312,1.41410830844,0.1974,0.0683,0.1889,0.1941,0.2471,0.3741,
      ↪ 0.2471,0.87294 & I-4m2 D_{2d}^{9} #119 (bdefghi^2) & tI44 &
      ↪ None & Cd2O7Re2 & Cd2O7Re2 & S.-W. Huang et al., J. Phys.:
      ↪ Condens. Matter 21, 195602(2009)
   1.00000000000000
  -3.61560000000000    3.61560000000000    5.11285000000000
   3.61560000000000   -3.61560000000000    5.11285000000000
   3.61560000000000    3.61560000000000   -5.11285000000000
    Cd     O     Re
     4    14      4
Direct
   0.37410000000000    0.62120000000000    0.24710000000000   Cd   (8i)
   0.37410000000000    0.12700000000000   -0.24710000000000   Cd   (8i)
  -0.62120000000000   -0.37410000000000   -0.24710000000000   Cd   (8i)
  -0.12700000000000   -0.37410000000000    0.24710000000000   Cd   (8i)
   0.50000000000000    0.50000000000000    0.00000000000000    O   (2b)
   0.25000000000000    0.75000000000000    0.50000000000000    O   (2d)
   0.19740000000000    0.19740000000000    0.00000000000000    O   (4e)
  -0.19740000000000   -0.19740000000000    0.00000000000000    O   (4e)
   0.56830000000000    0.06830000000000    0.50000000000000    O   (4f)
  -0.06830000000000    0.43170000000000    0.50000000000000    O   (4f)
   0.18890000000000    0.18890000000000    0.37780000000000    O   (8g)
  -0.18890000000000   -0.18890000000000   -0.37780000000000    O   (8g)
  -0.18890000000000    0.18890000000000    0.00000000000000    O   (8g)
   0.18890000000000   -0.18890000000000    0.00000000000000    O   (8g)
   0.94410000000000    0.44410000000000    0.88820000000000    O   (8h)
   0.55590000000000    0.05590000000000    0.11180000000000    O   (8h)
   0.55590000000000    0.44410000000000    0.50000000000000    O   (8h)
   0.94410000000000    0.05590000000000    0.50000000000000    O   (8h)
   0.87294000000000    1.12004000000000    0.24710000000000   Re   (8i)
   0.87294000000000    0.62584000000000   -0.24710000000000   Re   (8i)
  -1.12004000000000   -0.87294000000000   -0.24710000000000   Re   (8i)
  -0.62584000000000   -0.87294000000000    0.24710000000000   Re   (8i)
```

Tetragonal $TlFeS_2$: AB2C_tI8_119_c_e_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'FeS2Tl'
_chemical_formula_sum 'Fe S2 Tl'
```

```
loop_
_publ_author_name
 'A. Kutoglu'
_journal_name_full_name
;
 Naturwissenschaften
;
_journal_volume 61
_journal_year 1974
_journal_page_first 125
_journal_page_last 126
_publ_Section_title
;
 Synthese und Kristallstrukturen von TlFeS$_{2}$ und TlFeSe$_{2}$
;

# Found in Neutron-Diffraction Study in TlFeS$_{2}$ and TlFeSe$_{2}$ at
      ↪ Low Temperatures, 2014

_aflow_title 'Tetragonal TlFeS$_{2}$ Structure'
_aflow_proto 'AB2C_tI8_119_c_e_a'
_aflow_params 'a,c/a,z_{3}'
_aflow_params_values '3.753,3.55502264855,0.35126'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tI8'

_symmetry_space_group_name_H-M "I -4 m 2"
_symmetry_Int_Tables_number 119

_cell_length_a     3.75300
_cell_length_b     3.75300
_cell_length_c     13.34200
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,z
3 -y,-x,-z
4 y,x,-z
5 -x,y,z
6 x,-y,z
7 y,-x,-z
8 -y,x,-z
9 x+1/2,y+1/2,z+1/2
10 -x+1/2,-y+1/2,z+1/2
11 -y+1/2,-x+1/2,-z+1/2
12 y+1/2,x+1/2,-z+1/2
13 -x+1/2,y+1/2,z+1/2
14 x+1/2,-y+1/2,z+1/2
15 y+1/2,-x+1/2,-z+1/2
16 -y+1/2,x+1/2,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Tl1 Tl 2 a 0.00000 0.00000 0.00000 1.00000
Fe1 Fe 2 c 0.00000 0.50000 0.25000 1.00000
S1 S 4 e 0.00000 0.00000 0.35126 1.00000
```

Tetragonal $TlFeS_2$: AB2C_tI8_119_c_e_a - POSCAR

```
AB2C_tI8_119_c_e_a & a,c/a,z3 --params=3.753,3.55502264855,0.35126 &
      ↪ I-4m2 D_{2d}^{9} #119 (ace) & tI8 & None & FeS2Tl & FeS2Tl & A.
      ↪  Kutoglu, Naturwissenschaften 61, 125-126 (1974)
   1.00000000000000
  -1.87650000000000    1.87650000000000    6.67100000000000
   1.87650000000000   -1.87650000000000    6.67100000000000
   1.87650000000000    1.87650000000000   -6.67100000000000
    Fe     S     Tl
     1     2     1
Direct
   0.75000000000000    0.25000000000000    0.50000000000000    Fe    (2c)
   0.35126000000000    0.35126000000000    0.00000000000000    S     (4e)
  -0.35126000000000   -0.35126000000000    0.00000000000000    S     (4e)
   0.00000000000000    0.00000000000000    0.00000000000000    Tl    (2a)
```

$BeSO_4{\cdot}4H_2O$ ($H4_3$): AB8C8D_tI72_120_c_2i_2i_b - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'BeH8O8S'
_chemical_formula_sum 'Be H8 O8 S'

loop_
_publ_author_name
 'S. K. Sikka'
 'R. Chidambaram'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 25
_journal_year 1969
_journal_page_first 310
_journal_page_last 315
_publ_Section_title
;
 A neutron diffraction determination of the structure of beryllium
      ↪ sulphate tetrahydrate, BeSO$_{4}$$\cdot$4H$_{2}$O
;

_aflow_title 'BeSO$_{4}$$\cdot$4H$_{2}$O ($H4_{3}$) Structure'
_aflow_proto 'AB8C8D_tI72_120_c_2i_2i_b'
_aflow_params 'a,c/a,x_{3},y_{3},z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{
      ↪ 5},x_{6},y_{6},z_{6}'
_aflow_params_values '7.99,1.33767209011,0.22729,0.45696,0.13106,0.1229,
      ↪ 0.28418,0.13628,0.12563,0.08408,0.07733,0.13758,0.3956,0.17093'
_aflow_Strukturbericht '$H4_{3}$'
_aflow_Pearson 'tI72'

_symmetry_space_group_name_H-M "I -4 c 2"
_symmetry_Int_Tables_number 120

_cell_length_a     7.99000
_cell_length_b     7.99000
_cell_length_c     10.68800
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -x,-y,z
3 -y,-x,-z+1/2
4 y,x,-z+1/2
5 -x,y,z+1/2
6 x,-y,z+1/2
7 y,-x,-z
8 -y,x,-z
9 x+1/2,y+1/2,z+1/2
10 -x+1/2,-y+1/2,z+1/2
11 -y+1/2,-x+1/2,-z
12 y+1/2,x+1/2,-z
13 -x+1/2,y+1/2,z
14 x+1/2,-y+1/2,z
15 y+1/2,-x+1/2,-z+1/2
16 -y+1/2,x+1/2,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
S1 S 4 b 0.00000 0.00000 0.00000 1.00000
Be1 Be 4 c 0.00000 0.50000 0.25000 1.00000
H1 H 16 i 0.22729 0.45696 0.13106 1.00000
H2 H 16 i 0.12290 0.28418 0.13628 1.00000
O1 O 16 i 0.12563 0.08408 0.07733 1.00000
O2 O 16 i 0.13758 0.39560 0.17093 1.00000
```

$BeSO_4{\cdot}4H_2O$ ($H4_3$): AB8C8D_tI72_120_c_2i_2i_b - POSCAR

```
AB8C8D_tI72_120_c_2i_2i_b & a,c/a,x3,y3,z3,x4,y4,z4,x5,y5,z5,x6,y6,z6 --
      ↪ params=7.99,1.33767209011,0.22729,0.45696,0.13106,0.1229,
      ↪ 0.28418,0.13628,0.12563,0.08408,0.07733,0.13758,0.3956,0.17093
      ↪ & I-4c2 D_{2d}^{10} #120 (bci^4) & tI72 & $H4_{3}$ & BeH8O8S &
      ↪ BeH8O8S & S. K. Sikka and R. Chidambaram, Acta Crystallogr.
      ↪ Sect. B Struct. Sci. 25, 310-315 (1969)
   1.00000000000000
  -3.99500000000000    3.99500000000000    5.34400000000000
   3.99500000000000   -3.99500000000000    5.34400000000000
   3.99500000000000    3.99500000000000   -5.34400000000000
    Be     H     O     S
     2    16    16     2
Direct
   0.75000000000000    0.25000000000000    0.50000000000000    Be    (4c)
   0.25000000000000    0.75000000000000    0.50000000000000    Be    (4c)
   0.58802000000000    0.35835000000000    0.68425000000000    H    (16i)
  -0.32590000000000   -0.09623000000000   -0.68425000000000    H    (16i)
  -0.35835000000000    0.32590000000000    0.22967000000000    H    (16i)
   0.09623000000000   -0.58802000000000   -0.22967000000000    H    (16i)
   0.17410000000000    0.85835000000000   -0.22967000000000    H    (16i)
   1.08802000000000    0.40377000000000    0.22967000000000    H    (16i)
   0.59623000000000    0.82590000000000    0.68425000000000    H    (16i)
   0.14165000000000   -0.08802000000000   -0.68425000000000    H    (16i)
   0.42046000000000    0.25918000000000    0.40708000000000    H    (16i)
  -0.14790000000000    0.01338000000000   -0.40708000000000    H    (16i)
  -0.25918000000000    0.14790000000000    0.16128000000000    H    (16i)
  -0.01338000000000   -0.42046000000000   -0.16128000000000    H    (16i)
   0.35210000000000    0.75918000000000   -0.16128000000000    H    (16i)
   0.92046000000000    0.51338000000000    0.16128000000000    H    (16i)
   0.48662000000000    0.64790000000000    0.40708000000000    H    (16i)
   0.24082000000000    0.07954000000000   -0.40708000000000    H    (16i)
   0.16141000000000    0.20296000000000    0.20971000000000    O    (16i)
  -0.00675000000000   -0.04830000000000   -0.20971000000000    O    (16i)
  -0.20296000000000    0.00675000000000   -0.04155000000000    O    (16i)
   0.04830000000000   -0.16141000000000    0.04155000000000    O    (16i)
   0.49325000000000    0.70296000000000    0.04155000000000    O    (16i)
   0.66141000000000    0.45170000000000   -0.04155000000000    O    (16i)
   0.54830000000000    0.50675000000000    0.20971000000000    O    (16i)
   0.29704000000000    0.33859000000000   -0.20971000000000    O    (16i)
   0.56653000000000    0.30851000000000    0.53318000000000    O    (16i)
```

```
 −0.22467000000000   0.03335000000000  −0.53318000000000   O   (16i)
 −0.30851000000000   0.22467000000000   0.25802000000000   O   (16i)
 −0.03335000000000  −0.56653000000000  −0.25802000000000   O   (16i)
  0.27533000000000   0.80851000000000  −0.25802000000000   O   (16i)
  1.06653000000000   0.53335000000000   0.25802000000000   O   (16i)
  0.46665000000000   0.72467000000000   0.53318000000000   O   (16i)
  0.19149000000000  −0.06653000000000  −0.53318000000000   O   (16i)
  0.00000000000000   0.00000000000000   0.00000000000000   S    (4b)
  0.50000000000000   0.50000000000000   0.00000000000000   S    (4b)
```

$SrCu_2(BO_3)_2$: A2B2C6D_tI44_121_i_i_ij_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'B2Cu2O6Sr'
_chemical_formula_sum 'B2 Cu2 O6 Sr'

loop_
_publ_author_name
 'R. W. Smith'
 'D. A. Keszler'
_journal_name_full_name
;
 Journal of Solid State Chemistry
;
_journal_volume 93
_journal_year 1991
_journal_page_first 430
_journal_page_last 435
_publ_Section_title
;
 Synthesis, structure, and properties of the orthoborate SrCu$_{2}$(BO$_
    ↪ {3}$)$_{2}$
;

# Found in Exact Dimer Ground State and Quantized Magnetization Plateaus
    ↪ in the Two-Dimensional Spin System SrCu$_{2}$(BO$_{3}$)$_{2}$,
    ↪ 1999

_aflow_title 'SrCu$_{2}$(BO$_{3}$)$_{2}$ Structure'
_aflow_proto 'A2B2C6D_tI44_121_i_i_ij_c'
_aflow_params 'a,c/a,x_{2},z_{2},x_{3},z_{3},x_{4},z_{4},x_{5},y_{5},z_{
    ↪ 5}'
_aflow_params_values '8.995,0.739188438021,0.2953,0.243,0.11412,0.2783,
    ↪ 0.4004,0.212,0.3276,0.1456,0.254'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tI44'

_symmetry_space_group_name_H-M "I -4 2 m"
_symmetry_Int_Tables_number 121

_cell_length_a    8.99500
_cell_length_b    8.99500
_cell_length_c    6.64900
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 y,x,z
6 y,-x,-z
7 -y,x,-z
8 -y,-x,z
9 x+1/2,y+1/2,z+1/2
10 x+1/2,-y+1/2,-z+1/2
11 -x+1/2,y+1/2,-z+1/2
12 -x+1/2,-y+1/2,z+1/2
13 y+1/2,x+1/2,z+1/2
14 y+1/2,-x+1/2,-z+1/2
15 -y+1/2,x+1/2,-z+1/2
16 -y+1/2,-x+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Sr1 Sr 4 c 0.00000 0.50000 0.00000 1.00000
B1 B 8 i 0.29530 0.29530 0.24300 1.00000
Cu1 Cu 8 i 0.11412 0.11412 0.27830 1.00000
O1 O 8 i 0.40040 0.40040 0.21200 1.00000
O2 O 16 j 0.32760 0.14560 0.25400 1.00000
```

$SrCu_2(BO_3)_2$: A2B2C6D_tI44_121_i_i_ij_c - POSCAR

```
A2B2C6D_tI44_121_i_i_ij_c & a,c/a,x2,z2,x3,z3,x4,z4,x5,y5,z5 --params=
    ↪ 8.995,0.739188438021,0.2953,0.243,0.11412,0.2783,0.4004,0.212,
    ↪ 0.3276,0.1456,0.254 & I-42m D_{2d}^{11} #121 (ci^3j) & tI44 &
    ↪ None & B2Cu2O6Sr & B2Cu2O6Sr & R. W. Smith and D. A. Keszler,
    ↪ J. Solid State Chem. 93, 430-435 (1991)
   1.00000000000000
  -4.49750000000000   4.49750000000000   3.32450000000000
   4.49750000000000  -4.49750000000000   3.32450000000000
   4.49750000000000   4.49750000000000  -3.32450000000000
      B     Cu      O     Sr
      4      4     12      2
Direct
   0.53830000000000   0.53830000000000   0.59060000000000   B   (8i)
  -0.05230000000000  -0.05230000000000  -0.59060000000000   B   (8i)
  -0.53830000000000   0.05230000000000   0.00000000000000   B   (8i)
   0.05230000000000  -0.53830000000000   0.00000000000000   B   (8i)
   0.39242000000000   0.39242000000000   0.22824000000000   Cu  (8i)
   0.16418000000000   0.16418000000000  -0.22824000000000   Cu  (8i)
  -0.39242000000000  -0.16418000000000   0.00000000000000   Cu  (8i)
  -0.16418000000000  -0.39242000000000   0.00000000000000   Cu  (8i)
   0.61240000000000   0.61240000000000   0.80080000000000   O   (8i)
  -0.18840000000000  -0.18840000000000  -0.80080000000000   O   (8i)
  -0.61240000000000   0.18840000000000   0.00000000000000   O   (8i)
   0.18840000000000  -0.61240000000000   0.00000000000000   O   (8i)
   0.39960000000000   0.58160000000000   0.47320000000000   O  (16j)
   0.10840000000000  -0.07360000000000  -0.47320000000000   O  (16j)
  -0.58160000000000  -0.10840000000000  -0.18200000000000   O  (16j)
   0.07360000000000  -0.39960000000000   0.18200000000000   O  (16j)
  -0.10840000000000  -0.58160000000000  -0.18200000000000   O  (16j)
  -0.39960000000000   0.07360000000000   0.18200000000000   O  (16j)
  -0.07360000000000   0.10840000000000  -0.47320000000000   O  (16j)
   0.58160000000000   0.39960000000000   0.47320000000000   O  (16j)
   0.50000000000000   0.00000000000000   0.50000000000000   Sr  (4c)
   0.00000000000000   0.50000000000000   0.50000000000000   Sr  (4c)
```

$C17$ (Fe_2B) (*obsolete*): AB2_tI12_121_ab_i - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'BFe2'
_chemical_formula_sum 'B Fe2'

loop_
_publ_author_name
 'F. Wever'
 'A. M\"{u}ller'
_journal_name_full_name
;
 Zeitschrift fur Anorganische und Allgemeine Chemie
;
_journal_volume 192
_journal_year 1930
_journal_page_first 317
_journal_page_last 336
_publ_Section_title
;
 \"{U}ber das Zweistoffsystem Eisen-Bor und \"{u}ber die Struktur des
    ↪ Eisenborides Fe$_{4}$B$_{2}$
;

_aflow_title '$C17$ (Fe$_{2}$B) ({\em{obsolete}}) Structure'
_aflow_proto 'AB2_tI12_121_ab_i'
_aflow_params 'a,c/a,x_{3},z_{3}'
_aflow_params_values '5.078,0.831626624655,0.16667,0.2'
_aflow_Strukturbericht '$C17$'
_aflow_Pearson 'tI12'

_symmetry_space_group_name_H-M "I -4 2 m"
_symmetry_Int_Tables_number 121

_cell_length_a    5.07800
_cell_length_b    5.07800
_cell_length_c    4.22300
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 y,x,z
6 y,-x,-z
7 -y,x,-z
8 -y,-x,z
9 x+1/2,y+1/2,z+1/2
10 x+1/2,-y+1/2,-z+1/2
11 -x+1/2,y+1/2,-z+1/2
12 -x+1/2,-y+1/2,z+1/2
13 y+1/2,x+1/2,z+1/2
14 y+1/2,-x+1/2,-z+1/2
15 -y+1/2,x+1/2,-z+1/2
16 -y+1/2,-x+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
B1 B 2 a 0.00000 0.00000 0.00000 1.00000
B2 B 2 b 0.00000 0.00000 0.50000 1.00000
Fe1 Fe 8 i 0.16667 0.16667 0.20000 1.00000
```

$C17$ (Fe_2B) (*obsolete*): AB2_tI12_121_ab_i - POSCAR

```
AB2_tI12_121_ab_i & a,c/a,x3,z3 --params=5.078,0.831626624655,0.16667,
    ↪ 0.2 & I-42m D_{2d}^{11} #121 (abi) & tI12 & $C17$ & BFe2 & BFe2
    ↪ & F. Wever and A. M\"{u}ller, Z. Anorg. Allg. Chem. 192,
    ↪ 317-336 (1930)
  1.00000000000000
 -2.53900000000000   2.53900000000000   2.11150000000000
  2.53900000000000  -2.53900000000000   2.11150000000000
  2.53900000000000   2.53900000000000  -2.11150000000000
    B     Fe
    2     4
Direct
  0.00000000000000   0.00000000000000   0.00000000000000   B    (2a)
  0.50000000000000   0.50000000000000   0.00000000000000   B    (2b)
  0.36667000000000   0.36667000000000   0.33334000000000   Fe   (8i)
  0.03333000000000   0.03333000000000  -0.33334000000000   Fe   (8i)
 -0.36667000000000  -0.03333000000000   0.00000000000000   Fe   (8i)
 -0.03333000000000  -0.36667000000000   0.00000000000000   Fe   (8i)
```

K_3CrO_8: AB3C8_tI24_121_a_bd_2i - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'CrK3O8'
_chemical_formula_sum 'Cr K3 O8'

loop_
_publ_author_name
 'R. Stomberg'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 17
_journal_year 1963
_journal_page_first 1563
_journal_page_last 1566
_publ_Section_title
;
 Least-Squares Refinement of the Crystal Structure of Potassium
    ↪ Peroxochromate
;

_aflow_title 'K$_{3}$CrO$_{8}$ Structure'
_aflow_proto 'AB3C8_tI24_121_a_bd_2i'
_aflow_params 'a,c/a,x_{4},z_{4},x_{5},z_{5}'
_aflow_params_values '6.703,1.13859465911,0.1355,0.1788,0.2079,0.0082'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tI24'

_symmetry_space_group_name_H-M "I -4 2 m"
_symmetry_Int_Tables_number 121

_cell_length_a    6.70300
_cell_length_b    6.70300
_cell_length_c    7.63200
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 y,x,z
6 y,-x,-z
7 -y,x,-z
8 -y,-x,z
9 x+1/2,y+1/2,z+1/2
10 x+1/2,-y+1/2,-z+1/2
11 -x+1/2,y+1/2,-z+1/2
12 -x+1/2,-y+1/2,z+1/2
13 y+1/2,x+1/2,z+1/2
14 y+1/2,-x+1/2,-z+1/2
15 -y+1/2,x+1/2,-z+1/2
16 -y+1/2,-x+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cr1 Cr 2 a 0.00000 0.00000 0.00000 1.00000
K1 K 2 b 0.00000 0.00000 0.50000 1.00000
K2 K 4 d 0.00000 0.50000 0.25000 1.00000
O1 O 8 i 0.13550 0.13550 0.17880 1.00000
O2 O 8 i 0.20790 0.20790 0.00820 1.00000
```

K_3CrO_8: AB3C8_tI24_121_a_bd_2i - POSCAR

```
AB3C8_tI24_121_a_bd_2i & a,c/a,x4,z4,x5,z5 --params=6.703,1.13859465911,
    ↪ 0.1355,0.1788,0.2079,0.0082 & I-42m D_{2d}^{11} #121 (abdi^2) &
    ↪  tI24 & None & CrK3O8 & CrK3O8 & R. Stomberg, Acta Chem. Scand.
    ↪ 17, 1563-1566 (1963)
  1.00000000000000
 -3.35150000000000   3.35150000000000   3.81600000000000
  3.35150000000000  -3.35150000000000   3.81600000000000
  3.35150000000000   3.35150000000000  -3.81600000000000
   Cr    K    O
    1    3    8
Direct
  0.00000000000000   0.00000000000000   0.00000000000000   Cr   (2a)
  0.50000000000000   0.50000000000000   0.00000000000000   K    (2b)
  0.75000000000000   0.25000000000000   0.50000000000000   K    (4d)
  0.25000000000000   0.75000000000000   0.50000000000000   K    (4d)
  0.31430000000000   0.31430000000000   0.27100000000000   O    (8i)
  0.04330000000000   0.04330000000000  -0.27100000000000   O    (8i)
 -0.31430000000000  -0.04330000000000   0.00000000000000   O    (8i)
 -0.04330000000000  -0.31430000000000   0.00000000000000   O    (8i)
  0.21610000000000   0.21610000000000   0.41580000000000   O    (8i)
 -0.19970000000000  -0.19970000000000  -0.41580000000000   O    (8i)
 -0.21610000000000   0.19970000000000   0.00000000000000   O    (8i)
  0.19970000000000  -0.21610000000000   0.00000000000000   O    (8i)
```

α-V_3S: AB3_tI32_121_g_f2i - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'SV3'
_chemical_formula_sum 'S V3'

loop_
_publ_author_name
 'B. Pedersen'
 'F. Gr{\o}nvold'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 12
_journal_year 1959
_journal_page_first 1022
_journal_page_last 1027
_publ_Section_title
;
 The Crystal Structures of $\alpha$-V$_{3}$S and $\beta$-V$_{3}$S
;

_aflow_title '$\alpha$-V$_{3}$S Structure'
_aflow_proto 'AB3_tI32_121_g_f2i'
_aflow_params 'a,c/a,x_{1},x_{2},x_{3},z_{3},x_{4},z_{4}'
_aflow_params_values '9.47,0.484582893347,0.355,0.2851,0.5932,0.25,0.2,
    ↪ 0.25'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tI32'

_symmetry_space_group_name_H-M "I -4 2 m"
_symmetry_Int_Tables_number 121

_cell_length_a    9.47000
_cell_length_b    9.47000
_cell_length_c    4.58900
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 y,x,z
6 y,-x,-z
7 -y,x,-z
8 -y,-x,z
9 x+1/2,y+1/2,z+1/2
10 x+1/2,-y+1/2,-z+1/2
11 -x+1/2,y+1/2,-z+1/2
12 -x+1/2,-y+1/2,z+1/2
13 y+1/2,x+1/2,z+1/2
14 y+1/2,-x+1/2,-z+1/2
15 -y+1/2,x+1/2,-z+1/2
16 -y+1/2,-x+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
V1 V 8 f 0.35500 0.00000 0.00000 1.00000
S1 S 8 g 0.28510 0.00000 0.50000 1.00000
V2 V 8 i 0.59320 0.59320 0.25000 1.00000
V3 V 8 i 0.20000 0.20000 0.25000 1.00000
```

α-V_3S: AB3_tI32_121_g_f2i - POSCAR

```
AB3_tI32_121_g_f2i & a,c/a,x1,x2,x3,z3,x4,z4 --params=9.47,
    ↪ 0.484582893347,0.355,0.2851,0.5932,0.25,0.2,0.25 & I-42m D_{2d
    ↪ }^{11} #121 (fgi^2) & tI32 & None & SV3 & SV3 & B. Pedersen and
    ↪  F. Gr{\o}nvold, Acta Cryst. 12, 1022-1027 (1959)
  1.00000000000000
 -4.73500000000000   4.73500000000000   2.29450000000000
  4.73500000000000  -4.73500000000000   2.29450000000000
```

```
    4.73500000000000    4.73500000000000   -2.29450000000000
      S       V
      4      12
Direct
    0.50000000000000    0.78510000000000    0.28510000000000   S   (8g)
    0.50000000000000    0.21490000000000   -0.28510000000000   S   (8g)
    0.21490000000000    0.50000000000000   -0.28510000000000   S   (8g)
    0.78510000000000    0.50000000000000    0.28510000000000   S   (8g)
    0.00000000000000    0.35500000000000    0.35500000000000   V   (8f)
    0.00000000000000   -0.35500000000000   -0.35500000000000   V   (8f)
   -0.35500000000000    0.00000000000000   -0.35500000000000   V   (8f)
    0.35500000000000    0.00000000000000    0.35500000000000   V   (8f)
    0.84320000000000    0.84320000000000    1.18640000000000   V   (8i)
   -0.34320000000000   -0.34320000000000   -1.18640000000000   V   (8i)
   -0.84320000000000    0.34320000000000    0.00000000000000   V   (8i)
    0.34320000000000   -0.84320000000000    0.00000000000000   V   (8i)
    0.45000000000000    0.45000000000000    0.40000000000000   V   (8i)
    0.05000000000000    0.05000000000000   -0.40000000000000   V   (8i)
   -0.45000000000000   -0.05000000000000    0.00000000000000   V   (8i)
   -0.05000000000000   -0.45000000000000    0.00000000000000   V   (8i)
```

Mercury Cyanide [$Hg(CN)_2$, $F1_1$]: A2BC2_tI40_122_e_d_e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Mercury cyanide'
_chemical_formula_sum 'C2 Hg N2'

loop_
_publ_author_name
 'O. Reckeweg'
 'A. Simon'
_journal_name_full_name
;
 Zeitschrift f\"u r Naturforschung B
;
_journal_volume 57
_journal_year 2002
_journal_page_first 895
_journal_page_last 900
_publ_Section_title
;
 X-Ray and Raman Investigations on Cyanides of Mono- and Divalent Metals
      ↪  and Synthesis, Crystal Structure and Raman Spectrum of Tl$_{5
      ↪ }$(CO$_{3}$)$_{2}$(CN)
;

# Found in Hg(CN)$_{2}$ (Hg[CN]$_{2}$) Crystal Structure, {PAULING FILE
      ↪ in: Inorganic Solid Phases, SpringerMaterials (online database)
      ↪ }, 2016

_aflow_title 'Mercury Cyanide [Hg(CN)$_{2}$, $F1_{1}$] Structure'
_aflow_proto 'A2BC2_tI40_122_e_d_e'
_aflow_params 'a,c/a,x_{1},x_{2},y_{2},z_{2},x_{3},y_{3},z_{3}'
_aflow_params_values '9.6922,0.918418934814,0.21203,0.2035,-0.0729,
      ↪ 0.1785,0.205,0.0464,0.1639'
_aflow_Strukturbericht '$F1_{1}$'
_aflow_Pearson 'tI40'

_symmetry_space_group_name_H-M "I -4 2 d"
_symmetry_Int_Tables_number 122

_cell_length_a     9.69220
_cell_length_b     9.69220
_cell_length_c     8.90150
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y+1/2,-z+1/4
3 -x,y+1/2,-z+1/4
4 -x,-y,z
5 y,x+1/2,z+1/4
6 y,-x,-z
7 -y,x,-z
8 -y,-x+1/2,z+1/4
9 x+1/2,y+1/2,z+1/2
10 x+1/2,-y,-z+3/4
11 -x+1/2,y,-z+3/4
12 -x+1/2,-y+1/2,z+1/2
13 y+1/2,x,z+3/4
14 y+1/2,-x+1/2,-z+1/2
15 -y+1/2,x+1/2,-z+1/2
16 -y+1/2,-x,z+3/4

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Hg1 Hg 8 d 0.21203 0.25000 0.12500 1.00000
C1 C 16 e 0.20350 -0.07290 0.17850 1.00000
N1 N 16 e 0.20500 0.04640 0.16390 1.00000
```

Mercury Cyanide [$Hg(CN)_2$, $F1_1$]: A2BC2_tI40_122_e_d_e - POSCAR

```
A2BC2_tI40_122_e_d_e & a,c/a,x1,x2,y2,z2,x3,y3,z3 --params=9.6922,
      ↪ 0.918418934814,0.21203,0.2035,-0.0729,0.1785,0.205,0.0464,
      ↪ 0.1639 & I-42d D_{2d}^{12} #122 (de^2) & tI40 & $F1_{1}$ &
      ↪ C2HgN2 & Mercury cyanide & O. Reckeweg and A. Simon, Z.
      ↪ Naturforsch. B 57, 895-900 (2002)
   1.00000000000000
  -4.84610000000000    4.84610000000000    4.45075000000000
   4.84610000000000   -4.84610000000000    4.45075000000000
   4.84610000000000    4.84610000000000   -4.45075000000000
     C      Hg      N
     8       4       8
Direct
    0.10560000000000    0.38200000000000    0.13060000000000   C   (16e)
    0.25140000000000   -0.02500000000000   -0.13060000000000   C   (16e)
   -0.38200000000000   -0.25140000000000   -0.27640000000000   C   (16e)
    0.02500000000000   -0.10560000000000    0.27640000000000   C   (16e)
    0.49860000000000   -0.13200000000000    0.22360000000000   C   (16e)
    0.64440000000000    0.27500000000000    0.77640000000000   C   (16e)
    0.72500000000000    0.50140000000000    0.36940000000000   C   (16e)
    1.13200000000000    0.35560000000000    0.63060000000000   C   (16e)
    0.37500000000000    0.33703000000000    0.46203000000000   Hg   (8d)
    0.87500000000000   -0.08703000000000    0.53797000000000   Hg   (8d)
    0.66297000000000    0.12500000000000    0.03797000000000   Hg   (8d)
    1.08703000000000    0.62500000000000    0.96203000000000   Hg   (8d)
    0.21030000000000    0.36890000000000    0.25140000000000   N   (16e)
    0.11750000000000   -0.04110000000000   -0.25140000000000   N   (16e)
   -0.36890000000000   -0.11750000000000   -0.15860000000000   N   (16e)
    0.04110000000000   -0.21030000000000    0.15860000000000   N   (16e)
    0.63250000000000   -0.11890000000000    0.34140000000000   N   (16e)
    0.53970000000000    0.29110000000000    0.65860000000000   N   (16e)
    0.70890000000000    0.36750000000000    0.24860000000000   N   (16e)
    1.11890000000000    0.46030000000000    0.75140000000000   N   (16e)
```

KH_2PO_4 ($H2_2$): A4BC4D_tI40_122_e_b_e_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'KH2PO4'
_chemical_formula_sum 'H4 K O4 P'

loop_
_publ_author_name
 'R. J. Nelmes'
 'G. M. Meyer'
 'J. E. Tibballs'
_journal_name_full_name
;
 Journal of Physics C: Solid State Physics
;
_journal_volume 15
_journal_year 1982
_journal_page_first 59
_journal_page_last 75
_publ_Section_title
;
 The crystal structure of tetragonal KH$_{2}$PO$_{4}$ and KD$_{2}$PO$_{4
      ↪ }$ as a function of temperature
;

_aflow_title 'KH$_{2}$PO$_{4}$ ($H2_{2}$) Structure'
_aflow_proto 'A4BC4D_tI40_122_e_b_e_a'
_aflow_params 'a,c/a,x_{3},y_{3},z_{3},x_{4},y_{4},z_{4}'
_aflow_params_values '7.4264,0.933292039211,0.14867,0.22713,0.12266,
      ↪ 0.1933,0.08283,0.12675'
_aflow_Strukturbericht '$H2_{2}$'
_aflow_Pearson 'tI40'

_symmetry_space_group_name_H-M "I -4 2 d"
_symmetry_Int_Tables_number 122

_cell_length_a     7.42640
_cell_length_b     7.42640
_cell_length_c     6.93100
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y+1/2,-z+1/4
3 -x,y+1/2,-z+1/4
4 -x,-y,z
5 y,x+1/2,z+1/4
6 y,-x,-z
7 -y,x,-z
8 -y,-x+1/2,z+1/4
9 x+1/2,y+1/2,z+1/2
10 x+1/2,-y,-z+3/4
11 -x+1/2,y,-z+3/4
12 -x+1/2,-y+1/2,z+1/2
13 y+1/2,x,z+3/4
14 y+1/2,-x+1/2,-z+1/2
15 -y+1/2,x+1/2,-z+1/2
16 -y+1/2,-x,z+3/4

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
```

```
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
P1 P 4 a 0.00000 0.00000 0.00000 1.00000
K1 K 4 b 0.00000 0.00000 0.50000 1.00000
H1 H 16 e 0.14867 0.22713 0.12266 0.50000
O1 O 16 e 0.19330 0.08283 0.12675 1.00000
```

KH_2PO_4 ($H2_2$): A4BC4D_tI40_122_e_b_e_a - POSCAR

```
A4BC4D_tI40_122_e_b_e_a & a,c/a,x3,y3,z3,x4,y4,z4 --params=7.4264,
      ↪ 0.933292039211,0.14867,0.22713,0.12266,0.1933,0.08283,0.12675 &
      ↪  I-42d D_{2d}^{12} #122 (abe^2) & tI40 & $H2_{2}$ & KH2PO4 &
      ↪ KH2PO4 & R. J. Nelmes and G. M. Meyer and J. E. Tibballs, J.
      ↪ Phys. C: Solid State Phys. 15, 59-75 (1982)
   1.00000000000000
  -3.71320000000000    3.71320000000000    3.46550000000000
   3.71320000000000   -3.71320000000000    3.46550000000000
   3.71320000000000    3.71320000000000   -3.46550000000000
     H     K     O     P
     8     2     8     2
Direct
   0.34979000000000    0.27133000000000    0.37580000000000    H   (16e)
  -0.10447000000000   -0.02601000000000   -0.37580000000000    H   (16e)
  -0.27133000000000    0.10447000000000    0.07846000000000    H   (16e)
   0.02601000000000   -0.34979000000000   -0.07846000000000    H   (16e)
   0.85447000000000   -0.02133000000000    0.57846000000000    H   (16e)
   0.40021000000000    0.27601000000000    0.42154000000000    H   (16e)
   0.72399000000000    0.14553000000000    0.12420000000000    H   (16e)
   1.02133000000000    0.59979000000000    0.87580000000000    H   (16e)
   0.50000000000000    0.50000000000000    0.00000000000000    K    (4b)
   0.25000000000000    0.75000000000000    0.50000000000000    K    (4b)
   0.20958000000000    0.32005000000000    0.27613000000000    O   (16e)
   0.04392000000000   -0.06655000000000   -0.27613000000000    O   (16e)
  -0.32005000000000   -0.04392000000000   -0.11047000000000    O   (16e)
   0.06655000000000   -0.20958000000000    0.11047000000000    O   (16e)
   0.70608000000000   -0.07005000000000    0.38953000000000    O   (16e)
   0.54042000000000    0.31655000000000    0.61047000000000    O   (16e)
   0.68345000000000    0.29392000000000    0.22387000000000    O   (16e)
   1.07005000000000    0.45958000000000    0.77613000000000    O   (16e)
   0.00000000000000    0.00000000000000    0.00000000000000    P    (4a)
   0.75000000000000    0.25000000000000    0.50000000000000    P    (4a)
```

$NH_4H_2PO_4$: A8BC4D_tI56_122_2e_b_e_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'H6NO4P'
_chemical_formula_sum 'H8 N O4 P'

loop_
_publ_author_name
 'A. A. Khan'
 'W. H. Baur'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 29
_journal_year 1973
_journal_page_first 2721
_journal_page_last 2726
_publ_Section_title
;
 Refinement of the crystal structures of ammonium dihydrogen phosphate
      ↪ and ammonium dihydrogen arsenate
;

# Found in Refinement of the Crystal Structure of NH$_{4}$H$_{2}$PO$_{4}
      ↪ $ above and below Antiferroelectric Phase Transition
      ↪ Temperature, 1987

_aflow_title 'NH$_{4}$H$_{2}$PO$_{4}$ Structure'
_aflow_proto 'A8BC4D_tI56_122_2e_b_e_a'
_aflow_params 'a,c/a,x_{3},y_{3},z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{
      ↪ 5}'
_aflow_params_values '7.4997,1.00662693174,0.25,0.15,0.125,0.498,0.589,
      ↪ 0.063,0.0843,0.1466,0.1151'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tI56'

_symmetry_space_group_name_H-M "I -4 2 d"
_symmetry_Int_Tables_number 122

_cell_length_a     7.49970
_cell_length_b     7.49970
_cell_length_c     7.54940
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y+1/2,-z+1/4
3 -x,y+1/2,-z+1/4
4 -x,-y,z
5 y,x+1/2,z+1/4
6 y,-x,-z
7 -y,x,-z
8 -y,-x+1/2,z+1/4
9 x+1/2,y+1/2,z+1/2
10 x+1/2,-y,-z+3/4
11 -x+1/2,y,-z+3/4
12 -x+1/2,-y+1/2,z+1/2
13 y+1/2,x,z+3/4
14 y+1/2,-x+1/2,-z+1/2
15 -y+1/2,x+1/2,-z+1/2
16 -y+1/2,-x,z+3/4

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
P1 P 4 a 0.00000 0.00000 0.00000 1.00000
N1 N 4 b 0.00000 0.00000 0.50000 1.00000
H1 H 16 e 0.25000 0.15000 0.12500 0.50000
H2 H 16 e 0.49800 0.58900 0.06300 1.00000
O1 O 16 e 0.08430 0.14660 0.11510 1.00000
```

$NH_4H_2PO_4$: A8BC4D_tI56_122_2e_b_e_a - POSCAR

```
A8BC4D_tI56_122_2e_b_e_a & a,c/a,x3,y3,z3,x4,y4,z4,x5,y5,z5 --params=
      ↪ 7.4997,1.00662693174,0.25,0.15,0.125,0.498,0.589,0.063,0.0843,
      ↪ 0.1466,0.1151 & I-42d D_{2d}^{12} #122 (abe^3) & tI56 & None &
      ↪ H6NO4P & H6NO4P & A. A. Khan and W. H. Baur, Acta Crystallogr.
      ↪ Sect. B Struct. Sci. 29, 2721-2726 (1973)
   1.00000000000000
  -3.74985000000000    3.74985000000000    3.77470000000000
   3.74985000000000   -3.74985000000000    3.77470000000000
   3.74985000000000    3.74985000000000   -3.77470000000000
     H     N     O     P
    16     2     8     2
Direct
   0.27500000000000    0.37500000000000    0.40000000000000    H   (16e)
  -0.02500000000000   -0.12500000000000   -0.40000000000000    H   (16e)
  -0.37500000000000    0.02500000000000   -0.10000000000000    H   (16e)
   0.12500000000000   -0.27500000000000    0.10000000000000    H   (16e)
   0.77500000000000   -0.12500000000000    0.40000000000000    H   (16e)
   0.47500000000000    0.37500000000000    0.60000000000000    H   (16e)
   0.62500000000000    0.22500000000000    0.10000000000000    H   (16e)
   1.12500000000000    0.52500000000000    0.90000000000000    H   (16e)
   0.65200000000000    0.56100000000000    1.08700000000000    H   (16e)
  -0.52600000000000   -0.43500000000000   -1.08700000000000    H   (16e)
  -0.56100000000000    0.52600000000000    0.09100000000000    H   (16e)
   0.43500000000000   -0.65200000000000   -0.09100000000000    H   (16e)
   1.27600000000000   -0.31100000000000    0.59100000000000    H   (16e)
   0.09800000000000    0.68500000000000    0.40900000000000    H   (16e)
   0.31500000000000   -0.27600000000000   -0.58700000000000    H   (16e)
   1.31100000000000    0.90200000000000    1.58700000000000    H   (16e)
   0.50000000000000    0.50000000000000    0.00000000000000    N    (4b)
   0.25000000000000    0.75000000000000    0.50000000000000    N    (4b)
   0.26170000000000    0.19940000000000    0.23090000000000    O   (16e)
  -0.03150000000000    0.03080000000000   -0.23090000000000    O   (16e)
  -0.19940000000000    0.03150000000000    0.06230000000000    O   (16e)
  -0.03080000000000   -0.26170000000000   -0.06230000000000    O   (16e)
   0.78150000000000    0.05060000000000    0.56230000000000    O   (16e)
   0.48830000000000    0.21920000000000    0.43770000000000    O   (16e)
   0.78080000000000    0.21850000000000    0.26910000000000    O   (16e)
   0.94940000000000    0.51170000000000    0.73090000000000    O   (16e)
   0.00000000000000    0.00000000000000    0.00000000000000    P    (4a)
   0.75000000000000    0.25000000000000    0.50000000000000    P    (4a)
```

NaS_2: AB2_tI48_122_cd_2e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'NaS2'
_chemical_formula_sum 'Na S2'

loop_
_publ_author_name
 'R. Tegman'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 29
_journal_year 1973
_journal_page_first 1463
_journal_page_last 1469
_publ_Section_title
;
 The Crystal Structure of Sodium Tetrasulphide, Na$_{2}$S$_{4}$
;

# Found in Pearson's Handbook of Crystallographic Data for Intermetallic
      ↪ Phases, 1991

_aflow_title 'NaS$_{2}$ Structure'
_aflow_proto 'AB2_tI48_122_cd_2e'
_aflow_params 'a,c/a,z_{1},x_{2},x_{3},y_{3},z_{3},x_{4},y_{4},z_{4}'
_aflow_params_values '9.5965,1.22841661022,0.16953,0.27647,0.03373,
      ↪ 0.24952,0.29644,0.14693,0.116,0.39956'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tI48'

_symmetry_space_group_name_H-M "I -4 2 d"
_symmetry_Int_Tables_number 122
```

```
_cell_length_a     9.59650
_cell_length_b     9.59650
_cell_length_c     11.78850
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y+1/2,-z+1/4
3 -x,y+1/2,-z+1/4
4 -x,-y,z
5 y,x+1/2,z+1/4
6 y,-x,-z
7 -y,x,-z
8 -y,-x+1/2,z+1/4
9 x+1/2,y+1/2,z+1/2
10 x+1/2,-y,-z+3/4
11 -x+1/2,y,-z+3/4
12 -x+1/2,-y+1/2,z+1/2
13 y+1/2,x,z+3/4
14 y+1/2,-x+1/2,-z+1/2
15 -y+1/2,x+1/2,-z+1/2
16 -y+1/2,-x,z+3/4

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Na1 Na 8 c 0.00000 0.00000 0.16953 1.00000
Na2 Na 8 d 0.27647 0.25000 0.12500 1.00000
S1 S 16 e 0.03373 0.24952 0.29644 1.00000
S2 S 16 e 0.14693 0.11600 0.39956 1.00000
```

NaS_2: AB2_tI48_122_cd_2e - POSCAR

```
AB2_tI48_122_cd_2e & a,c/a,z1,x2,x3,y3,z3,x4,y4,z4 --params=9.5965,
      ↪ 1.22841661022,0.16953,0.27647,0.03373,0.24952,0.29644,0.14693,
      ↪ 0.116,0.39956 & I-42d D_{2d}^{12} #122 (cde^2) & tI48 & None &
      ↪ NaS2 & NaS2 & R. Tegman, Acta Crystallogr. Sect. B Struct. Sci.
      ↪ 29, 1463-1469 (1973)
   1.00000000000000
  -4.79825000000000    4.79825000000000    5.89425000000000
   4.79825000000000   -4.79825000000000    5.89425000000000
   4.79825000000000    4.79825000000000   -5.89425000000000
     Na      S
      8     16
Direct
   0.16953000000000    0.16953000000000    0.00000000000000    Na    (8c)
  -0.16953000000000   -0.16953000000000    0.00000000000000    Na    (8c)
   0.58047000000000    0.08047000000000    0.50000000000000    Na    (8c)
   0.91953000000000    0.41953000000000    0.50000000000000    Na    (8c)
   0.37500000000000    0.40147000000000    0.52647000000000    Na    (8d)
   0.87500000000000   -0.15147000000000    0.47353000000000    Na    (8d)
   0.59853000000000    0.12500000000000   -0.02647000000000    Na    (8d)
   1.15147000000000    0.62500000000000    1.02647000000000    Na    (8d)
   0.54596000000000    0.33017000000000    0.28325000000000    S    (16e)
   0.04692000000000    0.26271000000000   -0.28325000000000    S    (16e)
  -0.33017000000000   -0.04692000000000    0.21579000000000    S    (16e)
  -0.26271000000000   -0.54596000000000   -0.21579000000000    S    (16e)
   0.70308000000000   -0.08017000000000    0.71579000000000    S    (16e)
   0.20404000000000   -0.01271000000000    0.28421000000000    S    (16e)
   1.01271000000000    0.29692000000000    0.21675000000000    S    (16e)
   1.08017000000000    0.79596000000000    0.78325000000000    S    (16e)
   0.51556000000000    0.54649000000000    0.26293000000000    S    (16e)
   0.28356000000000    0.25263000000000   -0.26293000000000    S    (16e)
  -0.54649000000000   -0.28356000000000   -0.03093000000000    S    (16e)
  -0.25263000000000   -0.51556000000000    0.03093000000000    S    (16e)
   0.46644000000000   -0.29649000000000    0.46907000000000    S    (16e)
   0.23444000000000   -0.00263000000000    0.53093000000000    S    (16e)
   1.00263000000000    0.53356000000000    0.23707000000000    S    (16e)
   1.29649000000000    0.76556000000000    0.76293000000000    S    (16e)
```

NH_4HgCl_3 ($E2_5$): A3BC_tP5_123_cg_a_d - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Cl3Hg(NH4)'
_chemical_formula_sum 'Cl3 Hg (NH4)'

loop_
_publ_author_name
 'E. J. Harmsen'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 100
_journal_year 1939
_journal_page_first 208
_journal_page_last 211
_publ_Section_title
;
 The Crystal Structure of NH$_{4}$HgCl$_{3}$
;
```

```
# Found in Strukturbericht Band VII 1939, 1943

_aflow_title 'NH$_{4}$HgCl$_{3}$ ($E2_{5}$) Structure'
_aflow_proto 'A3BC_tP5_123_cg_a_d'
_aflow_params 'a,c/a,z_{4}'
_aflow_params_values '4.19,1.89498806683,0.294'
_aflow_Strukturbericht '$E2_{5}$'
_aflow_Pearson 'tP5'

_symmetry_space_group_name_H-M "P 4/m 2/m 2/m"
_symmetry_Int_Tables_number 123

_cell_length_a     4.19000
_cell_length_b     4.19000
_cell_length_c     7.94000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 -y,-x,-z
6 -y,x,z
7 y,-x,z
8 y,x,-z
9 -x,-y,-z
10 -x,y,z
11 x,-y,z
12 x,y,-z
13 y,x,z
14 y,-x,-z
15 -y,x,-z
16 -y,-x,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Hg1 Hg 1 a 0.00000 0.00000 0.00000 1.00000
Cl1 Cl 1 c 0.50000 0.50000 0.00000 1.00000
NH41 NH4 1 d 0.50000 0.50000 0.50000 1.00000
Cl2 Cl 2 g 0.00000 0.00000 0.29400 1.00000
```

NH_4HgCl_3 ($E2_5$): A3BC_tP5_123_cg_a_d - POSCAR

```
A3BC_tP5_123_cg_a_d & a,c/a,z4 --params=4.19,1.89498806683,0.294 & P4/
      ↪ mmm D_{4h}^{1} #123 (acdg) & tP5 & $E2_{5}$ & Cl3Hg(NH4) &
      ↪ Cl3Hg(NH4) & E. J. Harmsen, Zeitschrift f{"u}r Kristallographie
      ↪ - Crystalline Materials 100, 208-211 (1939)
   1.00000000000000
   4.19000000000000    0.00000000000000    0.00000000000000
   0.00000000000000    4.19000000000000    0.00000000000000
   0.00000000000000    0.00000000000000    7.94000000000000
     Cl     Hg    NH4
      3      1      1
Direct
   0.50000000000000    0.50000000000000    0.00000000000000    Cl    (1c)
   0.00000000000000    0.00000000000000    0.29400000000000    Cl    (2g)
   0.00000000000000    0.00000000000000   -0.29400000000000    Cl    (2g)
   0.00000000000000    0.00000000000000    0.00000000000000    Hg    (1a)
   0.50000000000000    0.50000000000000    0.50000000000000    NH4   (1d)
```

K_2PtCl_4 ($H1_5$): A4B2C_tP7_123_j_e_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Cl4K2Pt'
_chemical_formula_sum 'Cl4 K2 Pt'

loop_
_publ_author_name
 'R. H. B. Mais'
 'P. G. Owston'
 'A. Wood'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 28
_journal_year 1972
_journal_page_first 393
_journal_page_last 399
_publ_Section_title
;
 The crystal structure of K$_{2}$PtCl$_{4}$ and K$_{2}$PdCl$_{4}$ with
      ↪ estimates of the factors affecting accuracy
;

_aflow_title 'K$_{2}$PtCl$_{4}$ ($H1_{5}$) Structure'
_aflow_proto 'A4B2C_tP7_123_j_e_a'
_aflow_params 'a,c/a,x_{3}'
_aflow_params_values '7.025,0.5898932384344,0.2324'
_aflow_Strukturbericht '$H1_{5}$'
```

```
_aflow_Pearson 'tP7'

_symmetry_space_group_name_H-M "P 4/m 2/m 2/m"
_symmetry_Int_Tables_number 123

_cell_length_a 7.02500
_cell_length_b 7.02500
_cell_length_c 4.14400
_cell_angle_alpha 90.00000
_cell_angle_beta 90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 -y,-x,-z
6 -y,x,z
7 y,-x,z
8 y,x,-z
9 -x,-y,-z
10 -x,y,z
11 x,-y,z
12 x,y,-z
13 y,x,z
14 y,-x,-z
15 -y,x,-z
16 -y,-x,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Pt1 Pt 1 a 0.00000 0.00000 0.00000 1.00000
K1 K 2 e 0.00000 0.50000 0.50000 1.00000
Cl1 Cl 4 j 0.23240 0.23240 0.00000 1.00000
```

K_2PtCl_4 ($H1_5$): A4B2C_tP7_123_j_e_a - POSCAR

```
A4B2C_tP7_123_j_e_a & a,c/a,x3 --params=7.025,0.589893238434,0.2324 & P4
    ↪ /mmm D_{4h}^{1} #123 (aej) & tP7 & $H1_{5}$ & Cl4K2Pt & Cl4K2Pt
    ↪ & R. H. B. Mais and P. G. Owston and A. Wood, Acta
    ↪ Crystallogr. Sect. B Struct. Sci. 28, 393-399 (1972)
   1.00000000000000
   7.02500000000000   0.00000000000000   0.00000000000000
   0.00000000000000   7.02500000000000   0.00000000000000
   0.00000000000000   0.00000000000000   4.14400000000000
    Cl  K  Pt
     4  2   1
Direct
   0.23240000000000   0.23240000000000   0.00000000000000  Cl  (4j)
  -0.23240000000000  -0.23240000000000   0.00000000000000  Cl  (4j)
  -0.23240000000000   0.23240000000000   0.00000000000000  Cl  (4j)
   0.23240000000000  -0.23240000000000   0.00000000000000  Cl  (4j)
   0.00000000000000   0.50000000000000   0.50000000000000  K  (2e)
   0.50000000000000   0.00000000000000   0.50000000000000  K  (2e)
   0.00000000000000   0.00000000000000   0.00000000000000  Pt  (1a)
```

$E6_1$ ($Sr(OH)_2(H_2O)_8$) (*obsolete*): A8B2C_tP11_123_r_f_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Sr(OH)2(H2O)8'
_chemical_formula_sum '(H2O)8 (OH)2 Sr'

loop_
_publ_author_name
 'G. Natta'
_journal_name_full_name
;
 Gazzetta Chimica Italiana
;
_journal_volume 58
_journal_year 1928
_journal_page_first 870
_journal_page_last 872
_publ_Section_title
;
 Constitution of hydroxides and of hydrates. III. Octahydrated strontium
    ↪ hydroxide
;

# Found in Strukturbericht Band II 1928-1932, 1937

_aflow_title '$E6_{1}$ (Sr(OH)$_{2}$(H$_{2}$O)$_{8}$) ({\em{obsolete}})
    ↪ Structure'
_aflow_proto 'A8B2C_tP11_123_r_f_a'
_aflow_params 'a,c/a,x_{3},z_{3}'
_aflow_params_values '6.41,0.905928237129,0.29,0.25'
_aflow_Strukturbericht '$E6_{1}$'
_aflow_Pearson 'tP11'

_symmetry_space_group_name_H-M "P 4/m 2/m 2/m"
_symmetry_Int_Tables_number 123

_cell_length_a 6.41000
_cell_length_b 6.41000
_cell_length_c 5.80700
_cell_angle_alpha 90.00000
_cell_angle_beta 90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 -y,-x,-z
6 -y,x,z
7 y,-x,z
8 y,x,-z
9 -x,-y,-z
10 -x,y,z
11 x,-y,z
12 x,y,-z
13 y,x,z
14 y,-x,-z
15 -y,x,-z
16 -y,-x,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Sr1 Sr 1 a 0.00000 0.00000 0.00000 1.00000
OH1 OH 2 f 0.00000 0.50000 0.00000 1.00000
H2O1 H2O 8 r 0.29000 0.29000 0.25000 1.00000
```

$E6_1$ ($Sr(OH)_2(H_2O)_8$) (*obsolete*): A8B2C_tP11_123_r_f_a - POSCAR

```
A8B2C_tP11_123_r_f_a & a,c/a,x3,z3 --params=6.41,0.905928237129,0.29,
    ↪ 0.25 & P4/mmm D_{4h}^{1} #123 (afr) & tP11 & $E6_{1}$ & Sr(OH)2
    ↪ (H2O)8 & Sr(OH)2(H2O)8 & G. Natta, Gazz. Chim. Ital. 58,
    ↪ 870-872 (1928)
   1.00000000000000
   6.41000000000000   0.00000000000000   0.00000000000000
   0.00000000000000   6.41000000000000   0.00000000000000
   0.00000000000000   0.00000000000000   5.80700000000000
   H2O  OH  Sr
     8   2   1
Direct
   0.29000000000000   0.29000000000000   0.25000000000000  H2O  (8r)
  -0.29000000000000  -0.29000000000000   0.25000000000000  H2O  (8r)
  -0.29000000000000   0.29000000000000   0.25000000000000  H2O  (8r)
   0.29000000000000  -0.29000000000000   0.25000000000000  H2O  (8r)
  -0.29000000000000   0.29000000000000  -0.25000000000000  H2O  (8r)
   0.29000000000000  -0.29000000000000  -0.25000000000000  H2O  (8r)
   0.29000000000000   0.29000000000000  -0.25000000000000  H2O  (8r)
  -0.29000000000000  -0.29000000000000  -0.25000000000000  H2O  (8r)
   0.00000000000000   0.50000000000000   0.00000000000000  OH  (2f)
   0.50000000000000   0.00000000000000   0.00000000000000  OH  (2f)
   0.00000000000000   0.00000000000000   0.00000000000000  Sr  (1a)
```

$E6_2$ [$SrO_2(H_2O)_8$] (*possibly obsolete*): A8B2C_tP11_123_r_h_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral '(H2O)8SrO2'
_chemical_formula_sum '(H2O)8 O2 Sr'

loop_
_publ_author_name
 'G. Natta'
_journal_name_full_name
;
 Gazzetta Chimica Italiana
;
_journal_volume 62
_journal_year 1932
_journal_page_first 444
_journal_page_last 444
_publ_Section_title
;
 ~
;

# Found in The space group of calcium peroxide octahydrate, 1951

_aflow_title '$E6_{2}$ [SrO$_{2}$(H$_{2}$O)$_{8}$] ({\em{possibly
    ↪ obsolete}}) Structure'
_aflow_proto 'A8B2C_tP11_123_r_h_a'
_aflow_params 'a,c/a,z_{2},x_{3},z_{3}'
_aflow_params_values '6.32,0.879746835443,0.1,0.2,0.25'
_aflow_Strukturbericht '$E6_{2}$'
_aflow_Pearson 'tP11'

_symmetry_space_group_name_H-M "P 4/m 2/m 2/m"
_symmetry_Int_Tables_number 123

_cell_length_a 6.32000
_cell_length_b 6.32000
```

```
_cell_length_c     5.56000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 -y,-x,-z
6 -y,x,z
7 y,-x,z
8 y,x,-z
9 -x,-y,-z
10 -x,y,z
11 x,-y,z
12 x,y,-z
13 y,x,z
14 y,-x,-z
15 -y,x,-z
16 -y,-x,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Sr1 Sr 1 a 0.00000 0.00000 0.00000 1.00000
O1 O 2 h 0.50000 0.50000 0.10000 1.00000
H2O1 H2O 8 r 0.20000 0.20000 0.25000 1.00000
```

$E6_2$ [$SrO_2(H_2O)_8$] (*possibly obsolete*): A8B2C_tP11_123_r_h_a - POSCAR

```
A8B2C_tP11_123_r_h_a & a,c/a,z2,x3,z3 --params=6.32,0.879746835443,0.1,
    ↪ 0.2,0.25 & P4/mmm D_{4h}^{1} #123 (ahr) & tP11 & $E6_{2}$ & (
    ↪ H2O)8SrO2 & (H2O)8SrO2 & G. Natta, Gazz. Chim. Ital. 62, 444(
    ↪ 1932)
   1.00000000000000
   6.32000000000000   0.00000000000000   0.00000000000000
   0.00000000000000   6.32000000000000   0.00000000000000
   0.00000000000000   0.00000000000000   5.56000000000000
   H2O     O     Sr
     8     2     1
Direct
   0.20000000000000   0.20000000000000   0.25000000000000  H2O   (8r)
  -0.20000000000000  -0.20000000000000   0.25000000000000  H2O   (8r)
  -0.20000000000000   0.20000000000000   0.25000000000000  H2O   (8r)
   0.20000000000000  -0.20000000000000   0.25000000000000  H2O   (8r)
  -0.20000000000000   0.20000000000000  -0.25000000000000  H2O   (8r)
   0.20000000000000  -0.20000000000000  -0.25000000000000  H2O   (8r)
   0.20000000000000   0.20000000000000  -0.25000000000000  H2O   (8r)
  -0.20000000000000  -0.20000000000000  -0.25000000000000  H2O   (8r)
   0.50000000000000   0.50000000000000   0.10000000000000   O    (2h)
   0.50000000000000   0.50000000000000  -0.10000000000000   O    (2h)
   0.00000000000000   0.00000000000000   0.00000000000000   Sr   (1a)
```

$TlAlF_4$ ($H0_8$): AB4C_tP6_123_d_eh_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'AlF4Tl'
_chemical_formula_sum 'Al F4 Tl'

loop_
_publ_author_name
 'C. Brosset'
_journal_name_full_name
;
 Zeitschrift fur Anorganische und Allgemeine Chemie
;
_journal_volume 235
_journal_year 1937
_journal_page_first 139
_journal_page_last 147
_publ_Section_title
;
 Herstellung und Kristallbau der Verbindungen TlAlF$_{4}$ und Tl$_{2}
    ↪ $AlF$_{5}$
;

# Found in A Structural Classification of Fluoaluminates, 1950

_aflow_title 'TlAlF$_{4}$ ($H0_8$) Structure'
_aflow_proto 'AB4C_tP6_123_d_eh_a'
_aflow_params 'a,c/a,z_{4}'
_aflow_params_values '3.61,1.76454293629,0.215'
_aflow_Strukturbericht '$H0_{8}$'
_aflow_Pearson 'tP6'

_symmetry_space_group_name_H-M "P 4/m 2/m 2/m"
_symmetry_Int_Tables_number 123

_cell_length_a     3.61000
_cell_length_b     3.61000
_cell_length_c     6.37000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 -y,-x,-z
6 -y,x,z
7 y,-x,z
8 y,x,-z
9 -x,-y,-z
10 -x,y,z
11 x,-y,z
12 x,y,-z
13 y,x,z
14 y,-x,-z
15 -y,x,-z
16 -y,-x,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Tl1 Tl 1 a 0.00000 0.00000 0.00000 1.00000
Al1 Al 1 d 0.50000 0.50000 0.50000 1.00000
F1 F 2 e 0.00000 0.50000 0.50000 1.00000
F2 F 2 h 0.50000 0.50000 0.21500 1.00000
```

$TlAlF_4$ ($H0_8$): AB4C_tP6_123_d_eh_a - POSCAR

```
AB4C_tP6_123_d_eh_a & a,c/a,z4 --params=3.61,1.76454293629,0.215 & P4/
    ↪ mmm D_{4h}^{1} #123 (adeh) & tP6 & $H0_{8}$ & AlF4Tl & AlF4Tl &
    ↪ C. Brosset, Z. Anorg. Allg. Chem. 235, 139-147 (1937)
   1.00000000000000
   3.61000000000000   0.00000000000000   0.00000000000000
   0.00000000000000   3.61000000000000   0.00000000000000
   0.00000000000000   0.00000000000000   6.37000000000000
    Al     F     Tl
     1     4     1
Direct
   0.50000000000000   0.50000000000000   0.50000000000000  Al   (1d)
   0.00000000000000   0.50000000000000   0.50000000000000  F    (2e)
   0.50000000000000   0.00000000000000   0.50000000000000  F    (2e)
   0.50000000000000   0.50000000000000   0.21500000000000  F    (2h)
   0.50000000000000   0.50000000000000  -0.21500000000000  F    (2h)
   0.00000000000000   0.00000000000000   0.00000000000000  Tl   (1a)
```

δ-CuTi ($L2_a$): AB_tP2_123_a_d - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'CuTi'
_chemical_formula_sum 'Cu Ti'

loop_
_publ_author_name
 'N. Karlsson'
_journal_name_full_name
;
 Journal of the Institute of Metals
;
_journal_volume 79
_journal_year 1951
_journal_page_first 391
_journal_page_last 405
_publ_Section_title
;
 An X-Ray Study of the Phases in the Copper-Titanium System
;

# Found in A Handbook of Lattice Spacings and Structures of Metals and
    ↪ Alloys, {N.~R.~C. No. 4303}, 1958

_aflow_title '$\delta$-CuTi ($L2_{a}$) Structure'
_aflow_proto 'AB_tP2_123_a_d'
_aflow_params 'a,c/a'
_aflow_params_values '4.44,0.643243243243'
_aflow_Strukturbericht '$L2_{a}$'
_aflow_Pearson 'tP2'

_symmetry_space_group_name_H-M "P 4/m 2/m 2/m"
_symmetry_Int_Tables_number 123

_cell_length_a     4.44000
_cell_length_b     4.44000
_cell_length_c     2.85600
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
```

```
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 -y,-x,-z
6 -y,x,z
7 y,-x,z
8 y,x,-z
9 -x,-y,-z
10 -x,y,z
11 x,-y,z
12 x,y,-z
13 y,x,z
14 y,-x,-z
15 -y,x,-z
16 -y,-x,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cu1 Cu 1 a 0.00000 0.00000 0.00000 1.00000
Ti1 Ti 1 d 0.50000 0.50000 0.50000 1.00000
```

δ-CuTi ($L2_d$): AB_tP2_123_a_d - POSCAR

```
AB_tP2_123_a_d & a,c/a --params=4.44,0.6432432432432 & P4/mmm D_{4h}^{1}
    ↪ #123 (ad) & tP2 & $L2_{a}$ & CuTi & CuTi & N. Karlsson, J.
    ↪ Inst. Met. 79, 391-405 (1951)
   1.00000000000000
   4.44000000000000    0.00000000000000    0.00000000000000
   0.00000000000000    4.44000000000000    0.00000000000000
   0.00000000000000    0.00000000000000    2.85600000000000
    Cu    Ti
     1     1
Direct
   0.00000000000000    0.00000000000000    0.00000000000000    Cu    (1a)
   0.50000000000000    0.50000000000000    0.50000000000000    Ti    (1d)
```

$CaO_2(H_2O)_8$: AB8C2_tP22_124_a_n_h - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Ca(H2O)8O2'
_chemical_formula_sum 'Ca (H2O)8 O2'

loop_
_publ_author_name
 'R. S. Shineman'
 'A. J. King'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 4
_journal_year 1951
_journal_page_first 67
_journal_page_last 68
_publ_Section_title
;
 The space group of calcium peroxide octahydrate
;

_aflow_title 'CaO$_{2}$(H$_{2}$O)$_{8}$ Structure'
_aflow_proto 'AB8C2_tP22_124_a_n_h'
_aflow_params 'a,c/a,z_{2},x_{3},y_{3},z_{3}'
_aflow_params_values '6.21,1.77133655395,0.19,0.3,0.11,0.13'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tP22'

_symmetry_space_group_name_H-M "P 4/m 2/c 2/c"
_symmetry_Int_Tables_number 124

_cell_length_a    6.21000
_cell_length_b    6.21000
_cell_length_c    11.00000
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z+1/2
3 -x,y,-z+1/2
4 -x,-y,z
5 -y,-x,-z+1/2
6 -y,x,z
7 y,-x,z
8 y,x,-z+1/2
9 -x,-y,-z
10 -x,y,z+1/2
11 x,-y,z+1/2
12 x,y,-z
13 y,x,z+1/2
14 y,-x,-z
15 -y,x,-z
16 -y,-x,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ca1 Ca 2 a 0.00000 0.00000 0.25000 1.00000
O1 O 4 h 0.50000 0.50000 0.19000 1.00000
H2O1 H2O 16 n 0.30000 0.11000 0.13000 1.00000
```

$CaO_2(H_2O)_8$: AB8C2_tP22_124_a_n_h - POSCAR

```
AB8C2_tP22_124_a_n_h & a,c/a,z2,x3,y3,z3 --params=6.21,1.77133655395,
    ↪ 0.19,0.3,0.11,0.13 & P4/mcc D_{4h}^{2} #124 (ahn) & tP22 & None
    ↪ & Ca(H2O)8O2 & Ca(H2O)8O2 & R. S. Shineman and A. J. King,
    ↪ Acta Cryst. 4, 67-68 (1951)
   1.00000000000000
   6.21000000000000    0.00000000000000    0.00000000000000
   0.00000000000000    6.21000000000000    0.00000000000000
   0.00000000000000    0.00000000000000   11.00000000000000
    Ca   H2O    O
     2    16    4
Direct
   0.00000000000000    0.00000000000000    0.25000000000000    Ca    (2a)
   0.00000000000000    0.00000000000000    0.75000000000000    Ca    (2a)
   0.30000000000000    0.11000000000000    0.13000000000000    H2O   (16n)
  -0.30000000000000   -0.11000000000000    0.13000000000000    H2O   (16n)
  -0.11000000000000    0.30000000000000    0.13000000000000    H2O   (16n)
   0.11000000000000   -0.30000000000000    0.13000000000000    H2O   (16n)
  -0.30000000000000    0.11000000000000    0.37000000000000    H2O   (16n)
   0.30000000000000   -0.11000000000000    0.37000000000000    H2O   (16n)
   0.11000000000000    0.30000000000000    0.37000000000000    H2O   (16n)
  -0.11000000000000   -0.30000000000000    0.37000000000000    H2O   (16n)
  -0.30000000000000   -0.11000000000000   -0.13000000000000    H2O   (16n)
   0.30000000000000    0.11000000000000   -0.13000000000000    H2O   (16n)
   0.11000000000000   -0.30000000000000   -0.13000000000000    H2O   (16n)
  -0.11000000000000    0.30000000000000   -0.13000000000000    H2O   (16n)
   0.30000000000000   -0.11000000000000    0.63000000000000    H2O   (16n)
  -0.30000000000000    0.11000000000000    0.63000000000000    H2O   (16n)
  -0.11000000000000   -0.30000000000000    0.63000000000000    H2O   (16n)
   0.11000000000000    0.30000000000000    0.63000000000000    H2O   (16n)
   0.50000000000000    0.50000000000000    0.19000000000000    O     (4h)
   0.50000000000000    0.50000000000000    0.31000000000000    O     (4h)
   0.50000000000000    0.50000000000000   -0.19000000000000    O     (4h)
   0.50000000000000    0.50000000000000    0.69000000000000    O     (4h)
```

Vesuvianite ($Ca_{10}Al_4(Mg,Fe)_2Si_9O_{34}(OH)_4$, $S2_3$): A4B10C2D34E4F9_tP252_126_k_ce2k_f_h8k_k_d2k - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Vesuvianite'
_chemical_formula_sum 'Ag4 Ca10 Mg2 O34 (OH)4 Si9'

loop_
_publ_author_name
 'B. E. Warren'
 'D. I. Modell'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 78
_journal_year 1931
_journal_page_first 422
_journal_page_last 432
_publ_Section_title
;
 The Structure of Vesuvianite Ca$_{10}$Al$_{4}$(Mg,Fe)$_{2}$Si$_{9}$O$_{
    ↪ 34}$(OH)$_{4}$
;

# Found in Strukturbericht Band II 1928-1932, 1937

_aflow_title 'Vesuvianite (Ca$_{10}$Al$_{4}$(Mg,Fe)$_{2}$Si$_{9}$O$_{34}
    ↪ $(OH)$_{4}$, $S2_{3}$) Structure'
_aflow_proto 'A4B10C2D34E4F9_tP252_126_k_ce2k_f_h8k_k_d2k'
_aflow_params 'a,c/a,z_{3},x_{5},x_{6},y_{6},z_{6},x_{7},y_{7},z_{7},x_{
    ↪ 8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{10},y_{10},z_{10},x_{11},y_
    ↪ {11},z_{11},x_{12},y_{12},z_{12},x_{13},y_{13},z_{13},x_{14},y_
    ↪ {14},z_{14},x_{15},y_{15},z_{15},x_{16},y_{16},z_{16},x_{17},y_
    ↪ {17},z_{17},x_{18},y_{18},z_{18},x_{19},y_{19},z_{19}'
_aflow_params_values '15.63,0.756877799104,0.13,0.055,0.11,0.89,0.87,
    ↪ 0.81,0.05,0.86,0.17,0.09,0.88,0.83,0.22,0.08,0.87,0.16,0.78,
    ↪ 0.22,-0.06,-0.08,-0.07,0.13,-0.02,0.01,0.83,0.82,0.78,0.87,-
    ↪ 0.07,-0.05,0.82,0.18,-0.08,-0.1,-0.07,0.06,-0.01,0.83,-0.05,
    ↪ 0.19,0.87,-0.09,0.83,0.87'
_aflow_Strukturbericht '$S2_{3}$'
_aflow_Pearson 'tP252'

_symmetry_space_group_name_H-M "P 4/n 2/n 2/c (origin choice 2)"
_symmetry_Int_Tables_number 126

_cell_length_a    15.63000
_cell_length_b    15.63000
_cell_length_c    11.83000
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000
```

```
loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y+1/2,-z+1/2
3 -x+1/2,y,-z+1/2
4 -x+1/2,-y+1/2,z
5 -y+1/2,-x+1/2,-z+1/2
6 -y+1/2,x,z
7 y,-x+1/2,z
8 y,x,-z+1/2
9 -x,-y,-z
10 -x,y+1/2,z+1/2
11 x+1/2,-y,z+1/2
12 x+1/2,y+1/2,-z
13 y+1/2,x+1/2,z+1/2
14 y+1/2,-x,-z
15 -y,x+1/2,-z
16 -y,-x,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ca1 Ca 4 c 0.25000 0.75000 0.75000 1.00000
Si1 Si 4 d 0.25000 0.75000 0.00000 1.00000
Ca2 Ca 4 e 0.25000 0.25000 0.13000 1.00000
Mg1 Mg 8 f 0.00000 0.00000 0.00000 1.00000
O1 O 8 h 0.05500 0.05500 0.25000 1.00000
Ag1 Ag 16 k 0.11000 0.89000 0.87000 1.00000
Ca3 Ca 16 k 0.81000 0.05000 0.86000 1.00000
Ca4 Ca 16 k 0.17000 0.09000 0.88000 1.00000
O2 O 16 k 0.83000 0.22000 0.08000 1.00000
O3 O 16 k 0.87000 0.16000 0.78000 1.00000
O4 O 16 k 0.22000 -0.06000 -0.08000 1.00000
O5 O 16 k -0.07000 0.13000 -0.02000 1.00000
O6 O 16 k 0.01000 0.83000 0.82000 1.00000
O7 O 16 k 0.78000 0.87000 -0.07000 1.00000
O8 O 16 k -0.05000 0.82000 0.18000 1.00000
O9 O 16 k -0.08000 -0.10000 -0.07000 1.00000
OH1 OH 16 k 0.06000 -0.01000 0.83000 1.00000
Si2 Si 16 k -0.05000 0.19000 0.87000 1.00000
Si3 Si 16 k -0.09000 0.83000 0.87000 1.00000
```

Vesuvianite ($Ca_{10}Al_4(Mg,Fe)_2Si_9O_{34}(OH)_4$, $S2_3$): A4B10C2D34E4F9_tP252_126_k_ce2k_f_h8k_k_d2k - POSCAR

```
A4B10C2D34E4F9_tP252_126_k_ce2k_f_h8k_k_d2k & a,c/a,z3,x5,x6,y6,z6,x7,y7
    ↪ ,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10,x11,y11,z11,x12,y12,z12,x13,
    ↪ y13,z13,x14,y14,z14,x15,y15,z15,x16,y16,z16,x17,y17,z17,x18,y18
    ↪ ,z18,x19,y19,z19 --params=15.63,0.756877799104,0.13,0.055,0.11,
    ↪ 0.89,0.87,0.81,0.05,0.86,0.17,0.09,0.88,0.83,0.22,0.08,0.87,
    ↪ 0.16,0.78,0.22,-0.06,-0.08,-0.07,0.13,-0.02,0.01,0.83,0.82,0.78
    ↪ ,0.87,-0.07,-0.05,0.82,0.18,-0.08,-0.1,-0.07,0.06,-0.01,0.83,-
    ↪ 0.05,0.19,0.87,-0.09,0.83,0.87 & P4/nnc D_{4h}^{4} #126 (cdefhk
    ↪ ^14) & tP252 & $S2_{3}$ & Al4Ca10(Mg,Fe)2O34(OH)4Si9 &
    ↪ Vesuvianite & B. E. Warren and D. I. Modell, Zeitschrift f"{u}r
    ↪ Kristallographie - Crystalline Materials 78, 422-432 (1931)
   1.00000000000000
  15.63000000000000   0.00000000000000   0.00000000000000
   0.00000000000000  15.63000000000000   0.00000000000000
   0.00000000000000   0.00000000000000  11.83000000000000
    Ag   Ca   Mg    O   OH   Si
    16   40    8  136   16   36
Direct
   0.11000000000000   0.89000000000000   0.87000000000000  Ag  (16k)
   0.39000000000000  -0.39000000000000   0.87000000000000  Ag  (16k)
  -0.39000000000000   0.11000000000000   0.87000000000000  Ag  (16k)
   0.89000000000000   0.39000000000000   0.87000000000000  Ag  (16k)
   0.39000000000000   0.89000000000000  -0.37000000000000  Ag  (16k)
   0.11000000000000  -0.39000000000000  -0.37000000000000  Ag  (16k)
   0.89000000000000   0.11000000000000  -0.37000000000000  Ag  (16k)
  -0.39000000000000   0.39000000000000  -0.37000000000000  Ag  (16k)
  -0.11000000000000  -0.89000000000000  -0.87000000000000  Ag  (16k)
   0.61000000000000   1.39000000000000  -0.87000000000000  Ag  (16k)
   1.39000000000000  -0.11000000000000  -0.87000000000000  Ag  (16k)
  -0.89000000000000   0.61000000000000  -0.87000000000000  Ag  (16k)
   0.61000000000000  -0.89000000000000   1.37000000000000  Ag  (16k)
  -0.11000000000000   1.39000000000000   1.37000000000000  Ag  (16k)
  -0.89000000000000  -0.11000000000000   1.37000000000000  Ag  (16k)
   1.39000000000000   0.61000000000000   1.37000000000000  Ag  (16k)
   0.25000000000000   0.75000000000000   0.75000000000000  Ca   (4c)
   0.75000000000000   0.25000000000000   0.75000000000000  Ca   (4c)
   0.75000000000000   0.25000000000000   0.25000000000000  Ca   (4c)
   0.25000000000000   0.75000000000000   0.25000000000000  Ca   (4c)
   0.25000000000000   0.25000000000000   0.13000000000000  Ca   (4e)
   0.25000000000000   0.25000000000000   0.37000000000000  Ca   (4e)
   0.75000000000000   0.75000000000000  -0.13000000000000  Ca   (4e)
   0.75000000000000   0.75000000000000   0.63000000000000  Ca   (4e)
   0.81000000000000   0.05000000000000   0.86000000000000  Ca  (16k)
  -0.31000000000000   0.45000000000000   0.86000000000000  Ca  (16k)
   0.45000000000000   0.81000000000000   0.86000000000000  Ca  (16k)
   0.05000000000000  -0.31000000000000   0.86000000000000  Ca  (16k)
  -0.31000000000000   0.05000000000000  -0.36000000000000  Ca  (16k)
   0.81000000000000   0.45000000000000  -0.36000000000000  Ca  (16k)
   0.05000000000000   0.81000000000000  -0.36000000000000  Ca  (16k)
   0.45000000000000  -0.31000000000000  -0.36000000000000  Ca  (16k)
  -0.81000000000000  -0.05000000000000  -0.86000000000000  Ca  (16k)
   1.31000000000000   0.55000000000000  -0.86000000000000  Ca  (16k)
   0.55000000000000  -0.81000000000000  -0.86000000000000  Ca  (16k)
  -0.05000000000000   1.31000000000000  -0.86000000000000  Ca  (16k)
   1.31000000000000  -0.05000000000000   1.36000000000000  Ca  (16k)
  -0.81000000000000   0.55000000000000   1.36000000000000  Ca  (16k)
  -0.05000000000000  -0.81000000000000   1.36000000000000  Ca  (16k)
   0.55000000000000   1.31000000000000   1.36000000000000  Ca  (16k)
   0.17000000000000   0.09000000000000   0.88000000000000  Ca  (16k)
   0.33000000000000   0.41000000000000   0.88000000000000  Ca  (16k)
   0.41000000000000   0.17000000000000   0.88000000000000  Ca  (16k)
   0.09000000000000   0.33000000000000   0.88000000000000  Ca  (16k)
   0.33000000000000   0.09000000000000  -0.38000000000000  Ca  (16k)
   0.17000000000000   0.41000000000000  -0.38000000000000  Ca  (16k)
   0.09000000000000   0.17000000000000  -0.38000000000000  Ca  (16k)
   0.41000000000000   0.33000000000000  -0.38000000000000  Ca  (16k)
  -0.17000000000000  -0.09000000000000  -0.88000000000000  Ca  (16k)
   0.67000000000000   0.59000000000000  -0.88000000000000  Ca  (16k)
   0.59000000000000  -0.17000000000000  -0.88000000000000  Ca  (16k)
  -0.09000000000000   0.67000000000000  -0.88000000000000  Ca  (16k)
   0.67000000000000  -0.09000000000000   1.38000000000000  Ca  (16k)
  -0.17000000000000   0.59000000000000   1.38000000000000  Ca  (16k)
  -0.09000000000000  -0.17000000000000   1.38000000000000  Ca  (16k)
   0.59000000000000   0.67000000000000   1.38000000000000  Ca  (16k)
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   0.50000000000000   0.50000000000000   0.00000000000000  Mg   (8f)
   0.50000000000000   0.00000000000000   0.00000000000000  Mg   (8f)
   0.00000000000000   0.50000000000000   0.00000000000000  Mg   (8f)
   0.50000000000000   0.00000000000000   0.50000000000000  Mg   (8f)
   0.00000000000000   0.50000000000000   0.50000000000000  Mg   (8f)
   0.00000000000000   0.00000000000000   0.50000000000000  Mg   (8f)
   0.50000000000000   0.50000000000000   0.50000000000000  Mg   (8f)
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   0.44500000000000   0.44500000000000   0.25000000000000  O   (8h)
   0.44500000000000   0.05500000000000   0.25000000000000  O   (8h)
   0.05500000000000   0.44500000000000   0.25000000000000  O   (8h)
  -0.05500000000000  -0.05500000000000   0.75000000000000  O   (8h)
   0.55500000000000   0.55500000000000   0.75000000000000  O   (8h)
   0.55500000000000  -0.05500000000000   0.75000000000000  O   (8h)
  -0.05500000000000   0.55500000000000   0.75000000000000  O   (8h)
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  -0.33000000000000   0.28000000000000   0.08000000000000  O  (16k)
   0.28000000000000   0.83000000000000   0.08000000000000  O  (16k)
   0.22000000000000  -0.33000000000000   0.08000000000000  O  (16k)
  -0.33000000000000   0.22000000000000   0.42000000000000  O  (16k)
   0.83000000000000   0.28000000000000   0.42000000000000  O  (16k)
   0.22000000000000   0.83000000000000   0.42000000000000  O  (16k)
   0.28000000000000  -0.33000000000000   0.42000000000000  O  (16k)
  -0.83000000000000  -0.22000000000000  -0.08000000000000  O  (16k)
   1.33000000000000   0.72000000000000  -0.08000000000000  O  (16k)
   0.72000000000000  -0.83000000000000  -0.08000000000000  O  (16k)
  -0.22000000000000   1.33000000000000  -0.08000000000000  O  (16k)
   1.33000000000000  -0.22000000000000   0.58000000000000  O  (16k)
  -0.83000000000000   0.72000000000000   0.58000000000000  O  (16k)
  -0.22000000000000  -0.83000000000000   0.58000000000000  O  (16k)
   0.72000000000000   1.33000000000000   0.58000000000000  O  (16k)
   0.87000000000000   0.16000000000000   0.78000000000000  O  (16k)
  -0.37000000000000   0.34000000000000   0.78000000000000  O  (16k)
   0.34000000000000   0.87000000000000   0.78000000000000  O  (16k)
   0.16000000000000  -0.37000000000000   0.78000000000000  O  (16k)
  -0.37000000000000   0.16000000000000  -0.28000000000000  O  (16k)
   0.87000000000000   0.34000000000000  -0.28000000000000  O  (16k)
   0.16000000000000   0.87000000000000  -0.28000000000000  O  (16k)
   0.34000000000000  -0.37000000000000  -0.28000000000000  O  (16k)
  -0.87000000000000  -0.16000000000000  -0.78000000000000  O  (16k)
   1.37000000000000   0.66000000000000  -0.78000000000000  O  (16k)
   0.66000000000000  -0.87000000000000  -0.78000000000000  O  (16k)
  -0.16000000000000   1.37000000000000  -0.78000000000000  O  (16k)
   1.37000000000000  -0.16000000000000   1.28000000000000  O  (16k)
  -0.87000000000000   0.66000000000000   1.28000000000000  O  (16k)
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   0.66000000000000   1.37000000000000   1.28000000000000  O  (16k)
   0.22000000000000  -0.06000000000000  -0.08000000000000  O  (16k)
   0.28000000000000   0.56000000000000  -0.08000000000000  O  (16k)
   0.56000000000000   0.22000000000000  -0.08000000000000  O  (16k)
  -0.06000000000000   0.28000000000000  -0.08000000000000  O  (16k)
   0.28000000000000  -0.06000000000000   0.58000000000000  O  (16k)
   0.22000000000000   0.56000000000000   0.58000000000000  O  (16k)
  -0.06000000000000   0.22000000000000   0.58000000000000  O  (16k)
   0.56000000000000   0.28000000000000   0.58000000000000  O  (16k)
  -0.22000000000000   0.06000000000000   0.08000000000000  O  (16k)
   0.72000000000000   0.44000000000000   0.08000000000000  O  (16k)
   0.44000000000000  -0.22000000000000   0.08000000000000  O  (16k)
   0.06000000000000   0.72000000000000   0.08000000000000  O  (16k)
   0.72000000000000   0.06000000000000   0.42000000000000  O  (16k)
  -0.22000000000000   0.44000000000000   0.42000000000000  O  (16k)
   0.06000000000000  -0.22000000000000   0.42000000000000  O  (16k)
   0.44000000000000   0.72000000000000   0.42000000000000  O  (16k)
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   0.57000000000000   0.37000000000000  -0.02000000000000  O  (16k)
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   0.57000000000000   0.13000000000000   0.52000000000000  O  (16k)
  -0.07000000000000   0.37000000000000   0.52000000000000  O  (16k)
   0.13000000000000  -0.07000000000000   0.52000000000000  O  (16k)
   0.37000000000000   0.57000000000000   0.52000000000000  O  (16k)
   0.07000000000000  -0.13000000000000   0.02000000000000  O  (16k)
   0.43000000000000   0.63000000000000   0.02000000000000  O  (16k)
   0.63000000000000   0.07000000000000   0.02000000000000  O  (16k)
  -0.13000000000000   0.43000000000000   0.02000000000000  O  (16k)
   0.43000000000000  -0.13000000000000   0.48000000000000  O  (16k)
   0.07000000000000   0.63000000000000   0.48000000000000  O  (16k)
  -0.13000000000000   0.07000000000000   0.48000000000000  O  (16k)
   0.63000000000000   0.43000000000000   0.48000000000000  O  (16k)
   0.01000000000000   0.83000000000000   0.82000000000000  O  (16k)
   0.49000000000000  -0.33000000000000   0.82000000000000  O  (16k)
  -0.33000000000000   0.01000000000000   0.82000000000000  O  (16k)
```

```
 0.83000000000000   0.49000000000000   0.82000000000000   O   (16k)
 0.49000000000000   0.83000000000000  -0.32000000000000   O   (16k)
 0.01000000000000  -0.33000000000000  -0.32000000000000   O   (16k)
 0.83000000000000   0.01000000000000  -0.32000000000000   O   (16k)
-0.33000000000000   0.49000000000000  -0.32000000000000   O   (16k)
-0.01000000000000  -0.83000000000000  -0.82000000000000   O   (16k)
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 1.33000000000000  -0.01000000000000  -0.82000000000000   O   (16k)
-0.83000000000000   0.51000000000000  -0.82000000000000   O   (16k)
 0.51000000000000  -0.83000000000000   1.32000000000000   O   (16k)
-0.01000000000000   1.33000000000000   1.32000000000000   O   (16k)
-0.83000000000000  -0.01000000000000   1.32000000000000   O   (16k)
 1.33000000000000   0.51000000000000   1.32000000000000   O   (16k)
 0.78000000000000   0.87000000000000  -0.07000000000000   O   (16k)
-0.28000000000000  -0.37000000000000  -0.07000000000000   O   (16k)
-0.37000000000000   0.78000000000000  -0.07000000000000   O   (16k)
 0.87000000000000  -0.28000000000000  -0.07000000000000   O   (16k)
-0.28000000000000   0.87000000000000   0.57000000000000   O   (16k)
 0.78000000000000  -0.37000000000000   0.57000000000000   O   (16k)
 0.87000000000000   0.78000000000000   0.57000000000000   O   (16k)
-0.37000000000000  -0.28000000000000   0.57000000000000   O   (16k)
-0.78000000000000  -0.87000000000000   0.07000000000000   O   (16k)
 1.28000000000000   1.37000000000000   0.07000000000000   O   (16k)
 1.37000000000000  -0.78000000000000   0.07000000000000   O   (16k)
-0.87000000000000   1.28000000000000   0.07000000000000   O   (16k)
 1.28000000000000  -0.87000000000000   0.43000000000000   O   (16k)
-0.78000000000000   1.37000000000000   0.43000000000000   O   (16k)
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-0.01000000000000   0.44000000000000   0.83000000000000   OH  (16k)
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 0.01000000000000   0.56000000000000  -0.83000000000000   OH  (16k)
 0.56000000000000   0.01000000000000   1.33000000000000   OH  (16k)
-0.06000000000000   0.49000000000000   1.33000000000000   OH  (16k)
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 0.49000000000000   0.56000000000000   1.33000000000000   OH  (16k)
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 0.75000000000000   0.25000000000000   0.00000000000000   Si   (4d)
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-0.09000000000000   0.83000000000000   0.87000000000000   Si  (16k)
 0.59000000000000  -0.33000000000000   0.87000000000000   Si  (16k)
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 0.09000000000000   1.33000000000000   1.37000000000000   Si  (16k)
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 1.33000000000000   0.41000000000000   1.37000000000000   Si  (16k)
```

$Ag[Co(NH_3)_2(NO_2)_4]$ ($J1_9$): ABC4D2E8_tP32_126_a_b_h_e_k - CIF

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loop_
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 'A. F. Wells'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 95
_journal_year 1936
_journal_page_first 74
_journal_page_last 82
_publ_Section_title
;
 The Crystal Structure of Silver Diamminotetranitro-cobaltiate Ag[Co(
    ↪ NH$_{3}$)$_{2}$(NO$_{2}$)$_{4}$]
;

# Found in Strukturbericht Band IV 1936, 1938

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_symmetry_space_group_name_H-M "P 4/n 2/n 2/c (origin choice 2)"
_symmetry_Int_Tables_number 126

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_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
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2 x,-y+1/2,-z+1/2
3 -x+1/2,y,-z+1/2
4 -x+1/2,-y+1/2,z
5 -y+1/2,-x+1/2,-z+1/2
6 -y+1/2,x,z
7 y,-x+1/2,z
8 y,x,-z+1/2
9 -x,-y,-z
10 -x,y+1/2,z+1/2
11 x+1/2,-y,z+1/2
12 x+1/2,y+1/2,-z
13 y+1/2,x+1/2,z+1/2
14 y+1/2,-x,-z
15 -y,x+1/2,-z
16 -y,-x,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ag1 Ag 2 a 0.25000 0.25000 0.25000 1.00000
Co1 Co 2 b 0.25000 0.25000 0.75000 1.00000
NH31 NH3 4 e 0.25000 0.25000 0.53000 1.00000
N1 N 8 h -0.05500 -0.05500 0.25000 1.00000
O1 O 16 k 0.06000 -0.03000 0.34000 1.00000
```

$Ag[Co(NH_3)_2(NO_2)_4]$ ($J1_9$): ABC4D2E8_tP32_126_a_b_h_e_k - POSCAR

```
ABC4D2E8_tP32_126_a_b_h_e_k & a,c/a,z3,x4,x5,y5,z5 --params=6.67,
    ↪ 1.56371814093,0.53,-0.055,0.06,-0.03,0.34 & P4/nnc D_{4h}^{4} #
    ↪ 126 (abehk) & tP32 & $J1_{9}$ & AgCoN4(NH3)2O8 & AgCoN4(NH3)2O8
    ↪ & A. F. Wells, Zeitschrift f{"u}r Kristallographie -
    ↪ Crystalline Materials 95, 74-82 (1936)
   1.00000000000000
   6.67000000000000    0.00000000000000    0.00000000000000
   0.00000000000000    6.67000000000000    0.00000000000000
   0.00000000000000    0.00000000000000   10.43000000000000
    Ag     Co     N    NH3     O
     2      2      8      4     16
Direct
   0.25000000000000    0.25000000000000    0.25000000000000    Ag    (2a)
```

```
  0.75000000000000   0.75000000000000   0.75000000000000   Ag   (2a)
  0.25000000000000   0.25000000000000   0.75000000000000   Co   (2b)
  0.75000000000000   0.75000000000000   0.25000000000000   Co   (2b)
 -0.05550000000000  -0.05550000000000   0.25000000000000   N    (8h)
  0.55550000000000   0.55550000000000   0.25000000000000   N    (8h)
  0.55550000000000  -0.05550000000000   0.25000000000000   N    (8h)
 -0.05550000000000   0.55550000000000   0.25000000000000   N    (8h)
  0.05550000000000   0.05550000000000   0.75000000000000   N    (8h)
  0.44450000000000   0.44450000000000   0.75000000000000   N    (8h)
  0.44450000000000   0.05550000000000   0.75000000000000   N    (8h)
  0.05550000000000   0.44450000000000   0.75000000000000   N    (8h)
  0.25000000000000   0.25000000000000   0.53000000000000   NH3  (4e)
  0.25000000000000   0.25000000000000  -0.03000000000000   NH3  (4e)
  0.75000000000000   0.75000000000000  -0.53000000000000   NH3  (4e)
  0.75000000000000   0.75000000000000   1.03000000000000   NH3  (4e)
  0.06000000000000  -0.03000000000000   0.34000000000000   O    (16k)
  0.44000000000000   0.53000000000000   0.34000000000000   O    (16k)
  0.53000000000000   0.06000000000000   0.34000000000000   O    (16k)
 -0.03000000000000   0.44000000000000   0.34000000000000   O    (16k)
  0.44000000000000  -0.03000000000000   0.16000000000000   O    (16k)
  0.06000000000000   0.53000000000000   0.16000000000000   O    (16k)
 -0.03000000000000   0.06000000000000   0.16000000000000   O    (16k)
  0.53000000000000   0.44000000000000   0.16000000000000   O    (16k)
 -0.06000000000000   0.03000000000000  -0.34000000000000   O    (16k)
  0.56000000000000   0.47000000000000  -0.34000000000000   O    (16k)
  0.47000000000000  -0.06000000000000  -0.34000000000000   O    (16k)
  0.03000000000000   0.56000000000000  -0.34000000000000   O    (16k)
  0.56000000000000   0.03000000000000   0.84000000000000   O    (16k)
 -0.06000000000000   0.47000000000000   0.84000000000000   O    (16k)
  0.03000000000000  -0.06000000000000   0.84000000000000   O    (16k)
  0.47000000000000   0.56000000000000   0.84000000000000   O    (16k)
```

$Pd(NH_3)_4Cl_2{\cdot}H_2O$ ($H4_9$): A2BC4D_tP16_127_h_d_i_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Cl2(H2O)(NH3)4Pd'
_chemical_formula_sum 'Cl2 (H2O) (NH3)4 Pd'

loop_
_publ_author_name
 'B. N. Dickinson'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 88
_journal_year 1934
_journal_page_first 281
_journal_page_last 297
_publ_Section_title
;
 The Crystal Structure of Tetramminopalladous Chloride Pd(NH$_{3}$)$_{4}
     ↪ $Cl$_{2}$$\cdot$H$_{2}$O
;

_aflow_title 'Pd(NH$_{3}$)$_{4}$Cl$_{2}$$\cdot$H$_{2}$O ($H4_{9}$)
     ↪ Structure'
_aflow_proto 'A2BC4D_tP16_127_h_d_i_a'
_aflow_params 'a,c/a,x_{3},x_{4},y_{4}'
_aflow_params_values '10.302,0.42127742186,0.285,0.194,0.027'
_aflow_Strukturbericht '$H4_{9}$'
_aflow_Pearson 'tP16'

_symmetry_space_group_name_H-M "P 4/m 21/b 2/m"
_symmetry_Int_Tables_number 127

_cell_length_a    10.30200
_cell_length_b    10.30200
_cell_length_c    4.34000
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z
3 -x+1/2,y+1/2,-z
4 -x,-y,z
5 -y+1/2,-x+1/2,-z
6 -y,x,z
7 y,-x,z
8 y+1/2,x+1/2,-z
9 -x,-y,-z
10 -x+1/2,y+1/2,z
11 x+1/2,-y+1/2,z
12 x,y,-z
13 y+1/2,x+1/2,z
14 y,-x,-z
15 -y,x,-z
16 -y+1/2,-x+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Pd1 Pd 2 a 0.00000 0.00000 0.00000 1.00000
H2O1 H2O 2 d 0.00000 0.50000 0.00000 1.00000
Cl1 Cl 4 h 0.28500 0.78500 0.50000 1.00000
NH31 NH3 8 i 0.19400 0.02700 0.00000 1.00000
```

$Pd(NH_3)_4Cl_2{\cdot}H_2O$ ($H4_9$): A2BC4D_tP16_127_h_d_i_a - POSCAR

```
A2BC4D_tP16_127_h_d_i_a & a,c/a,x3,x4,y4 --params=10.302,0.42127742186,
     ↪ 0.285,0.194,0.027 & P4/mbm D_{4h}^{5} #127 (adhi) & tP16 & $H4_
     ↪ {9}$ & Cl2(H2O)(NH3)4Pd & Cl2(H2O)(NH3)4Pd & B. N. Dickinson,
     ↪ Zeitschrift f{"u}r Kristallographie - Crystalline Materials 88,
     ↪ 281-297 (1934)
   1.00000000000000
  10.30200000000000   0.00000000000000   0.00000000000000
   0.00000000000000  10.30200000000000   0.00000000000000
   0.00000000000000   0.00000000000000   4.34000000000000
    Cl   H2O   NH3    Pd
     4     2     8     2
Direct
   0.28500000000000   0.78500000000000   0.50000000000000   Cl   (4h)
  -0.28500000000000   0.21500000000000   0.50000000000000   Cl   (4h)
   0.21500000000000   0.28500000000000   0.50000000000000   Cl   (4h)
   0.78500000000000  -0.28500000000000   0.50000000000000   Cl   (4h)
   0.00000000000000   0.50000000000000   0.00000000000000   H2O  (2d)
   0.50000000000000   0.00000000000000   0.00000000000000   H2O  (2d)
   0.19400000000000   0.02700000000000   0.00000000000000   NH3  (8i)
  -0.19400000000000  -0.02700000000000   0.00000000000000   NH3  (8i)
  -0.02700000000000   0.19400000000000   0.00000000000000   NH3  (8i)
   0.02700000000000  -0.19400000000000   0.00000000000000   NH3  (8i)
   0.30600000000000   0.52700000000000   0.00000000000000   NH3  (8i)
   0.69400000000000   0.47300000000000   0.00000000000000   NH3  (8i)
   0.52700000000000   0.69400000000000   0.00000000000000   NH3  (8i)
   0.47300000000000   0.30600000000000   0.00000000000000   NH3  (8i)
   0.00000000000000   0.00000000000000   0.00000000000000   Pd   (2a)
   0.50000000000000   0.50000000000000   0.00000000000000   Pd   (2a)
```

Phosgenite $[Pb_2Cl_2(CO_3)]$: AB2C3D2_tP32_127_g_eh_gk_k - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Phosgenite'
_chemical_formula_sum 'C Cl2 O3 Pb2'

loop_
_publ_author_name
 'G. Giuseppetti'
 'C. Tadini'
_journal_name_full_name
;
 Tschermaks mineralogische und petrographische Mitteilungen
;
_journal_volume 21
_journal_year 1974
_journal_page_first 101
_journal_page_last 109
_publ_Section_title
;
 Reexamination of the crystal structure of phosgenite, Pb$_{2}$Cl$_{2}$(
     ↪ CO$_{3}$)
;

_aflow_title 'Phosgenite [Pb$_{2}$Cl$_{2}$(CO$_{3}$)] Structure'
_aflow_proto 'AB2C3D2_tP32_127_g_eh_gk_k'
_aflow_params 'a,c/a,z_{1},x_{2},x_{3},x_{4},x_{5},z_{5},x_{6},z_{6}'
_aflow_params_values '8.16,1.08860294118,0.7572,0.3257,0.211,0.3521,
     ↪ 0.3726,0.1269,0.1659,0.2594'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tP32'

_symmetry_space_group_name_H-M "P 4/m 21/b 2/m"
_symmetry_Int_Tables_number 127

_cell_length_a    8.16000
_cell_length_b    8.16000
_cell_length_c    8.88300
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z
3 -x+1/2,y+1/2,-z
4 -x,-y,z
5 -y+1/2,-x+1/2,-z
6 -y,x,z
7 y,-x,z
8 y+1/2,x+1/2,-z
9 -x,-y,-z
10 -x+1/2,y+1/2,z
11 x+1/2,-y+1/2,z
12 x,y,-z
13 y+1/2,x+1/2,z
14 y,-x,-z
15 -y,x,-z
16 -y+1/2,-x+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
```

```
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cl1 Cl 4 e 0.00000 0.00000 0.75720 1.00000
C1 C 4 g 0.32570 0.82570 0.00000 1.00000
O1 O 4 g 0.21100 0.71100 0.00000 1.00000
Cl2 Cl 4 h 0.35210 0.85210 0.50000 1.00000
O2 O 8 k 0.37260 0.87260 0.12690 1.00000
Pb1 Pb 8 k 0.16590 0.66590 0.25940 1.00000
```

Phosgenite [$Pb_2Cl_2(CO_3)$]: AB2C3D2_tP32_127_g_eh_gk_k - POSCAR

```
AB2C3D2_tP32_127_g_eh_gk_k & a,c/a,z1,x2,x3,x4,x5,z5,x6,z6 --params=8.16
    ↪ ,1.08860294118,0.7572,0.3257,0.211,0.3521,0.3726,0.1269,0.1659,
    ↪ 0.2594 & P4/mbm D_{4h}^{5} #127 (eg^2hk^2) & tP32 & None &
    ↪ CCl2O3Pb2 & Phosgenite & G. Giuseppetti and C. Tadini,
    ↪ Tschermaks Min. Petr. Mitt. 21, 101-109 (1974)
   1.00000000000000
   8.16000000000000   0.00000000000000   0.00000000000000
   0.00000000000000   8.16000000000000   0.00000000000000
   0.00000000000000   0.00000000000000   8.88300000000000
     C     Cl     O     Pb
     4      8     12      8
Direct
   0.32570000000000   0.82570000000000   0.00000000000000   C    (4g)
  -0.32570000000000   0.17430000000000   0.00000000000000   C    (4g)
   0.17430000000000   0.32570000000000   0.00000000000000   C    (4g)
   0.82570000000000  -0.32570000000000   0.00000000000000   C    (4g)
   0.00000000000000   0.00000000000000   0.75720000000000   Cl   (4e)
   0.50000000000000   0.50000000000000  -0.75720000000000   Cl   (4e)
   0.00000000000000   0.00000000000000  -0.75720000000000   Cl   (4e)
   0.50000000000000   0.50000000000000   0.75720000000000   Cl   (4e)
   0.35210000000000   0.85210000000000   0.50000000000000   Cl   (4h)
  -0.35210000000000   0.14790000000000   0.50000000000000   Cl   (4h)
   0.14790000000000   0.35210000000000   0.50000000000000   Cl   (4h)
   0.85210000000000  -0.35210000000000   0.50000000000000   Cl   (4h)
   0.21100000000000   0.71100000000000   0.00000000000000   O    (4g)
  -0.21100000000000   0.28900000000000   0.00000000000000   O    (4g)
   0.28900000000000   0.21100000000000   0.00000000000000   O    (4g)
   0.71100000000000  -0.21100000000000   0.00000000000000   O    (4g)
   0.37260000000000   0.87260000000000   0.12690000000000   O    (8k)
  -0.37260000000000   0.12740000000000   0.12690000000000   O    (8k)
   0.12740000000000   0.37260000000000   0.12690000000000   O    (8k)
   0.87260000000000  -0.37260000000000   0.12690000000000   O    (8k)
   0.12740000000000   0.37260000000000  -0.12690000000000   O    (8k)
   0.87260000000000  -0.37260000000000  -0.12690000000000   O    (8k)
   0.37260000000000   0.87260000000000  -0.12690000000000   O    (8k)
  -0.37260000000000   0.12740000000000  -0.12690000000000   O    (8k)
   0.16590000000000   0.66590000000000   0.25940000000000   Pb   (8k)
  -0.16590000000000   0.33410000000000   0.25940000000000   Pb   (8k)
   0.33410000000000   0.16590000000000   0.25940000000000   Pb   (8k)
   0.66590000000000  -0.16590000000000   0.25940000000000   Pb   (8k)
   0.33410000000000   0.16590000000000  -0.25940000000000   Pb   (8k)
   0.66590000000000  -0.16590000000000  -0.25940000000000   Pb   (8k)
   0.16590000000000   0.66590000000000  -0.25940000000000   Pb   (8k)
  -0.16590000000000   0.33410000000000  -0.25940000000000   Pb   (8k)
```

Chiolite ($Na_5Al_3F_{14}$, $K7_5$): A3B14C5_tP44_128_ac_ehi_bg - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Chiolite'
_chemical_formula_sum 'Al3 F14 Na5'

loop_
_publ_author_name
 'C. Jacoboni'
 'A. Leble'
 'J. J. Rousseau'
_journal_name_full_name
;
 Journal of Solid State Chemistry
;
_journal_volume 36
_journal_year 1981
_journal_page_first 297
_journal_page_last 304
_publ_Section_title
;
 D\\'{e}termination pr\\'{e}cise de la structure de la chiolite Na$_{5}
    ↪ $Al$_{3}$F$_{14}$ et \\'{e}tude par R.P.E. de Na$_{5}$Al$_{3}
    ↪ $F$_{14}$:Cr$^{3+}$
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'Chiolite (Na$_{5}$Al$_{3}$F$_{14}$, $K7_{5}$) Structure'
_aflow_proto 'A3B14C5_tP44_128_ac_ehi_bg'
_aflow_params 'a,c/a,z_{4},x_{5},x_{6},y_{6},x_{7},y_{7},z_{7}'
_aflow_params_values '7.0138,1.48307622116,0.1711,0.2768,0.0642,0.2477,
    ↪ 0.1794,0.5364,0.1198'
_aflow_Strukturbericht '$K7_{5}$'
_aflow_Pearson 'tP44'

_symmetry_space_group_name_H-M "P 4/m 21/n 2/c"
_symmetry_Int_Tables_number 128

_cell_length_a    7.01380
_cell_length_b    7.01380
_cell_length_c    10.40200
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x+1/2,y+1/2,-z+1/2
4 -x,-y,z
5 -y+1/2,-x+1/2,-z+1/2
6 -y,x,z
7 y,-x,z
8 y+1/2,x+1/2,-z+1/2
9 -x,-y,-z
10 -x+1/2,y+1/2,z+1/2
11 x+1/2,-y+1/2,z+1/2
12 x,y,-z
13 y+1/2,x+1/2,z+1/2
14 y,-x,-z
15 -y,x,-z
16 -y+1/2,-x+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Al1 Al 2 a 0.00000 0.00000 0.00000 1.00000
Na1 Na 2 b 0.00000 0.00000 0.50000 1.00000
Al2 Al 4 c 0.00000 0.50000 0.00000 1.00000
F1 F 4 e 0.00000 0.00000 0.17110 1.00000
Na2 Na 8 g 0.27680 0.77680 0.25000 1.00000
F2 F 8 h 0.06420 0.24770 0.00000 1.00000
F3 F 16 i 0.17940 0.53640 0.11980 1.00000
```

Chiolite ($Na_5Al_3F_{14}$, $K7_5$): A3B14C5_tP44_128_ac_ehi_bg - POSCAR

```
A3B14C5_tP44_128_ac_ehi_bg & a,c/a,z4,x5,x6,y6,x7,y7,z7 --params=7.0138,
    ↪ 1.48307622116,0.1711,0.2768,0.0642,0.2477,0.1794,0.5364,0.1198
    ↪ & P4/mnc D_{4h}^{6} #128 (abceghi) & tP44 & $K7_{5}$ &
    ↪ Al3F14Na5 & Chiolite & C. Jacoboni and A. Leble and J. J.
    ↪ Rousseau, J. Solid State Chem. 36, 297-304 (1981)
   1.00000000000000
   7.01380000000000   0.00000000000000   0.00000000000000
   0.00000000000000   7.01380000000000   0.00000000000000
   0.00000000000000   0.00000000000000  10.40200000000000
    Al     F     Na
     6    28     10
Direct
   0.00000000000000   0.00000000000000   0.00000000000000   Al   (2a)
   0.50000000000000   0.50000000000000   0.50000000000000   Al   (2a)
   0.00000000000000   0.50000000000000   0.00000000000000   Al   (4c)
   0.50000000000000   0.00000000000000   0.00000000000000   Al   (4c)
   0.50000000000000   0.00000000000000   0.50000000000000   Al   (4c)
   0.00000000000000   0.50000000000000   0.50000000000000   Al   (4c)
   0.00000000000000   0.00000000000000   0.17110000000000   F    (4e)
   0.50000000000000   0.50000000000000   0.32890000000000   F    (4e)
   0.00000000000000   0.00000000000000  -0.17110000000000   F    (4e)
   0.50000000000000   0.50000000000000   0.67110000000000   F    (4e)
   0.06420000000000   0.24770000000000   0.00000000000000   F    (8h)
  -0.06420000000000  -0.24770000000000   0.00000000000000   F    (8h)
  -0.24770000000000   0.06420000000000   0.00000000000000   F    (8h)
   0.24770000000000  -0.06420000000000   0.00000000000000   F    (8h)
   0.43580000000000   0.74770000000000   0.50000000000000   F    (8h)
   0.56420000000000   0.25230000000000   0.50000000000000   F    (8h)
   0.74770000000000   0.56420000000000   0.50000000000000   F    (8h)
   0.25230000000000   0.43580000000000   0.50000000000000   F    (8h)
   0.17940000000000   0.53640000000000   0.11980000000000   F   (16i)
  -0.17940000000000  -0.53640000000000   0.11980000000000   F   (16i)
  -0.53640000000000   0.17940000000000   0.11980000000000   F   (16i)
   0.53640000000000  -0.17940000000000   0.11980000000000   F   (16i)
   0.32060000000000   1.03640000000000   0.38020000000000   F   (16i)
   0.67940000000000  -0.03640000000000   0.38020000000000   F   (16i)
   1.03640000000000   0.67940000000000   0.38020000000000   F   (16i)
  -0.03640000000000   0.32060000000000   0.38020000000000   F   (16i)
  -0.17940000000000  -0.53640000000000  -0.11980000000000   F   (16i)
   0.17940000000000   0.53640000000000  -0.11980000000000   F   (16i)
   0.53640000000000  -0.17940000000000  -0.11980000000000   F   (16i)
  -0.53640000000000   0.17940000000000  -0.11980000000000   F   (16i)
   0.67940000000000  -0.03640000000000   0.61980000000000   F   (16i)
   0.32060000000000   1.03640000000000   0.61980000000000   F   (16i)
  -0.03640000000000   0.32060000000000   0.61980000000000   F   (16i)
   1.03640000000000   0.67940000000000   0.61980000000000   F   (16i)
   0.00000000000000   0.00000000000000   0.50000000000000   Na   (2b)
   0.50000000000000   0.50000000000000   0.00000000000000   Na   (2b)
   0.27680000000000   0.77680000000000   0.25000000000000   Na   (8g)
  -0.27680000000000   0.22320000000000   0.25000000000000   Na   (8g)
   0.22320000000000   0.27680000000000   0.25000000000000   Na   (8g)
   0.77680000000000  -0.27680000000000   0.25000000000000   Na   (8g)
  -0.27680000000000   0.22320000000000   0.75000000000000   Na   (8g)
   0.27680000000000   0.77680000000000   0.75000000000000   Na   (8g)
   0.77680000000000  -0.27680000000000   0.75000000000000   Na   (8g)
   0.22320000000000   0.27680000000000   0.75000000000000   Na   (8g)
```

Apophyllite ($KCa_4Si_8O_{20}F{\cdot}8H_2O$, $S5_2$): A4BC16DE28F8_tP116_128_h_a_2i_b_g3i_i - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM
```

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_chemical_name_mineral 'Apophyllite'
_chemical_formula_sum 'Ca4 F H16 K O28 Si8'

loop_
_publ_author_name
 'G. Y. Chao'
_journal_name_full_name
;
 American Mineralogist
;
_journal_volume 56
_journal_year 1971
_journal_page_first 1234
_journal_page_last 1242
_publ_Section_title
;
 The refinement of the crystal structure of apophyllite: II.
      ↪ Determination of the hydrogen positions by X-ray diffraction
;

_aflow_title 'Apophyllite (KCa$_{4}$Si$_{8}$O$_{20}$F$\cdot$8H$_{2}$O,
      ↪ $S5_{2}$) Structure'
_aflow_proto 'A4BC16DE28F8_tP116_128_h_a_2i_b_g3i_i'
_aflow_params 'a,c/a,x_{3},x_{4},y_{4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{
      ↪ 6},x_{7},y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{10
      ↪ },y_{10},z_{10}'
_aflow_params_values '8.965,1.75872838818,0.1369,0.1094,0.2466,0.4515,
      ↪ 0.177,0.0775,0.2362,0.4706,0.1198,0.0846,0.1891,0.2178,0.2636,
      ↪ 0.1026,0.0923,0.2131,0.4491,0.0898,0.2256,0.0865,0.19'
_aflow_Strukturbericht '$S5_{2}$'
_aflow_Pearson 'tP116'

_symmetry_space_group_name_H-M "P 4/m 21/n 2/c"
_symmetry_Int_Tables_number 128

_cell_length_a      8.96500
_cell_length_b      8.96500
_cell_length_c      15.76700
_cell_angle_alpha   90.00000
_cell_angle_beta    90.00000
_cell_angle_gamma   90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x+1/2,y+1/2,-z+1/2
4 -x,-y,z
5 -y+1/2,-x+1/2,-z+1/2
6 -y,x,z
7 y,-x,z
8 y+1/2,x+1/2,-z+1/2
9 -x,-y,-z
10 -x+1/2,y+1/2,z+1/2
11 x+1/2,-y+1/2,z+1/2
12 x,y,-z
13 y+1/2,x+1/2,z+1/2
14 y,-x,-z
15 -y,x,-z
16 -y+1/2,-x+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
F1 F 2 a 0.00000 0.00000 0.00000 1.00000
K1 K 2 b 0.00000 0.00000 0.50000 1.00000
O1 O 8 g 0.13690 0.63690 0.25000 1.00000
Ca1 Ca 8 h 0.10940 0.24660 0.00000 1.00000
H1 H 16 i 0.45150 0.17700 0.07750 1.00000
H2 H 16 i 0.23620 0.47060 0.11980 1.00000
O2 O 16 i 0.08460 0.18910 0.21780 1.00000
O3 O 16 i 0.26360 0.10260 0.09230 1.00000
O4 O 16 i 0.21310 0.44910 0.08980 1.00000
Si1 Si 16 i 0.22560 0.08650 0.19000 1.00000
```

Apophyllite ($KCa_4Si_8O_{20}F{\cdot}8H_2O$, $S5_2$): A4BC16DE28F8_tP116_128_h_a_2i_b_g3i_i - POSCAR

```
A4BC16DE28F8_tP116_128_h_a_2i_b_g3i_i & a,c/a,x3,x4,y4,x5,y5,z5,x6,y6,z6
      ↪ ,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10 --params=8.965,
      ↪ 1.75872838818,0.1369,0.1094,0.2466,0.4515,0.177,0.0775,0.2362,
      ↪ 0.4706,0.1198,0.0846,0.1891,0.2178,0.2636,0.1026,0.0923,0.2131,
      ↪ 0.4491,0.0898,0.2256,0.0865,0.19 & P4/mnc D_{4h}^{6} #128 (
      ↪ abghi^6) & tP116 & $S5_{2}$ & Ca4FH16KO28Si8 & Apophyllite & G.
      ↪ Y. Chao, Am. Mineral. 56, 1234-1242 (1971)
   1.00000000000000
     8.96500000000000    0.00000000000000    0.00000000000000
     0.00000000000000    8.96500000000000    0.00000000000000
     0.00000000000000    0.00000000000000   15.76700000000000
      Ca      F      H      K      O     Si
       8      2     32      2     56     16
Direct
    0.10940000000000    0.24660000000000    0.00000000000000   Ca    (8h)
   -0.10940000000000   -0.24660000000000    0.00000000000000   Ca    (8h)
   -0.24660000000000    0.10940000000000    0.00000000000000   Ca    (8h)
    0.24660000000000   -0.10940000000000    0.00000000000000   Ca    (8h)
    0.39060000000000    0.74660000000000    0.50000000000000   Ca    (8h)
    0.60940000000000    0.25340000000000    0.50000000000000   Ca    (8h)
    0.74660000000000    0.60940000000000    0.50000000000000   Ca    (8h)
    0.25340000000000    0.39060000000000    0.50000000000000   Ca    (8h)
    0.00000000000000    0.00000000000000    0.00000000000000   F     (2a)
    0.50000000000000    0.50000000000000    0.50000000000000   F     (2a)
    0.45150000000000    0.17700000000000    0.07750000000000   H    (16i)
   -0.45150000000000   -0.17700000000000    0.07750000000000   H    (16i)
   -0.17700000000000    0.45150000000000    0.07750000000000   H    (16i)
    0.17700000000000   -0.45150000000000    0.07750000000000   H    (16i)
    0.04850000000000    0.67700000000000    0.42250000000000   H    (16i)
    0.95150000000000    0.32300000000000    0.42250000000000   H    (16i)
    0.67700000000000    0.95150000000000    0.42250000000000   H    (16i)
    0.32300000000000    0.04850000000000    0.42250000000000   H    (16i)
   -0.45150000000000   -0.17700000000000   -0.07750000000000   H    (16i)
    0.45150000000000    0.17700000000000   -0.07750000000000   H    (16i)
    0.17700000000000   -0.45150000000000   -0.07750000000000   H    (16i)
   -0.17700000000000    0.45150000000000   -0.07750000000000   H    (16i)
    0.95150000000000    0.32300000000000    0.57750000000000   H    (16i)
    0.04850000000000    0.67700000000000    0.57750000000000   H    (16i)
    0.32300000000000    0.04850000000000    0.57750000000000   H    (16i)
    0.67700000000000    0.95150000000000    0.57750000000000   H    (16i)
    0.23620000000000    0.47060000000000    0.11980000000000   H    (16i)
   -0.23620000000000   -0.47060000000000    0.11980000000000   H    (16i)
   -0.47060000000000    0.23620000000000    0.11980000000000   H    (16i)
    0.47060000000000   -0.23620000000000    0.11980000000000   H    (16i)
    0.26380000000000    0.97060000000000    0.38020000000000   H    (16i)
    0.73620000000000    0.02940000000000    0.38020000000000   H    (16i)
    0.97060000000000    0.73620000000000    0.38020000000000   H    (16i)
    0.02940000000000    0.26380000000000    0.38020000000000   H    (16i)
   -0.23620000000000   -0.47060000000000   -0.11980000000000   H    (16i)
    0.23620000000000    0.47060000000000   -0.11980000000000   H    (16i)
    0.47060000000000   -0.23620000000000   -0.11980000000000   H    (16i)
   -0.47060000000000    0.23620000000000   -0.11980000000000   H    (16i)
    0.73620000000000    0.02940000000000    0.61980000000000   H    (16i)
    0.26380000000000    0.97060000000000    0.61980000000000   H    (16i)
    0.02940000000000    0.26380000000000    0.61980000000000   H    (16i)
    0.97060000000000    0.73620000000000    0.61980000000000   H    (16i)
    0.00000000000000    0.00000000000000    0.50000000000000   K     (2b)
    0.50000000000000    0.50000000000000    0.00000000000000   K     (2b)
    0.13690000000000    0.63690000000000    0.25000000000000   O     (8g)
   -0.13690000000000    0.36310000000000    0.25000000000000   O     (8g)
    0.36310000000000    0.13690000000000    0.25000000000000   O     (8g)
    0.63690000000000   -0.13690000000000    0.25000000000000   O     (8g)
   -0.13690000000000    0.36310000000000    0.75000000000000   O     (8g)
    0.13690000000000    0.63690000000000    0.75000000000000   O     (8g)
    0.63690000000000   -0.13690000000000    0.75000000000000   O     (8g)
    0.36310000000000    0.13690000000000    0.75000000000000   O     (8g)
    0.08460000000000    0.18910000000000    0.21780000000000   O    (16i)
   -0.08460000000000   -0.18910000000000    0.21780000000000   O    (16i)
   -0.18910000000000    0.08460000000000    0.21780000000000   O    (16i)
    0.18910000000000   -0.08460000000000    0.21780000000000   O    (16i)
    0.41540000000000    0.68910000000000    0.28220000000000   O    (16i)
    0.58460000000000    0.31090000000000    0.28220000000000   O    (16i)
    0.68910000000000    0.58460000000000    0.28220000000000   O    (16i)
    0.31090000000000    0.41540000000000    0.28220000000000   O    (16i)
   -0.08460000000000   -0.18910000000000   -0.21780000000000   O    (16i)
    0.08460000000000    0.18910000000000   -0.21780000000000   O    (16i)
    0.18910000000000   -0.08460000000000   -0.21780000000000   O    (16i)
   -0.18910000000000    0.08460000000000   -0.21780000000000   O    (16i)
    0.58460000000000    0.31090000000000    0.71780000000000   O    (16i)
    0.41540000000000    0.68910000000000    0.71780000000000   O    (16i)
    0.31090000000000    0.41540000000000    0.71780000000000   O    (16i)
    0.68910000000000    0.58460000000000    0.71780000000000   O    (16i)
    0.26360000000000    0.10260000000000    0.09230000000000   O    (16i)
   -0.26360000000000   -0.10260000000000    0.09230000000000   O    (16i)
   -0.10260000000000    0.26360000000000    0.09230000000000   O    (16i)
    0.10260000000000   -0.26360000000000    0.09230000000000   O    (16i)
    0.23640000000000    0.60260000000000    0.40770000000000   O    (16i)
    0.76360000000000    0.39740000000000    0.40770000000000   O    (16i)
    0.60260000000000    0.76360000000000    0.40770000000000   O    (16i)
    0.39740000000000    0.23640000000000    0.40770000000000   O    (16i)
   -0.26360000000000   -0.10260000000000   -0.09230000000000   O    (16i)
    0.26360000000000    0.10260000000000   -0.09230000000000   O    (16i)
    0.10260000000000   -0.26360000000000   -0.09230000000000   O    (16i)
   -0.10260000000000    0.26360000000000   -0.09230000000000   O    (16i)
    0.76360000000000    0.39740000000000    0.59230000000000   O    (16i)
    0.23640000000000    0.60260000000000    0.59230000000000   O    (16i)
    0.39740000000000    0.23640000000000    0.59230000000000   O    (16i)
    0.60260000000000    0.76360000000000    0.59230000000000   O    (16i)
    0.21310000000000    0.44910000000000    0.08980000000000   O    (16i)
   -0.21310000000000   -0.44910000000000    0.08980000000000   O    (16i)
   -0.44910000000000    0.21310000000000    0.08980000000000   O    (16i)
    0.44910000000000   -0.21310000000000    0.08980000000000   O    (16i)
    0.28690000000000    0.94910000000000    0.41020000000000   O    (16i)
    0.71310000000000    0.05090000000000    0.41020000000000   O    (16i)
    0.94910000000000    0.71310000000000    0.41020000000000   O    (16i)
    0.05090000000000    0.28690000000000    0.41020000000000   O    (16i)
   -0.21310000000000   -0.44910000000000   -0.08980000000000   O    (16i)
    0.21310000000000    0.44910000000000   -0.08980000000000   O    (16i)
    0.44910000000000   -0.21310000000000   -0.08980000000000   O    (16i)
   -0.44910000000000    0.21310000000000   -0.08980000000000   O    (16i)
    0.71310000000000    0.05090000000000    0.58980000000000   O    (16i)
    0.28690000000000    0.94910000000000    0.58980000000000   O    (16i)
    0.05090000000000    0.28690000000000    0.58980000000000   O    (16i)
    0.94910000000000    0.71310000000000    0.58980000000000   O    (16i)
    0.22560000000000    0.08650000000000    0.19000000000000   Si   (16i)
   -0.22560000000000   -0.08650000000000    0.19000000000000   Si   (16i)
   -0.08650000000000    0.22560000000000    0.19000000000000   Si   (16i)
    0.08650000000000   -0.22560000000000    0.19000000000000   Si   (16i)
    0.27440000000000    0.58650000000000    0.31000000000000   Si   (16i)
    0.72560000000000    0.41350000000000    0.31000000000000   Si   (16i)
    0.58650000000000    0.72560000000000    0.31000000000000   Si   (16i)
    0.41350000000000    0.27440000000000    0.31000000000000   Si   (16i)
   -0.22560000000000   -0.08650000000000   -0.19000000000000   Si   (16i)
    0.22560000000000    0.08650000000000   -0.19000000000000   Si   (16i)
    0.08650000000000   -0.22560000000000   -0.19000000000000   Si   (16i)
   -0.08650000000000    0.22560000000000   -0.19000000000000   Si   (16i)
```

```
    0.72560000000000    0.41350000000000    0.69000000000000    Si    (16i)
    0.27440000000000    0.58650000000000    0.69000000000000    Si    (16i)
    0.41350000000000    0.27440000000000    0.69000000000000    Si    (16i)
    0.58650000000000    0.72560000000000    0.69000000000000    Si    (16i)
```

$CaBe_2Ge_2$: A2BC2_tP10_129_ac_c_bc - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Be2CaGe2'
_chemical_formula_sum 'Be2 Ca Ge2'

loop_
_publ_author_name
 'B. Eisenmann'
 'N. May'
 'W. M\"{u}ller'
 'H. Sch\"{a}fer'
_journal_year 1972
_publ_Section_title
;
 Eine neue strukturelle Variante des BaAl$_{4}$-Typs: Der CaBe$_{2}$Ge$_
    ↪ {2}$-Typ
;

_aflow_title 'CaBe$_{2}$Ge$_{2}$ Structure'
_aflow_proto 'A2BC2_tP10_129_ac_c_bc'
_aflow_params 'a,c/a,z_{3},z_{4},z_{5}'
_aflow_params_values '4.02,2.46766169154,0.608,0.249,0.868'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tP10'

_symmetry_space_group_name_H-M "P 4/n 21/m 2/m (origin choice 2)"
_symmetry_Int_Tables_number 129

_cell_length_a    4.02000
_cell_length_b    4.02000
_cell_length_c    9.92000
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y,-z
3 -x,y+1/2,-z
4 -x+1/2,-y+1/2,z
5 -y,-x,-z
6 -y+1/2,x,z
7 y,-x+1/2,z
8 y+1/2,x+1/2,-z
9 -x,-y,-z
10 -x+1/2,y,z
11 x,-y+1/2,z
12 x+1/2,y+1/2,-z
13 y,x,z
14 y+1/2,-x,-z
15 -y,x+1/2,-z
16 -y+1/2,-x+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Be1 Be 2 a 0.75000 0.25000 0.00000 1.00000
Ge1 Ge 2 b 0.75000 0.25000 0.50000 1.00000
Be2 Be 2 c 0.25000 0.25000 0.60800 1.00000
Ca1 Ca 2 c 0.25000 0.25000 0.24900 1.00000
Ge2 Ge 2 c 0.25000 0.25000 0.86800 1.00000
```

$CaBe_2Ge_2$: A2BC2_tP10_129_ac_c_bc - POSCAR

```
A2BC2_tP10_129_ac_c_bc & a,c/a,z3,z4,z5 --params=4.02,2.46766169154,
    ↪ 0.608,0.249,0.868 & P4/nmm D_{4h}^{7} #129 (abc^3) & tP10 &
    ↪ None & Be2CaGe2 & Be2CaGe2 & B. Eisenmann et al., (1972)
   1.00000000000000
   4.02000000000000    0.00000000000000    0.00000000000000
   0.00000000000000    4.02000000000000    0.00000000000000
   0.00000000000000    0.00000000000000    9.92000000000000
     Be     Ca     Ge
      4      2      4
Direct
   0.75000000000000    0.25000000000000    0.00000000000000    Be    (2a)
   0.25000000000000    0.75000000000000    0.00000000000000    Be    (2a)
   0.25000000000000    0.25000000000000    0.60800000000000    Be    (2c)
   0.75000000000000    0.75000000000000   -0.60800000000000    Be    (2c)
   0.25000000000000    0.25000000000000    0.24900000000000    Ca    (2c)
   0.75000000000000    0.75000000000000   -0.24900000000000    Ca    (2c)
   0.75000000000000    0.25000000000000    0.50000000000000    Ge    (2b)
   0.25000000000000    0.75000000000000    0.50000000000000    Ge    (2b)
   0.25000000000000    0.25000000000000    0.86800000000000    Ge    (2c)
   0.75000000000000    0.75000000000000   -0.86800000000000    Ge    (2c)
```

Meta-autunite (I) $[Ca(UO_2)_2(PO_4)_2{\cdot}6H_2O, H5_{10}]$: AB4C6DE_tP26_129_c_j_2ci_a_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Meta-autunite (i)'
_chemical_formula_sum 'Ca (H2O)4 O6 P U'

loop_
_publ_author_name
 'Y. S. Makarov'
 'V. I. Ivanov'
_journal_name_full_name
;
 Doklady Akademii Nauk SSSR
;
_journal_volume 132
_journal_year 1960
_journal_page_first 601
_journal_page_last 603
_publ_Section_title
;
 The crystal structure of meta-autunite, Ca(UO$_{2}$)$_{2}$(PO$_{4}$)$_{
    ↪ 2}$*6H$_{2}$O
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'Meta-autunite (I) [Ca(UO$_{2}$)$_{2}$(PO$_{4}$)$_{2}$$\
    ↪ cdot$6H$_{2}$O, $H5_{10}$] Structure'
_aflow_proto 'AB4C6DE_tP26_129_c_j_2ci_a_c'
_aflow_params 'a,c/a,z_{2},z_{3},z_{4},z_{5},y_{6},z_{6},x_{7},z_{7}'
_aflow_params_values '6.98,1.20630372493,0.612,0.343,0.893,0.106,0.584,
    ↪ 0.106,0.486,0.392'
_aflow_Strukturbericht '$H5_{10}$'
_aflow_Pearson 'tP26'

_symmetry_space_group_name_H-M "P 4/n 21/m 2/m (origin choice 2)"
_symmetry_Int_Tables_number 129

_cell_length_a    6.98000
_cell_length_b    6.98000
_cell_length_c    8.42000
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y,-z
3 -x,y+1/2,-z
4 -x+1/2,-y+1/2,z
5 -y,-x,-z
6 -y+1/2,x,z
7 y,-x+1/2,z
8 y+1/2,x+1/2,-z
9 -x,-y,-z
10 -x+1/2,y,z
11 x,-y+1/2,z
12 x+1/2,y+1/2,-z
13 y,x,z
14 y+1/2,-x,-z
15 -y,x+1/2,-z
16 -y+1/2,-x+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
P1 P 2 a 0.75000 0.25000 0.00000 1.00000
Ca1 Ca 2 c 0.25000 0.25000 0.61200 0.50000
O1 O 2 c 0.25000 0.25000 0.34300 1.00000
O2 O 2 c 0.25000 0.25000 0.89300 1.00000
U1 U 2 c 0.25000 0.25000 0.10600 1.00000
O3 O 8 i 0.25000 0.58400 0.10600 1.00000
H2O1 H2O 8 j 0.48600 0.48600 0.39200 0.75000
```

Meta-autunite (I) $[Ca(UO_2)_2(PO_4)_2{\cdot}6H_2O, H5_{10}]$: AB4C6DE_tP26_129_c_j_2ci_a_c - POSCAR

```
AB4C6DE_tP26_129_c_j_2ci_a_c & a,c/a,z2,z3,z4,z5,y6,z6,x7,z7 --params=
    ↪ 6.98,1.20630372493,0.612,0.343,0.893,0.106,0.584,0.106,0.486,
    ↪ 0.392 & P4/nmm D_{4h}^{7} #129 (ac^4ij) & tP26 & $H5_{10}$ &
    ↪ CaH12O18P2U2 & Meta-autunite (i) & Y. S. Makarov and V. I.
    ↪ Ivanov, Doklady Akademii Nauk SSSR 132, 601-603 (1960)
   1.00000000000000
   6.98000000000000    0.00000000000000    0.00000000000000
   0.00000000000000    6.98000000000000    0.00000000000000
   0.00000000000000    0.00000000000000    8.42000000000000
     Ca    H2O     O     P     U
      2      8     12     2     2
Direct
   0.25000000000000    0.25000000000000    0.61200000000000    Ca    (2c)
   0.75000000000000    0.75000000000000   -0.61200000000000    Ca    (2c)
   0.48600000000000    0.48600000000000    0.39200000000000    H2O   (8j)
   0.01400000000000    0.01400000000000    0.39200000000000    H2O   (8j)
   0.01400000000000    0.48600000000000    0.39200000000000    H2O   (8j)
   0.48600000000000    0.01400000000000    0.39200000000000    H2O   (8j)
  -0.48600000000000    0.98600000000000   -0.39200000000000    H2O   (8j)
   0.98600000000000   -0.48600000000000   -0.39200000000000    H2O   (8j)
```

```
   0.98600000000000   0.98600000000000  -0.39200000000000  H2O   ( 8j )
  -0.48600000000000  -0.48600000000000  -0.39200000000000  H2O   ( 8j )
   0.25000000000000   0.25000000000000   0.34300000000000  O     ( 2c )
   0.75000000000000   0.75000000000000  -0.34300000000000  O     ( 2c )
   0.25000000000000   0.25000000000000   0.89300000000000  O     ( 2c )
   0.75000000000000   0.75000000000000  -0.89300000000000  O     ( 2c )
   0.25000000000000   0.58400000000000   0.10600000000000  O     ( 8i )
   0.25000000000000  -0.08400000000000   0.10600000000000  O     ( 8i )
  -0.08400000000000   0.25000000000000   0.10600000000000  O     ( 8i )
   0.58400000000000   0.25000000000000   0.10600000000000  O     ( 8i )
   0.75000000000000   1.08400000000000  -0.10600000000000  O     ( 8i )
   0.75000000000000  -0.58400000000000  -0.10600000000000  O     ( 8i )
   1.08400000000000   0.75000000000000  -0.10600000000000  O     ( 8i )
  -0.58400000000000   0.75000000000000  -0.10600000000000  O     ( 8i )
   0.75000000000000   0.25000000000000   0.00000000000000  P     ( 2a )
   0.25000000000000   0.75000000000000   0.00000000000000  P     ( 2a )
   0.25000000000000   0.25000000000000   0.10600000000000  U     ( 2c )
   0.75000000000000   0.75000000000000  -0.10600000000000  U     ( 2c )
```

NH_4Br ($B25$): AB4C_tP12_129_c_i_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'BrH4N'
_chemical_formula_sum 'Br H4 N'

loop_
_publ_author_name
 'H. A. Levy'
 'S. W. Peterson'
_journal_name_full_name
;
 Journal of the American Chemical Society
;
_journal_volume 75
_journal_year 1953
_journal_page_first 1536
_journal_page_last 1542
_publ_Section_title
;
 Neutron Diffraction Determination of the Crystal Structure of Ammonium
      ↪ Bromide in Four Phases
;

_aflow_title 'NH$_{4}$Br ($B25$) Structure'
_aflow_proto 'AB4C_tP12_129_c_i_a'
_aflow_params 'a,c/a,z_{2},y_{3},z_{3}'
_aflow_params_values '5.82,0.710652920962,0.48,0.897,0.147'
_aflow_Strukturbericht '$B25$'
_aflow_Pearson 'tP12'

_symmetry_space_group_name_H-M "P 4/n 21/m 2/m (origin choice 2)"
_symmetry_Int_Tables_number 129

_cell_length_a    5.82000
_cell_length_b    5.82000
_cell_length_c    4.13600
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y,-z
3 -x,y+1/2,-z
4 -x+1/2,-y+1/2,z
5 -y,-x,-z
6 -y+1/2,x,z
7 y,-x+1/2,z
8 y+1/2,x+1/2,-z
9 -x,-y,-z
10 -x+1/2,y,z
11 x,-y+1/2,z
12 x+1/2,y+1/2,-z
13 y,x,z
14 y+1/2,-x,-z
15 -y,x+1/2,-z
16 -y+1/2,-x+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
N1 N 2 a 0.25000 0.75000 0.00000 1.00000
Br1 Br 2 c 0.25000 0.25000 0.48000 1.00000
H1 H 8 i 0.25000 0.89700 0.14700 1.00000
```

NH_4Br ($B25$): AB4C_tP12_129_c_i_a - POSCAR

```
AB4C_tP12_129_c_i_a & a,c/a,z2,y3,z3 --params=5.82,0.710652920962,0.48,
      ↪ 0.897,0.147 & P4/nmm D_{4h}^{7} #129 (aci) & tP12 & $B25$ &
      ↪ BrH4N & BrH4N & H. A. Levy and S. W. Peterson, J. Am. Chem.
      ↪ Soc. 75, 1536-1542 (1953)
   1.00000000000000
   5.82000000000000   0.00000000000000   0.00000000000000
   0.00000000000000   5.82000000000000   0.00000000000000
   0.00000000000000   0.00000000000000   4.13600000000000
    Br    H    N
     2    8    2
Direct
   0.25000000000000   0.25000000000000   0.48000000000000  Br    ( 2c )
   0.75000000000000   0.75000000000000  -0.48000000000000  Br    ( 2c )
   0.25000000000000   0.89700000000000   0.14700000000000  H     ( 8i )
   0.25000000000000  -0.39700000000000   0.14700000000000  H     ( 8i )
  -0.39700000000000   0.25000000000000   0.14700000000000  H     ( 8i )
   0.89700000000000   0.25000000000000   0.14700000000000  H     ( 8i )
   0.75000000000000   1.39700000000000  -0.14700000000000  H     ( 8i )
   0.75000000000000  -0.89700000000000  -0.14700000000000  H     ( 8i )
   1.39700000000000   0.75000000000000  -0.14700000000000  H     ( 8i )
  -0.89700000000000   0.75000000000000  -0.14700000000000  H     ( 8i )
   0.75000000000000   0.25000000000000   0.00000000000000  N     ( 2a )
   0.25000000000000   0.75000000000000   0.00000000000000  N     ( 2a )
```

LaOAgS: ABCD_tP8_129_b_c_a_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'AgLaOS'
_chemical_formula_sum 'Ag La O S'

loop_
_publ_author_name
 'M. Palazzi'
 'S. Jaulmes'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 37
_journal_year 1981
_journal_page_first 1337
_journal_page_last 1339
_publ_Section_title
;
 Structure du Conducteur Ionique (LaO)AgS
;

# Found in Pnictides and Chalcogenides II, 2003

_aflow_title 'LaOAgS Structure'
_aflow_proto 'ABCD_tP8_129_b_c_a_c'
_aflow_params 'a,c/a,z_{3},z_{4}'
_aflow_params_values '4.05,2.23185185185,0.1356,0.6929'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tP8'

_symmetry_space_group_name_H-M "P 4/n 21/m 2/m (origin choice 2)"
_symmetry_Int_Tables_number 129

_cell_length_a    4.05000
_cell_length_b    4.05000
_cell_length_c    9.03900
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y,-z
3 -x,y+1/2,-z
4 -x+1/2,-y+1/2,z
5 -y,-x,-z
6 -y+1/2,x,z
7 y,-x+1/2,z
8 y+1/2,x+1/2,-z
9 -x,-y,-z
10 -x+1/2,y,z
11 x,-y+1/2,z
12 x+1/2,y+1/2,-z
13 y,x,z
14 y+1/2,-x,-z
15 -y,x+1/2,-z
16 -y+1/2,-x+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 2 a 0.75000 0.25000 0.00000 1.00000
Ag1 Ag 2 b 0.75000 0.25000 0.50000 1.00000
La1 La 2 c 0.25000 0.25000 0.13560 1.00000
S1 S 2 c 0.25000 0.25000 0.69290 1.00000
```

LaOAgS: ABCD_tP8_129_b_c_a_c - POSCAR

```
ABCD_tP8_129_b_c_a_c & a,c/a,z3,z4 --params=4.05,2.23185185185,0.1356,
      ↪ 0.6929 & P4/nmm D_{4h}^{7} #129 (abc^2) & tP8 & None & AgLaOS &
      ↪ AgLaOS & M. Palazzi and S. Jaulmes, Acta Crystallogr. Sect. B
      ↪ Struct. Sci. 37, 1337-1339 (1981)
   1.00000000000000
   4.05000000000000   0.00000000000000   0.00000000000000
   0.00000000000000   4.05000000000000   0.00000000000000
```

```
    0.00000000000000    0.00000000000000    9.03900000000000
     Ag      La      O       S
      2       2       2       2
Direct
    0.75000000000000    0.25000000000000    0.50000000000000    Ag    (2b)
    0.25000000000000    0.75000000000000    0.50000000000000    Ag    (2b)
    0.25000000000000    0.25000000000000    0.13560000000000    La    (2c)
    0.75000000000000    0.75000000000000   -0.13560000000000    La    (2c)
    0.75000000000000    0.25000000000000    0.00000000000000    O     (2a)
    0.25000000000000    0.75000000000000    0.00000000000000    O     (2a)
    0.25000000000000    0.25000000000000    0.69290000000000    S     (2c)
    0.75000000000000    0.75000000000000   -0.69290000000000    S     (2c)
```

$Sr(OH)_2(H_2O)_8$: A18B10C_tP116_130_2c4g_2c2g_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'H18O10Sr'
_chemical_formula_sum 'H18 O10 Sr'

loop_
_publ_author_name
 'J. S. Ricci'
 'R. C. Stevens'
 'R. K. {McMullan}'
 'W. T. Klooster'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 61
_journal_year 2005
_journal_page_first 381
_journal_page_last 386
_publ_Section_title
;
 Structure of strontium hydroxide octahydrate, Sr(OH)$_{2} \cdot $8H$_{2
    ↪ }$O, at 20, 100 and 200 K from neutron diffraction
;

_aflow_title 'Sr(OH)$_{2}$(H$_{2}$O)$_{8}$ Structure'
_aflow_proto 'A18B10C_tP116_130_2c4g_2c2g_a'
_aflow_params 'a,c/a,z_{2},z_{3},z_{4},z_{5},x_{6},y_{6},z_{6},x_{7},y_{
    ↪ 7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{10},y_{10},z_{
    ↪ 10},x_{11},y_{11},z_{11}'
_aflow_params_values '8.984,1.28194568121,0.4888,0.2426,0.40482,0.15904,
    ↪ 0.8044,0.5759,0.1201,-0.047,0.4874,0.1327,0.8054,0.5766,0.8623,
    ↪ 0.8392,0.4596,-0.0398,0.84543,0.47219,0.12549,0.84587,0.47366,
    ↪ 0.87655'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tP116'

_symmetry_space_group_name_H-M "P 4/n 21/c 2/c (origin choice 2)"
_symmetry_Int_Tables_number 130

_cell_length_a     8.98400
_cell_length_b     8.98400
_cell_length_c     11.51700
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y,-z+1/2
3 -x,y+1/2,-z+1/2
4 -x+1/2,-y+1/2,z
5 -y,-x,-z+1/2
6 -y+1/2,x,z
7 y,-x+1/2,z
8 y+1/2,x+1/2,-z+1/2
9 -x,-y,-z
10 -x+1/2,y,z+1/2
11 x,-y+1/2,z+1/2
12 x+1/2,y+1/2,-z
13 y,x,z+1/2
14 y+1/2,-x,-z
15 -y,x+1/2,-z
16 -y+1/2,-x+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Sr1 Sr 4 a 0.75000 0.25000 0.25000 1.00000
H1 H 4 c 0.25000 0.25000 0.48880 1.00000
H2 H 4 c 0.25000 0.25000 0.24260 1.00000
O1 O 4 c 0.25000 0.25000 0.40482 1.00000
O2 O 4 c 0.25000 0.25000 0.15904 1.00000
H3 H 16 g 0.80440 0.57590 0.12010 1.00000
H4 H 16 g -0.04700 0.48740 0.13270 1.00000
H5 H 16 g 0.80540 0.57660 0.86230 1.00000
H6 H 16 g 0.83920 0.45960 -0.03980 1.00000
O3 O 16 g 0.84543 0.47219 0.12549 1.00000
O4 O 16 g 0.84587 0.47366 0.87655 1.00000
```

$Sr(OH)_2(H_2O)_8$: A18B10C_tP116_130_2c4g_2c2g_a - POSCAR

```
A18B10C_tP116_130_2c4g_2c2g_a & a,c/a,z2,z3,z4,z5,x6,y6,z6,x7,y7,z7,x8,
    ↪ y8,z8,x9,y9,z9,x10,y10,z10,x11,y11,z11 --params=8.984,
    ↪ 1.28194568121,0.4888,0.2426,0.40482,0.15904,0.8044,0.5759,
    ↪ 0.1201,-0.047,0.4874,0.1327,0.8054,0.5766,0.8623,0.8392,0.4596
    ↪ ,-0.0398,0.84543,0.47219,0.12549,0.84587,0.47366,0.87655 & P4/
    ↪ ncc D_{4h}^{8} #130 (ac^4g^6) & tP116 & None & H18O10Sr &
    ↪ H18O10Sr & J. S. Ricci et al., Acta Crystallogr. Sect. B
    ↪ Struct. Sci. 61, 381-386 (2005)
   1.00000000000000
   8.98400000000000    0.00000000000000    0.00000000000000
   0.00000000000000    8.98400000000000    0.00000000000000
   0.00000000000000    0.00000000000000   11.51700000000000
     H      O      Sr
    72     40      4
Direct
   0.25000000000000    0.25000000000000    0.48880000000000    H    (4c)
   0.75000000000000    0.75000000000000    0.01120000000000    H    (4c)
   0.75000000000000    0.75000000000000   -0.48880000000000    H    (4c)
   0.25000000000000    0.25000000000000    0.98880000000000    H    (4c)
   0.25000000000000    0.25000000000000    0.24260000000000    H    (4c)
   0.75000000000000    0.75000000000000    0.25740000000000    H    (4c)
   0.75000000000000    0.75000000000000   -0.24260000000000    H    (4c)
   0.25000000000000    0.25000000000000    0.74260000000000    H    (4c)
   0.80440000000000    0.57590000000000    0.12010000000000    H   (16g)
  -0.30440000000000   -0.07590000000000    0.12010000000000    H   (16g)
  -0.07590000000000    0.80440000000000    0.12010000000000    H   (16g)
   0.57590000000000   -0.30440000000000    0.12010000000000    H   (16g)
  -0.80440000000000    1.07590000000000    0.37990000000000    H   (16g)
   1.30440000000000   -0.57590000000000    0.37990000000000    H   (16g)
   1.07590000000000    1.30440000000000    0.37990000000000    H   (16g)
  -0.57590000000000   -0.80440000000000    0.37990000000000    H   (16g)
  -0.80440000000000   -0.57590000000000   -0.12010000000000    H   (16g)
   1.30440000000000    1.07590000000000   -0.12010000000000    H   (16g)
   1.07590000000000   -0.80440000000000   -0.12010000000000    H   (16g)
  -0.57590000000000    1.30440000000000   -0.12010000000000    H   (16g)
   0.80440000000000   -0.07590000000000    0.62010000000000    H   (16g)
  -0.30440000000000    0.57590000000000    0.62010000000000    H   (16g)
  -0.07590000000000   -0.30440000000000    0.62010000000000    H   (16g)
   0.57590000000000    0.80440000000000    0.62010000000000    H   (16g)
  -0.04700000000000    0.48740000000000    0.13270000000000    H   (16g)
   0.54700000000000    0.01260000000000    0.13270000000000    H   (16g)
   0.01260000000000   -0.04700000000000    0.13270000000000    H   (16g)
   0.48740000000000    0.54700000000000    0.13270000000000    H   (16g)
   0.04700000000000    0.98740000000000    0.36730000000000    H   (16g)
   0.45300000000000   -0.48740000000000    0.36730000000000    H   (16g)
   0.98740000000000    0.45300000000000    0.36730000000000    H   (16g)
  -0.48740000000000    0.04700000000000    0.36730000000000    H   (16g)
   0.04700000000000   -0.48740000000000   -0.13270000000000    H   (16g)
   0.45300000000000    0.98740000000000   -0.13270000000000    H   (16g)
   0.98740000000000    0.04700000000000   -0.13270000000000    H   (16g)
  -0.48740000000000    0.45300000000000   -0.13270000000000    H   (16g)
  -0.04700000000000    0.01260000000000    0.63270000000000    H   (16g)
   0.54700000000000    0.48740000000000    0.63270000000000    H   (16g)
   0.01260000000000    0.54700000000000    0.63270000000000    H   (16g)
   0.48740000000000   -0.04700000000000    0.63270000000000    H   (16g)
   0.80540000000000    0.57660000000000    0.86230000000000    H   (16g)
  -0.30540000000000   -0.07660000000000    0.86230000000000    H   (16g)
  -0.07660000000000    0.80540000000000    0.86230000000000    H   (16g)
   0.57660000000000   -0.30540000000000    0.86230000000000    H   (16g)
  -0.80540000000000    1.07660000000000   -0.36230000000000    H   (16g)
   1.30540000000000   -0.57660000000000   -0.36230000000000    H   (16g)
   1.07660000000000    1.30540000000000   -0.36230000000000    H   (16g)
  -0.57660000000000   -0.80540000000000   -0.36230000000000    H   (16g)
  -0.80540000000000   -0.57660000000000   -0.86230000000000    H   (16g)
   1.30540000000000    1.07660000000000   -0.86230000000000    H   (16g)
   1.07660000000000   -0.80540000000000   -0.86230000000000    H   (16g)
  -0.57660000000000    1.30540000000000   -0.86230000000000    H   (16g)
   0.80540000000000   -0.07660000000000    1.36230000000000    H   (16g)
  -0.30540000000000    0.57660000000000    1.36230000000000    H   (16g)
  -0.07660000000000   -0.30540000000000    1.36230000000000    H   (16g)
   0.57660000000000    0.80540000000000    1.36230000000000    H   (16g)
   0.83920000000000    0.45960000000000   -0.03980000000000    H   (16g)
  -0.33920000000000    0.04040000000000   -0.03980000000000    H   (16g)
   0.04040000000000    0.83920000000000   -0.03980000000000    H   (16g)
   0.45960000000000   -0.33920000000000   -0.03980000000000    H   (16g)
  -0.83920000000000    0.95960000000000    0.53980000000000    H   (16g)
   1.33920000000000   -0.45960000000000    0.53980000000000    H   (16g)
   0.95960000000000    1.33920000000000    0.53980000000000    H   (16g)
  -0.45960000000000   -0.83920000000000    0.53980000000000    H   (16g)
  -0.83920000000000   -0.45960000000000    0.03980000000000    H   (16g)
   1.33920000000000    0.95960000000000    0.03980000000000    H   (16g)
   0.95960000000000   -0.83920000000000    0.03980000000000    H   (16g)
  -0.45960000000000    1.33920000000000    0.03980000000000    H   (16g)
   0.83920000000000    0.04040000000000    0.46020000000000    H   (16g)
  -0.33920000000000    0.45960000000000    0.46020000000000    H   (16g)
   0.04040000000000   -0.33920000000000    0.46020000000000    H   (16g)
   0.45960000000000    0.83920000000000    0.46020000000000    H   (16g)
   0.25000000000000    0.25000000000000    0.40482000000000    O    (4c)
   0.75000000000000    0.75000000000000    0.09518000000000    O    (4c)
   0.75000000000000    0.75000000000000   -0.40482000000000    O    (4c)
   0.25000000000000    0.25000000000000    0.90482000000000    O    (4c)
   0.25000000000000    0.25000000000000    0.15904000000000    O    (4c)
   0.75000000000000    0.75000000000000    0.34096000000000    O    (4c)
   0.75000000000000    0.75000000000000   -0.15904000000000    O    (4c)
   0.25000000000000    0.25000000000000    0.65904000000000    O    (4c)
   0.84543000000000    0.47219000000000    0.12549000000000    O   (16g)
  -0.34543000000000    0.02781000000000    0.12549000000000    O   (16g)
   0.02781000000000    0.84543000000000    0.12549000000000    O   (16g)
   0.47219000000000   -0.34543000000000    0.12549000000000    O   (16g)
  -0.84543000000000    0.97219000000000    0.37451000000000    O   (16g)
   1.34543000000000   -0.47219000000000    0.37451000000000    O   (16g)
   0.97219000000000    1.34543000000000    0.37451000000000    O   (16g)
  -0.47219000000000   -0.84543000000000    0.37451000000000    O   (16g)
  -0.84543000000000   -0.47219000000000   -0.12549000000000    O   (16g)
```

```
  1.34543000000000   0.97219000000000  −0.12549000000000   O  (16g)
  0.97219000000000  −0.84543000000000  −0.12549000000000   O  (16g)
 −0.47219000000000   1.34543000000000  −0.12549000000000   O  (16g)
  0.84543000000000   0.02781000000000   0.62549000000000   O  (16g)
 −0.34543000000000   0.47219000000000   0.62549000000000   O  (16g)
  0.02781000000000  −0.34543000000000   0.62549000000000   O  (16g)
  0.47219000000000   0.84543000000000   0.62549000000000   O  (16g)
  0.84587000000000   0.47366000000000   0.87655000000000   O  (16g)
 −0.34587000000000   0.02634000000000   0.87655000000000   O  (16g)
  0.02634000000000   0.84587000000000   0.87655000000000   O  (16g)
  0.47366000000000  −0.34587000000000   0.87655000000000   O  (16g)
 −0.84587000000000   0.97366000000000  −0.37655000000000   O  (16g)
  1.34587000000000  −0.47366000000000  −0.37655000000000   O  (16g)
  0.97366000000000   1.34587000000000  −0.37655000000000   O  (16g)
 −0.47366000000000  −0.84587000000000  −0.37655000000000   O  (16g)
 −0.84587000000000  −0.47366000000000  −0.87655000000000   O  (16g)
  1.34587000000000   0.97366000000000  −0.87655000000000   O  (16g)
  0.97366000000000  −0.84587000000000  −0.87655000000000   O  (16g)
 −0.47366000000000   1.34587000000000  −0.87655000000000   O  (16g)
  0.84587000000000   0.02634000000000   1.37655000000000   O  (16g)
 −0.34587000000000   0.47366000000000   1.37655000000000   O  (16g)
  0.02634000000000  −0.34587000000000   1.37655000000000   O  (16g)
  0.47366000000000   0.84587000000000   1.37655000000000   O  (16g)
  0.75000000000000   0.25000000000000   0.25000000000000  Sr  (4a)
  0.25000000000000   0.75000000000000   0.25000000000000  Sr  (4a)
  0.25000000000000   0.75000000000000   0.75000000000000  Sr  (4a)
  0.75000000000000   0.25000000000000   0.75000000000000  Sr  (4a)
```

α-WO_3: A3B_tP16_130_cf_c - CIF

```
# CIF file
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_audit_creation_method FINDSYM

_chemical_name_mineral 'O3W'
_chemical_formula_sum 'O3 W'

loop_
_publ_author_name
 'P. M. Woodward'
 'A. W. Sleight'
 'T. Vogt'
_journal_name_full_name
;
 Journal of Solid State Chemistry
;
_journal_volume 131
_journal_year 1997
_journal_page_first 9
_journal_page_last 17
_publ_Section_title
;
 Ferroelectric Tungsten Trioxide
;

_aflow_title '$\alpha$-WO$_{3}$ Structure'
_aflow_proto 'A3B_tP16_130_cf_c'
_aflow_params 'a,c/a,z_{1},z_{2},x_{3}'
_aflow_params_values '5.2759,1.48717754317,0.0028,0.2847,0.0287'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tP16'

_symmetry_space_group_name_H-M "P 4/n 21/c 2/c (origin choice 2)"
_symmetry_Int_Tables_number 130

_cell_length_a    5.27590
_cell_length_b    5.27590
_cell_length_c    7.84620
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y,-z+1/2
3 -x,y+1/2,-z+1/2
4 -x+1/2,-y+1/2,z
5 -y,-x,-z+1/2
6 -y+1/2,x,z
7 y,-x+1/2,z
8 y+1/2,x+1/2,-z+1/2
9 -x,-y,-z
10 -x+1/2,y,z+1/2
11 x,-y+1/2,z+1/2
12 x+1/2,y+1/2,-z
13 y,x,z+1/2
14 y+1/2,-x,-z
15 -y,x+1/2,-z
16 -y+1/2,-x+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 4 c 0.25000 0.25000 0.00280 1.00000
W1 W 4 c 0.25000 0.25000 0.28470 1.00000
O2 O 8 f 0.02870 -0.02870 0.25000 1.00000
```

α-WO_3: A3B_tP16_130_cf_c - POSCAR

```
A3B_tP16_130_cf_c & a,c/a,z1,z2,x3 --params=5.2759,1.48717754317,0.0028,
    ↪ 0.2847,0.0287 & P4/ncc D_{4h}^{8} #130 (c^2f) & tP16 & None &
    ↪ O3W & O3W & P. M. Woodward and A. W. Sleight and T. Vogt, J.
    ↪ Solid State Chem. 131, 9-17 (1997)
   1.00000000000000
   5.27590000000000   0.00000000000000   0.00000000000000
   0.00000000000000   5.27590000000000   0.00000000000000
   0.00000000000000   0.00000000000000   7.84620000000000
     O     W
    12     4
Direct
   0.25000000000000   0.25000000000000   0.00280000000000   O  (4c)
   0.75000000000000   0.75000000000000   0.49720000000000   O  (4c)
   0.75000000000000   0.75000000000000  -0.00280000000000   O  (4c)
   0.25000000000000   0.25000000000000   0.50280000000000   O  (4c)
   0.02870000000000  -0.02870000000000   0.25000000000000   O  (8f)
   0.47130000000000   0.52870000000000   0.25000000000000   O  (8f)
   0.52870000000000   0.02870000000000   0.25000000000000   O  (8f)
  -0.02870000000000   0.47130000000000   0.25000000000000   O  (8f)
  -0.02870000000000   0.02870000000000   0.75000000000000   O  (8f)
   0.52870000000000   0.47130000000000   0.75000000000000   O  (8f)
   0.47130000000000  -0.02870000000000   0.75000000000000   O  (8f)
   0.02870000000000   0.52870000000000   0.75000000000000   O  (8f)
   0.25000000000000   0.25000000000000   0.28470000000000   W  (4c)
   0.75000000000000   0.75000000000000   0.21530000000000   W  (4c)
   0.75000000000000   0.75000000000000  -0.28470000000000   W  (4c)
   0.25000000000000   0.25000000000000   0.78470000000000   W  (4c)
```

Zr_3Al_2: A2B3_tP20_136_j_dfg - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Al2Zr3'
_chemical_formula_sum 'Al2 Zr3'

loop_
_publ_author_name
 'C. G. Wilson'
 'F. J. Spooner'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 13
_journal_year 1960
_journal_page_first 358
_journal_page_last 359
_publ_Section_title
;
 The Crystal Structure of Zr$_{3}$Al$_{2}$
;

# Found in Crystal Structure Investigations on the Zr-Al and Hf-Al
    ↪ Systems, 1962

_aflow_title 'Zr$_{3}$Al$_{2}$ Structure'
_aflow_proto 'A2B3_tP20_136_j_dfg'
_aflow_params 'a,c/a,x_{2},x_{3},x_{4},z_{4}'
_aflow_params_values '7.63,0.917169069463,0.34,0.2,0.125,0.21'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tP20'

_symmetry_space_group_name_H-M "P 42/m 21/n 2/m"
_symmetry_Int_Tables_number 136

_cell_length_a    7.63000
_cell_length_b    7.63000
_cell_length_c    6.99800
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x+1/2,y+1/2,-z+1/2
4 -x,-y,z
5 -y,-x,-z
6 -y+1/2,x+1/2,z+1/2
7 y+1/2,-x+1/2,z+1/2
8 y,x,-z
9 -x,-y,-z
10 -x+1/2,y+1/2,z+1/2
11 x+1/2,-y+1/2,z+1/2
12 x,y,-z
13 y,x,z
14 y+1/2,-x+1/2,-z+1/2
15 -y+1/2,x+1/2,-z+1/2
16 -y,-x,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
```

```
Zr1  Zr  4 d  0.00000  0.50000  0.25000  1.00000
Zr2  Zr  4 f  0.34000  0.34000  0.00000  1.00000
Zr3  Zr  4 g  0.20000  0.80000  0.00000  1.00000
Al1  Al  8 j  0.12500  0.12500  0.21000  1.00000
```

Zr_3Al_2: A2B3_tP20_136_j_dfg - POSCAR

```
A2B3_tP20_136_j_dfg & a,c/a,x2,x3,x4,z4 --params=7.63,0.917169069463,
      ↪ 0.34,0.2,0.125,0.21 & P4_{2}/mmm D_{4h}^{14} #136 (dfgj) & tP20
      ↪  & None & Al2Zr3 & Al2Zr3 & C. G. Wilson and F. J. Spooner,
      ↪ Acta Cryst. 13, 358-359 (1960)
   1.00000000000000
   7.63000000000000    0.00000000000000    0.00000000000000
   0.00000000000000    7.63000000000000    0.00000000000000
   0.00000000000000    0.00000000000000    6.99800000000000
     Al     Zr
      8     12
Direct
   0.12500000000000    0.12500000000000    0.21000000000000    Al    (8j)
  -0.12500000000000   -0.12500000000000    0.21000000000000    Al    (8j)
   0.37500000000000    0.62500000000000    0.71000000000000    Al    (8j)
   0.62500000000000    0.37500000000000    0.71000000000000    Al    (8j)
   0.37500000000000    0.62500000000000    0.29000000000000    Al    (8j)
   0.62500000000000    0.37500000000000    0.29000000000000    Al    (8j)
   0.12500000000000    0.12500000000000   -0.21000000000000    Al    (8j)
  -0.12500000000000   -0.12500000000000   -0.21000000000000    Al    (8j)
   0.00000000000000    0.50000000000000    0.25000000000000    Zr    (4d)
   0.00000000000000    0.50000000000000    0.75000000000000    Zr    (4d)
   0.50000000000000    0.00000000000000    0.25000000000000    Zr    (4d)
   0.50000000000000    0.00000000000000    0.75000000000000    Zr    (4d)
   0.34000000000000    0.34000000000000    0.00000000000000    Zr    (4f)
  -0.34000000000000   -0.34000000000000    0.00000000000000    Zr    (4f)
   0.16000000000000    0.84000000000000    0.50000000000000    Zr    (4f)
   0.84000000000000    0.16000000000000    0.50000000000000    Zr    (4f)
   0.20000000000000   -0.20000000000000    0.00000000000000    Zr    (4g)
  -0.20000000000000    0.20000000000000    0.00000000000000    Zr    (4g)
   0.70000000000000    0.70000000000000    0.50000000000000    Zr    (4g)
   0.30000000000000    0.30000000000000    0.50000000000000    Zr    (4g)
```

$ZrFe_4Si_2$: A4B2C_tP14_136_i_g_b - CIF

```
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data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Fe4Si2Zr'
_chemical_formula_sum 'Fe4 Si2 Zr'

loop_
_publ_author_name
 'Y. P. Yarmolyuk'
 'L. A. Lysenko'
 'E. I. Gladyshevsky'
_journal_name_full_name
;
 Dopov. Akad. Nauk Ukr. RSR, Ser. A
;
_journal_volume 37
_journal_year 1975
_journal_page_first 281
_journal_page_last 284
_publ_Section_title
;
 Crystal Structure of ZrFe$_{4}$Si$_{2}$ -- A New Structure Type of
      ↪ Ternary Silicides
;

# Found in Simultaneous structural and magnetic transitions in YFe$_{4}
      ↪ $Ge$_{2}$ studied by neutron diffraction and magnetic
      ↪ measurements, 2001

_aflow_title 'ZrFe$_{4}$Si$_{2}$ Structure'
_aflow_proto 'A4B2C_tP14_136_i_g_b'
_aflow_params 'a,c/a,x_{2},x_{3},y_{3}'
_aflow_params_values '7.004,0.536122215877,0.2201,0.092,0.3468'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tP14'

_symmetry_space_group_name_H-M "P 42/m 21/n 2/m"
_symmetry_Int_Tables_number 136

_cell_length_a     7.00400
_cell_length_b     7.00400
_cell_length_c     3.75500
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x+1/2,y+1/2,-z+1/2
4 -x,-y,z
5 -y,-x,-z
6 -y+1/2,x+1/2,z+1/2
7 y+1/2,-x+1/2,z+1/2
8 y,x,-z
9 -x,-y,-z
10 -x+1/2,y+1/2,z+1/2
11 x+1/2,-y+1/2,z+1/2
12 x,y,-z
13 y,x,z
14 y+1/2,-x+1/2,-z+1/2
15 -y+1/2,x+1/2,-z+1/2
16 -y,-x,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Zr1  Zr  2 b  0.00000  0.00000  0.50000  1.00000
Si1  Si  4 g  0.22010  0.77990  0.00000  1.00000
Fe1  Fe  8 i  0.09200  0.34680  0.00000  1.00000
```

$ZrFe_4Si_2$: A4B2C_tP14_136_i_g_b - POSCAR

```
A4B2C_tP14_136_i_g_b & a,c/a,x2,x3,y3 --params=7.004,0.536122215877,
      ↪ 0.2201,0.092,0.3468 & P4_{2}/mmm D_{4h}^{14} #136 (bgi) & tP14
      ↪ & None & Fe4Si2Zr & Fe4Si2Zr & Y. P. Yarmolyuk and L. A.
      ↪ Lysenko and E. I. Gladyshevsky, Dopov. Akad. Nauk Ukr. RSR,
      ↪ Ser. A 37, 281-284 (1975)
   1.00000000000000
   7.00400000000000    0.00000000000000    0.00000000000000
   0.00000000000000    7.00400000000000    0.00000000000000
   0.00000000000000    0.00000000000000    3.75500000000000
     Fe     Si     Zr
      8      4      2
Direct
   0.09200000000000    0.34680000000000    0.00000000000000    Fe    (8i)
  -0.09200000000000   -0.34680000000000    0.00000000000000    Fe    (8i)
   0.15320000000000    0.59200000000000    0.50000000000000    Fe    (8i)
   0.84680000000000    0.40800000000000    0.50000000000000    Fe    (8i)
   0.40800000000000    0.84680000000000    0.50000000000000    Fe    (8i)
   0.59200000000000    0.15320000000000    0.50000000000000    Fe    (8i)
   0.34680000000000    0.09200000000000    0.00000000000000    Fe    (8i)
  -0.34680000000000   -0.09200000000000    0.00000000000000    Fe    (8i)
   0.22010000000000   -0.22010000000000    0.00000000000000    Si    (4g)
  -0.22010000000000    0.22010000000000    0.00000000000000    Si    (4g)
   0.72010000000000    0.72010000000000    0.50000000000000    Si    (4g)
   0.27990000000000    0.27990000000000    0.50000000000000    Si    (4g)
   0.00000000000000    0.00000000000000    0.50000000000000    Zr    (2b)
   0.50000000000000    0.50000000000000    0.00000000000000    Zr    (2b)
```

$K_2CuCl_4{\cdot}2H_2O$ ($H4_1$): A4BC4D2E2_tP26_136_fg_a_j_d_e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Cl4CuH4K2O2'
_chemical_formula_sum 'Cl4 Cu H4 K2 O2'

loop_
_publ_author_name
 'R. Chidambaram'
 'Q. O. Navarro'
 'A. Garcia'
 'K. Linggoatmodjo'
 'L. {Shi-Chien}'
 'I.-H. Suh'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 26
_journal_year 1970
_journal_page_first 827
_journal_page_last 830
_publ_Section_title
;
 Neutron diffraction refinement of the crystal structure of potassium
      ↪ copper chloride dihydrate, K$_{2}$CuCl$_{4}$ $\cdot$ 2H$_{2}$O
;

_aflow_title 'K$_{2}$CuCl$_{4}$$\cdot$2H$_{2}$O ($H4_{1}$) Structure'
_aflow_proto 'A4BC4D2E2_tP26_136_fg_a_j_d_e'
_aflow_params 'a,c/a,z_{3},x_{4},x_{5},x_{6},z_{6}'
_aflow_params_values '7.477,1.06125451384,0.2484,0.2161,0.7262,0.0739,
      ↪ 0.3178'
_aflow_Strukturbericht '$H4_{1}$'
_aflow_Pearson 'tP26'

_symmetry_space_group_name_H-M "P 42/m 21/n 2/m"
_symmetry_Int_Tables_number 136

_cell_length_a     7.47700
_cell_length_b     7.47700
_cell_length_c     7.93500
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x+1/2,y+1/2,-z+1/2
4 -x,-y,z
5 -y,-x,-z
6 -y+1/2,x+1/2,z+1/2
7 y+1/2,-x+1/2,z+1/2
8 y,x,-z
```

```
9 -x,-y,-z
10 -x+1/2,y+1/2,z+1/2
11 x+1/2,-y+1/2,z+1/2
12 x,y,-z
13 y,x,z
14 y+1/2,-x+1/2,-z+1/2
15 -y+1/2,x+1/2,-z+1/2
16 -y,-x,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cu1 Cu 2 a 0.00000 0.00000 0.00000 1.00000
K1 K 4 d 0.00000 0.50000 0.25000 1.00000
O1 O 4 e 0.00000 0.00000 0.24840 1.00000
Cl1 Cl 4 f 0.21610 0.21610 0.00000 1.00000
Cl2 Cl 4 g 0.72620 0.27380 0.00000 1.00000
H1 H 8 j 0.07390 0.07390 0.31780 1.00000
```

$K_2CuCl_4{\cdot}2H_2O$ ($H4_1$): A4BC4D2E2_tP26_136_fg_a_j_d_e - POSCAR

```
A4BC4D2E2_tP26_136_fg_a_j_d_e & a,c/a,z3,x4,x5,x6,z6 --params=7.477,
    ↪ 1.06125451384,0.2484,0.2161,0.7262,0.0739,0.3178 & P4_{2}/mnm
    ↪ D_{4h}^{14} #136 (adefgj) & tP26 & $H4_{1}$ & Cl4CuH4K2O2 &
    ↪ Cl4CuH4K2O2 & R. Chidambaram et al., Acta Crystallogr. Sect. B
    ↪ Struct. Sci. 26, 827-830 (1970)
  1.00000000000000
  7.47700000000000   0.00000000000000   0.00000000000000
  0.00000000000000   7.47700000000000   0.00000000000000
  0.00000000000000   0.00000000000000   7.93500000000000
   Cl   Cu   H   K   O
    8    2   8   4   4
Direct
  0.21610000000000   0.21610000000000   0.00000000000000  Cl  (4f)
 -0.21610000000000  -0.21610000000000   0.00000000000000  Cl  (4f)
  0.28390000000000   0.71610000000000   0.50000000000000  Cl  (4f)
  0.71610000000000   0.28390000000000   0.50000000000000  Cl  (4f)
  0.72620000000000  -0.72620000000000   0.00000000000000  Cl  (4g)
 -0.72620000000000   0.72620000000000   0.00000000000000  Cl  (4g)
  1.22620000000000   1.22620000000000   0.50000000000000  Cl  (4g)
 -0.22620000000000  -0.22620000000000   0.50000000000000  Cl  (4g)
  0.00000000000000   0.00000000000000   0.00000000000000  Cu  (2a)
  0.50000000000000   0.50000000000000   0.50000000000000  Cu  (2a)
  0.07390000000000   0.07390000000000   0.31780000000000  H  (8j)
 -0.07390000000000  -0.07390000000000   0.31780000000000  H  (8j)
  0.42610000000000   0.57390000000000   0.81780000000000  H  (8j)
  0.57390000000000   0.42610000000000   0.81780000000000  H  (8j)
  0.42610000000000   0.57390000000000   0.18220000000000  H  (8j)
  0.57390000000000   0.42610000000000   0.18220000000000  H  (8j)
  0.07390000000000   0.07390000000000  -0.31780000000000  H  (8j)
 -0.07390000000000  -0.07390000000000  -0.31780000000000  H  (8j)
  0.00000000000000   0.50000000000000   0.25000000000000  K  (4d)
  0.00000000000000   0.50000000000000   0.75000000000000  K  (4d)
  0.50000000000000   0.00000000000000   0.25000000000000  K  (4d)
  0.50000000000000   0.00000000000000   0.75000000000000  K  (4d)
  0.00000000000000   0.00000000000000   0.24840000000000  O  (4e)
  0.50000000000000   0.50000000000000   0.74840000000000  O  (4e)
  0.50000000000000   0.50000000000000   0.25160000000000  O  (4e)
  0.00000000000000   0.00000000000000  -0.24840000000000  O  (4e)
```

$Nd_2Fe_{14}B$: AB14C2_tP68_136_f_ce2j2k_fg - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'BFe14Nd2'
_chemical_formula_sum 'B Fe14 Nd2'

loop_
_publ_author_name
 'D. Givord'
 'H. S. Li'
 'J. M. Moreau'
_journal_name_full_name
;
 Solid State Communications
;
_journal_volume 50
_journal_year 1984
_journal_page_first 497
_journal_page_last 499
_publ_Section_title
;
 Magnetic properties and crystal structure of Nd${2}$Fe${14}$B
;

_aflow_title 'Nd${2}$Fe${14}$B Structure'
_aflow_proto 'AB14C2_tP68_136_f_ce2j2k_fg'
_aflow_params 'a,c/a,z_{2},x_{3},x_{4},x_{5},x_{6},z_{6},x_{7},z_{7},x_{
    ↪ 8},y_{8},z_{8},x_{9},y_{9},z_{9}'
_aflow_params_values '8.792,1.38648771611,0.116,0.124,0.3572,0.7698,
    ↪ 0.0978,0.2942,0.3184,0.255,0.567,0.2245,0.3735,0.1397,0.537,
    ↪ 0.1759'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tP68'

_symmetry_space_group_name_H-M "P 42/m 21/n 2/m"
_symmetry_Int_Tables_number 136

_cell_length_a    8.79200
_cell_length_b    8.79200
_cell_length_c    12.19000
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z+1/2
3 -x+1/2,y+1/2,-z+1/2
4 -x,-y,z
5 -y,-x,-z
6 -y+1/2,x+1/2,z+1/2
7 y+1/2,-x+1/2,z+1/2
8 y,x,-z
9 -x,-y,-z
10 -x+1/2,y+1/2,z+1/2
11 x+1/2,-y+1/2,z+1/2
12 x,y,-z
13 y,x,z
14 y+1/2,-x+1/2,-z+1/2
15 -y+1/2,x+1/2,-z+1/2
16 -y,-x,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Fe1 Fe 4 c 0.00000 0.50000 0.00000 1.00000
Fe2 Fe 4 e 0.00000 0.00000 0.11600 1.00000
B1 B 4 f 0.12400 0.12400 0.00000 1.00000
Nd1 Nd 4 f 0.35720 0.35720 0.00000 1.00000
Nd2 Nd 4 g 0.76980 0.23020 0.00000 1.00000
Fe3 Fe 8 j 0.09780 0.09780 0.29420 1.00000
Fe4 Fe 8 j 0.31840 0.31840 0.25500 1.00000
Fe5 Fe 16 k 0.56700 0.22450 0.37350 1.00000
Fe6 Fe 16 k 0.13970 0.53700 0.17590 1.00000
```

$Nd_2Fe_{14}B$: AB14C2_tP68_136_f_ce2j2k_fg - POSCAR

```
AB14C2_tP68_136_f_ce2j2k_fg & a,c/a,z2,x3,x4,x5,x6,z6,x7,z7,x8,y8,z8,x9,
    ↪ y9,z9 --params=8.792,1.38648771611,0.116,0.124,0.3572,0.7698,
    ↪ 0.0978,0.2942,0.3184,0.255,0.567,0.2245,0.3735,0.1397,0.537,
    ↪ 0.1759 & P4_{2}/mnm D_{4h}^{14} #136 (cef^2gj^2k^2) & tP68 &
    ↪ None & BFe14Nd2 & BFe14Nd2 & D. Givord and H. S. Li and J. M.
    ↪ Moreau, Solid State Commun. 50, 497-499 (1984)
  1.00000000000000
  8.79200000000000   0.00000000000000   0.00000000000000
  0.00000000000000   8.79200000000000   0.00000000000000
  0.00000000000000   0.00000000000000  12.19000000000000
    B   Fe   Nd
    4   56    8
Direct
  0.12400000000000   0.12400000000000   0.00000000000000  B  (4f)
 -0.12400000000000  -0.12400000000000   0.00000000000000  B  (4f)
  0.37600000000000   0.62400000000000   0.50000000000000  B  (4f)
  0.62400000000000   0.37600000000000   0.50000000000000  B  (4f)
  0.00000000000000   0.50000000000000   0.00000000000000  Fe  (4c)
  0.00000000000000   0.50000000000000   0.50000000000000  Fe  (4c)
  0.50000000000000   0.00000000000000   0.50000000000000  Fe  (4c)
  0.50000000000000   0.00000000000000   0.00000000000000  Fe  (4c)
  0.00000000000000   0.00000000000000   0.11600000000000  Fe  (4e)
  0.50000000000000   0.50000000000000   0.61600000000000  Fe  (4e)
  0.50000000000000   0.50000000000000   0.38400000000000  Fe  (4e)
  0.00000000000000   0.00000000000000  -0.11600000000000  Fe  (4e)
  0.09780000000000   0.09780000000000   0.29420000000000  Fe  (8j)
 -0.09780000000000  -0.09780000000000   0.29420000000000  Fe  (8j)
  0.40220000000000   0.59780000000000   0.79420000000000  Fe  (8j)
  0.59780000000000   0.40220000000000   0.79420000000000  Fe  (8j)
  0.40220000000000   0.59780000000000   0.20580000000000  Fe  (8j)
  0.59780000000000   0.40220000000000   0.20580000000000  Fe  (8j)
  0.09780000000000   0.09780000000000  -0.29420000000000  Fe  (8j)
 -0.09780000000000  -0.09780000000000  -0.29420000000000  Fe  (8j)
  0.31840000000000   0.31840000000000   0.25500000000000  Fe  (8j)
 -0.31840000000000  -0.31840000000000   0.25500000000000  Fe  (8j)
  0.18160000000000   0.81840000000000   0.75500000000000  Fe  (8j)
  0.81840000000000   0.18160000000000   0.75500000000000  Fe  (8j)
  0.18160000000000   0.81840000000000   0.24500000000000  Fe  (8j)
  0.81840000000000   0.18160000000000   0.24500000000000  Fe  (8j)
  0.31840000000000   0.31840000000000  -0.25500000000000  Fe  (8j)
 -0.31840000000000  -0.31840000000000  -0.25500000000000  Fe  (8j)
  0.56700000000000   0.22450000000000   0.37350000000000  Fe  (16k)
 -0.56700000000000  -0.22450000000000   0.37350000000000  Fe  (16k)
  0.27550000000000   1.06700000000000   0.87350000000000  Fe  (16k)
  0.72450000000000  -0.06700000000000   0.87350000000000  Fe  (16k)
 -0.06700000000000   0.72450000000000   0.12650000000000  Fe  (16k)
  1.06700000000000   0.27550000000000   0.12650000000000  Fe  (16k)
  0.22450000000000   0.56700000000000  -0.37350000000000  Fe  (16k)
 -0.22450000000000  -0.56700000000000  -0.37350000000000  Fe  (16k)
 -0.56700000000000  -0.22450000000000  -0.37350000000000  Fe  (16k)
  0.56700000000000   0.22450000000000  -0.37350000000000  Fe  (16k)
  0.72450000000000  -0.06700000000000   0.12650000000000  Fe  (16k)
  0.27550000000000   1.06700000000000   0.12650000000000  Fe  (16k)
  1.06700000000000   0.27550000000000   0.87350000000000  Fe  (16k)
 -0.06700000000000   0.72450000000000   0.87350000000000  Fe  (16k)
 -0.22450000000000  -0.56700000000000   0.37350000000000  Fe  (16k)
```

```
   0.22450000000000   0.56700000000000   0.37350000000000   Fe   (16k)
   0.13970000000000   0.53700000000000   0.17590000000000   Fe   (16k)
  -0.13970000000000  -0.53700000000000   0.17590000000000   Fe   (16k)
  -0.03700000000000   0.63970000000000   0.67590000000000   Fe   (16k)
   1.03700000000000   0.36030000000000   0.67590000000000   Fe   (16k)
   0.36030000000000   1.03700000000000   0.32410000000000   Fe   (16k)
   0.63970000000000  -0.03700000000000   0.32410000000000   Fe   (16k)
   0.53700000000000   0.13970000000000  -0.17590000000000   Fe   (16k)
  -0.53700000000000  -0.13970000000000  -0.17590000000000   Fe   (16k)
  -0.13970000000000  -0.53700000000000  -0.17590000000000   Fe   (16k)
   0.13970000000000   0.53700000000000  -0.17590000000000   Fe   (16k)
   1.03700000000000   0.36030000000000   0.32410000000000   Fe   (16k)
  -0.03700000000000   0.63970000000000   0.32410000000000   Fe   (16k)
   0.63970000000000  -0.03700000000000   0.67590000000000   Fe   (16k)
   0.36030000000000   1.03700000000000   0.67590000000000   Fe   (16k)
  -0.53700000000000  -0.13970000000000   0.17590000000000   Fe   (16k)
   0.53700000000000   0.13970000000000   0.17590000000000   Fe   (16k)
   0.35720000000000   0.35720000000000   0.00000000000000   Nd    (4f)
  -0.35720000000000  -0.35720000000000   0.00000000000000   Nd    (4f)
   0.14280000000000   0.85720000000000   0.50000000000000   Nd    (4f)
   0.85720000000000   0.14280000000000   0.50000000000000   Nd    (4f)
   0.76980000000000  -0.76980000000000   0.00000000000000   Nd    (4g)
  -0.76980000000000   0.76980000000000   0.00000000000000   Nd    (4g)
   1.26980000000000   1.26980000000000   0.50000000000000   Nd    (4g)
  -0.26980000000000  -0.26980000000000   0.50000000000000   Nd    (4g)
```

$Sr_4Ti_3O_{10}$: A10B4C3_tI34_139_c2eg_2e_ae - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'O10Sr4Ti3'
_chemical_formula_sum 'O10 Sr4 Ti3'

loop_
_publ_author_name
 'S. N. Ruddlesden'
 'P. Popper'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 11
_journal_year 1958
_journal_page_first 54
_journal_page_last 55
_publ_Section_title
;
 The compound Sr${3}$Ti${2}$O${7}$ and its structure
;

# Found in Ruddlesden-Popper phase, {A${3}$B${2}$X${7}$ series},

_aflow_title 'Sr${4}$Ti${3}$O${10}$ Structure'
_aflow_proto 'A10B4C3_tI34_139_c2eg_2e_ae'
_aflow_params 'a,c/a,z_{3},z_{4},z_{5},z_{6},z_{7},z_{8}'
_aflow_params_values '3.9,7.20512820513,0.068,0.204,0.432,0.296,0.136,
    ↪ 0.136'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tI34'

_symmetry_space_group_name_H-M "I 4/m 2/m 2/m"
_symmetry_Int_Tables_number 139

_cell_length_a    3.90000
_cell_length_b    3.90000
_cell_length_c    28.10000
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 -y,-x,-z
6 -y,x,z
7 y,-x,z
8 y,x,-z
9 -x,-y,-z
10 -x,y,z
11 x,-y,z
12 x,y,-z
13 y,x,z
14 y,-x,-z
15 -y,x,-z
16 -y,-x,z
17 x+1/2,y+1/2,z+1/2
18 x+1/2,-y+1/2,-z+1/2
19 -x+1/2,y+1/2,-z+1/2
20 -x+1/2,-y+1/2,z+1/2
21 -y+1/2,-x+1/2,-z+1/2
22 -y+1/2,x+1/2,z+1/2
23 y+1/2,-x+1/2,z+1/2
24 y+1/2,x+1/2,-z+1/2
25 -x+1/2,-y+1/2,-z+1/2
26 -x+1/2,y+1/2,z+1/2
27 x+1/2,-y+1/2,z+1/2
28 x+1/2,y+1/2,-z+1/2
29 y+1/2,x+1/2,z+1/2
30 y+1/2,-x+1/2,-z+1/2
31 -y+1/2,x+1/2,-z+1/2
32 -y+1/2,-x+1/2,z+1/2

loop_
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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ti1 Ti 2 a 0.00000 0.00000 0.00000 1.00000
O1 O 4 c 0.00000 0.50000 0.00000 1.00000
O2 O 4 e 0.00000 0.00000 0.06800 1.00000
O3 O 4 e 0.00000 0.00000 0.20400 1.00000
Sr1 Sr 4 e 0.00000 0.00000 0.43200 1.00000
Sr2 Sr 4 e 0.00000 0.00000 0.29600 1.00000
Ti2 Ti 4 e 0.00000 0.00000 0.13600 1.00000
O4 O 8 g 0.00000 0.50000 0.13600 1.00000
```

$Sr_4Ti_3O_{10}$: A10B4C3_tI34_139_c2eg_2e_ae - POSCAR

```
A10B4C3_tI34_139_c2eg_2e_ae & a,c/a,z3,z4,z5,z6,z7,z8 --params=3.9,
    ↪ 7.20512820513,0.068,0.204,0.432,0.296,0.136,0.136 & I4/mmm D_{
    ↪ 4h}^{17} #139 (ace^5g) & tI34 & None & O10Sr4Ti3 & O10Sr4Ti3 &
    ↪ S. N. Ruddlesden and P. Popper, Acta Cryst. 11, 54-55 (1958)
   1.00000000000000
  -1.95000000000000   1.95000000000000  14.05000000000000
   1.95000000000000  -1.95000000000000  14.05000000000000
   1.95000000000000   1.95000000000000 -14.05000000000000
     O    Sr    Ti
    10     4     3
Direct
   0.50000000000000   0.00000000000000   0.50000000000000   O   (4c)
   0.00000000000000   0.50000000000000   0.50000000000000   O   (4c)
   0.06800000000000   0.06800000000000   0.00000000000000   O   (4e)
  -0.06800000000000  -0.06800000000000   0.00000000000000   O   (4e)
   0.20400000000000   0.20400000000000   0.00000000000000   O   (4e)
  -0.20400000000000  -0.20400000000000   0.00000000000000   O   (4e)
   0.63600000000000   0.13600000000000   0.50000000000000   O   (8g)
   0.13600000000000   0.63600000000000   0.50000000000000   O   (8g)
   0.36400000000000  -0.13600000000000   0.50000000000000   O   (8g)
  -0.13600000000000   0.36400000000000   0.50000000000000   O   (8g)
   0.43200000000000   0.43200000000000   0.00000000000000   Sr  (4e)
  -0.43200000000000  -0.43200000000000   0.00000000000000   Sr  (4e)
   0.29600000000000   0.29600000000000   0.00000000000000   Sr  (4e)
  -0.29600000000000  -0.29600000000000   0.00000000000000   Sr  (4e)
   0.00000000000000   0.00000000000000   0.00000000000000   Ti  (2a)
   0.13600000000000   0.13600000000000   0.00000000000000   Ti  (4e)
  -0.13600000000000  -0.13600000000000   0.00000000000000   Ti  (4e)
```

$TlCo_2S_2$: A2B2C_tI10_139_d_e_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Co2S2Tl'
_chemical_formula_sum 'Co2 S2 Tl'

loop_
_publ_author_name
 'K. Klepp'
 'H. Boller'
_journal_name_full_name
;
 Monatshefte f{\"u}r Chemie - Chemical Monthly
;
_journal_volume 109
_journal_year 1978
_journal_page_first 1049
_journal_page_last 1057
_publ_Section_title
;
 Tern\"{a}re Thallium-\"{U}bergangsmetall-Chalkogenide mit ThCr$_{2}
    ↪ $Si$_{2}$-Struktur
;

# Found in TlCu$_{2}$Se$_{2}$: A p-type metal with a layer structure,
    ↪ 1984

_aflow_title 'TlCo$_{2}$S$_{2}$ Structure'
_aflow_proto 'A2B2C_tI10_139_d_e_a'
_aflow_params 'a,c/a,z_{3}'
_aflow_params_values '3.741,3.46324779471,0.3474'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tI10'

_symmetry_space_group_name_H-M "I 4/m 2/m 2/m"
_symmetry_Int_Tables_number 139

_cell_length_a    3.74100
_cell_length_b    3.74100
_cell_length_c    12.95601
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
```

```
4 -x,-y,z
5 -y,-x,-z
6 -y,x,z
7 y,-x,z
8 y,x,-z
9 -x,-y,-z
10 -x,y,z
11 x,-y,z
12 x,y,-z
13 y,x,z
14 y,-x,-z
15 -y,x,-z
16 -y,-x,z
17 x+1/2,y+1/2,z+1/2
18 x+1/2,-y+1/2,-z+1/2
19 -x+1/2,y+1/2,-z+1/2
20 -x+1/2,-y+1/2,z+1/2
21 -y+1/2,-x+1/2,-z+1/2
22 -y+1/2,x+1/2,z+1/2
23 y+1/2,-x+1/2,z+1/2
24 y+1/2,x+1/2,-z+1/2
25 -x+1/2,-y+1/2,-z+1/2
26 -x+1/2,y+1/2,z+1/2
27 x+1/2,-y+1/2,z+1/2
28 x+1/2,y+1/2,-z+1/2
29 y+1/2,x+1/2,z+1/2
30 y+1/2,-x+1/2,-z+1/2
31 -y+1/2,x+1/2,-z+1/2
32 -y+1/2,-x+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Tl1 Tl 2 a 0.00000 0.00000 0.00000 1.00000
Co1 Co 4 d 0.00000 0.50000 0.25000 1.00000
S1 S 4 e 0.00000 0.00000 0.34740 1.00000
```

$TlCo_2S_2$: A2B2C_tI10_139_d_e_a - POSCAR

```
A2B2C_tI10_139_d_e_a & a,c/a,z3 --params=3.741,3.46324779471,0.3474 & I4
      ↪ /mmm D_{4h}^{17} #139 (ade) & tI10 & None & Co2S2Tl & Co2S2Tl &
      ↪  K. Klepp and H. Boller, Monatshefte f{"u}r Chemie - Chemical
      ↪ Monthly 109, 1049-1057 (1978)
   1.00000000000000
  -1.87050000000000    1.87050000000000    6.47800500000000
   1.87050000000000   -1.87050000000000    6.47800500000000
   1.87050000000000    1.87050000000000   -6.47800500000000
    Co      S      Tl
     2      2       1
Direct
   0.75000000000000    0.25000000000000    0.50000000000000    Co    (4d)
   0.25000000000000    0.75000000000000    0.50000000000000    Co    (4d)
   0.34740000000000    0.34740000000000    0.00000000000000    S     (4e)
  -0.34740000000000   -0.34740000000000    0.00000000000000    S     (4e)
   0.00000000000000    0.00000000000000    0.00000000000000    Tl    (2a)
```

$ThCr_2Si_2$: A2B2C_tI10_139_d_e_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Cr2Si2Th'
_chemical_formula_sum 'Cr2 Si2 Th'

loop_
_publ_author_name
 'Z. Ban'
 'M. Sikirica'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 18
_journal_year 1965
_journal_page_first 594
_journal_page_last 599
_publ_Section_title
;
 The crystal structure of ternary silicides Th$M$$_{2}$Si$_{2}$ ($M$ =
      ↪ Cr, Mn, Fe, Co, Ni and Cu)
;

_aflow_title 'ThCr$_{2}$Si$_{2}$ Structure'
_aflow_proto 'A2B2C_tI10_139_d_e_a'
_aflow_params 'a,c/a,z_{3}'
_aflow_params_values '4.043,2.61612663863,0.374'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tI10'

_symmetry_space_group_name_H-M "I 4/m 2/m 2/m"
_symmetry_Int_Tables_number 139

_cell_length_a     4.04300
_cell_length_b     4.04300
_cell_length_c     10.57700
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 -y,-x,-z
6 -y,x,z
7 y,-x,z
8 y,x,-z
9 -x,-y,-z
10 -x,y,z
11 x,-y,z
12 x,y,-z
13 y,x,z
14 y,-x,-z
15 -y,x,-z
16 -y,-x,z
17 x+1/2,y+1/2,z+1/2
18 x+1/2,-y+1/2,-z+1/2
19 -x+1/2,y+1/2,-z+1/2
20 -x+1/2,-y+1/2,z+1/2
21 -y+1/2,-x+1/2,-z+1/2
22 -y+1/2,x+1/2,z+1/2
23 y+1/2,-x+1/2,z+1/2
24 y+1/2,x+1/2,-z+1/2
25 -x+1/2,-y+1/2,-z+1/2
26 -x+1/2,y+1/2,z+1/2
27 x+1/2,-y+1/2,z+1/2
28 x+1/2,y+1/2,-z+1/2
29 y+1/2,x+1/2,z+1/2
30 y+1/2,-x+1/2,-z+1/2
31 -y+1/2,x+1/2,-z+1/2
32 -y+1/2,-x+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Th1 Th 2 a 0.00000 0.00000 0.00000 1.00000
Cr1 Cr 4 d 0.00000 0.50000 0.25000 1.00000
Si1 Si 4 e 0.00000 0.00000 0.37400 1.00000
```

$ThCr_2Si_2$: A2B2C_tI10_139_d_e_a - POSCAR

```
A2B2C_tI10_139_d_e_a & a,c/a,z3 --params=4.043,2.61612663863,0.374 & I4/
      ↪ mmm D_{4h}^{17} #139 (ade) & tI10 & None & Cr2Si2Th & Cr2Si2Th
      ↪ & Z. Ban and M. Sikirica, Acta Cryst. 18, 594-599 (1965)
   1.00000000000000
  -2.02150000000000    2.02150000000000    5.28850000000000
   2.02150000000000   -2.02150000000000    5.28850000000000
   2.02150000000000    2.02150000000000   -5.28850000000000
    Cr      Si      Th
     2       2       1
Direct
   0.75000000000000    0.25000000000000    0.50000000000000    Cr    (4d)
   0.25000000000000    0.75000000000000    0.50000000000000    Cr    (4d)
   0.37400000000000    0.37400000000000    0.00000000000000    Si    (4e)
  -0.37400000000000   -0.37400000000000    0.00000000000000    Si    (4e)
   0.00000000000000    0.00000000000000    0.00000000000000    Th    (2a)
```

Au_2Nb_3: A2B3_tI10_139_e_ae - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Au2Nb3'
_chemical_formula_sum 'Au2 Nb3'

loop_
_publ_author_name
 'K. Schubert'
 'T. R. Anantharaman'
 'H. O. K. Ata'
 'H. G. Meissner'
 'M. P\"{o}tzschke'
 'W. Rossteutscher'
 'E. Stolz'
_journal_name_full_name
;
 Naturwissenschaften
;
_journal_volume 47
_journal_year 1960
_journal_page_first 512
_journal_page_last 512
_publ_Section_title
;
 Einige strukturelle Ergebnisse an metallischen Phasen (6)
;

# Found in The Crystal Structures of Os$_{2}$Al$_{3}$ and OsAl$_{2}$,
      ↪ 1965

_aflow_title 'Au$_{2}$Nb$_{3}$ Structure'
_aflow_proto 'A2B3_tI10_139_e_ae'
_aflow_params 'a,c/a,z_{2},z_{3}'
```

```
_aflow_params_values '3.36301,4.49555309083,0.2,0.4'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tI10'

_symmetry_space_group_name_H-M "I 4/m 2/m 2/m"
_symmetry_Int_Tables_number 139

_cell_length_a 3.36301
_cell_length_b 3.36301
_cell_length_c 15.11859
_cell_angle_alpha 90.00000
_cell_angle_beta 90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 -y,-x,-z
6 -y,x,z
7 y,-x,z
8 y,x,-z
9 -x,-y,-z
10 -x,y,z
11 x,-y,z
12 x,y,-z
13 y,x,z
14 y,-x,-z
15 -y,x,-z
16 -y,-x,z
17 x+1/2,y+1/2,z+1/2
18 x+1/2,-y+1/2,-z+1/2
19 -x+1/2,y+1/2,-z+1/2
20 -x+1/2,-y+1/2,z+1/2
21 -y+1/2,-x+1/2,-z+1/2
22 -y+1/2,x+1/2,z+1/2
23 y+1/2,-x+1/2,z+1/2
24 y+1/2,x+1/2,-z+1/2
25 -x+1/2,-y+1/2,-z+1/2
26 -x+1/2,y+1/2,z+1/2
27 x+1/2,-y+1/2,z+1/2
28 x+1/2,y+1/2,-z+1/2
29 y+1/2,x+1/2,z+1/2
30 y+1/2,-x+1/2,-z+1/2
31 -y+1/2,x+1/2,-z+1/2
32 -y+1/2,-x+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Nb1 Nb 2 a 0.00000 0.00000 0.00000 1.00000
Au1 Au 4 e 0.00000 0.00000 0.20000 1.00000
Nb2 Nb 4 e 0.00000 0.00000 0.40000 1.00000
```

Au_2Nb_3: A2B3_tI10_139_e_ae - POSCAR

```
A2B3_tI10_139_e_ae & a,c/a,z2,z3 --params=3.36301,4.49555309083,0.2,0.4
    ↪ & I4/mmm D_{4h}^{17} #139 (ae^2) & tI10 & None & Au2Nb3 &
    ↪ Au2Nb3 & K. Schubert et al., Naturwissenschaften 47, 512(1960)
   1.00000000000000
  -1.68150500000000   1.68150500000000   7.55929500000000
   1.68150500000000  -1.68150500000000   7.55929500000000
   1.68150500000000   1.68150500000000  -7.55929500000000
    Au  Nb
     2   3
Direct
   0.20000000000000   0.20000000000000   0.00000000000000  Au  (4e)
  -0.20000000000000  -0.20000000000000   0.00000000000000  Au  (4e)
   0.00000000000000   0.00000000000000   0.00000000000000  Nb  (2a)
   0.40000000000000   0.40000000000000   0.00000000000000  Nb  (4e)
  -0.40000000000000  -0.40000000000000   0.00000000000000  Nb  (4e)
```

CaC_2-I ($C11_a$): A2B_tI6_139_e_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'C2Ca'
_chemical_formula_sum 'C2 Ca'

loop_
_publ_author_name
 'M. {von Stackelberg}'
_journal_name_full_name
;
 Naturwissenschaften
;
_journal_volume 18
_journal_year 1930
_journal_page_first 305
_journal_page_last 306
_publ_Section_title
;
 Die Krystallstruktur des CaC$_{2}$
;

# Found in The Crystal Structure of Calcium Carbide III, 1961

_aflow_title 'CaC$_{2}$-I ($C11_{a}$) Structure'
_aflow_proto 'A2B_tI6_139_e_a'
_aflow_params 'a,c/a,z_{2}'
_aflow_params_values '3.87,1.64599483204,0.38'
_aflow_Strukturbericht '$C11_{a}$'
_aflow_Pearson 'tI6'

_symmetry_space_group_name_H-M "I 4/m 2/m 2/m"
_symmetry_Int_Tables_number 139

_cell_length_a 3.87000
_cell_length_b 3.87000
_cell_length_c 6.37000
_cell_angle_alpha 90.00000
_cell_angle_beta 90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 -y,-x,-z
6 -y,x,z
7 y,-x,z
8 y,x,-z
9 -x,-y,-z
10 -x,y,z
11 x,-y,z
12 x,y,-z
13 y,x,z
14 y,-x,-z
15 -y,x,-z
16 -y,-x,z
17 x+1/2,y+1/2,z+1/2
18 x+1/2,-y+1/2,-z+1/2
19 -x+1/2,y+1/2,-z+1/2
20 -x+1/2,-y+1/2,z+1/2
21 -y+1/2,-x+1/2,-z+1/2
22 -y+1/2,x+1/2,z+1/2
23 y+1/2,-x+1/2,z+1/2
24 y+1/2,x+1/2,-z+1/2
25 -x+1/2,-y+1/2,-z+1/2
26 -x+1/2,y+1/2,z+1/2
27 x+1/2,-y+1/2,z+1/2
28 x+1/2,y+1/2,-z+1/2
29 y+1/2,x+1/2,z+1/2
30 y+1/2,-x+1/2,-z+1/2
31 -y+1/2,x+1/2,-z+1/2
32 -y+1/2,-x+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ca1 Ca 2 a 0.00000 0.00000 0.00000 1.00000
C1 C 4 e 0.00000 0.00000 0.38000 1.00000
```

CaC_2-I ($C11_a$): A2B_tI6_139_e_a - POSCAR

```
A2B_tI6_139_e_a & a,c/a,z2 --params=3.87,1.64599483204,0.38 & I4/mmm D_{
    ↪ 4h}^{17} #139 (ae) & tI6 & $C11_{a}$ & C2Ca & C2Ca & M. {von
    ↪ Stackelberg}, Naturwissenschaften 18, 305-306 (1930)
   1.00000000000000
  -1.93500000000000   1.93500000000000   3.18500000000000
   1.93500000000000  -1.93500000000000   3.18500000000000
   1.93500000000000   1.93500000000000  -3.18500000000000
    C  Ca
    2   1
Direct
   0.38000000000000   0.38000000000000   0.00000000000000   C   (4e)
  -0.38000000000000  -0.38000000000000   0.00000000000000   C   (4e)
   0.00000000000000   0.00000000000000   0.00000000000000  Ca   (2a)
```

$K_2OsO_2Cl_4$ ($J1_5$): A4B2C2D_tI18_139_h_d_e_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Cl4K2O2Os'
_chemical_formula_sum 'Cl4 K2 O2 Os'

loop_
_publ_author_name
 'J. L. Hoard'
 'J. D. Grenko'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 87
_journal_year 1934
_journal_page_first 100
_journal_page_last 109
```

```
_publ_Section_title
;
 The Crystal Structure of Potassium Osmyl Chloride, K$_{2}$OsO$_{2}$Cl$_
      ↪ {4}$
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'K$_{2}$OsO$_{2}$Cl$_{4}$ ($J1_{5}$) Structure'
_aflow_proto 'A4B2C2D_tI18_139_h_d_e_a'
_aflow_params 'a,c/a,z_{3},x_{4}'
_aflow_params_values '6.99,1.25178826896,0.212,0.23'
_aflow_Strukturbericht '$J1_{5}$'
_aflow_Pearson 'tI18'

_symmetry_space_group_name_H-M "I 4/m 2/m 2/m"
_symmetry_Int_Tables_number 139

_cell_length_a     6.99000
_cell_length_b     6.99000
_cell_length_c     8.75000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 -y,-x,-z
6 -y,x,z
7 y,-x,z
8 y,x,-z
9 -x,-y,-z
10 -x,y,z
11 x,-y,z
12 x,y,-z
13 y,x,z
14 y,-x,-z
15 -y,x,-z
16 -y,-x,z
17 x+1/2,y+1/2,z+1/2
18 x+1/2,-y+1/2,-z+1/2
19 -x+1/2,y+1/2,-z+1/2
20 -x+1/2,-y+1/2,z+1/2
21 -y+1/2,-x+1/2,-z+1/2
22 -y+1/2,x+1/2,z+1/2
23 y+1/2,-x+1/2,z+1/2
24 y+1/2,x+1/2,-z+1/2
25 -x+1/2,-y+1/2,-z+1/2
26 -x+1/2,y+1/2,z+1/2
27 x+1/2,-y+1/2,z+1/2
28 x+1/2,y+1/2,-z+1/2
29 y+1/2,x+1/2,z+1/2
30 y+1/2,-x+1/2,-z+1/2
31 -y+1/2,x+1/2,-z+1/2
32 -y+1/2,-x+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Os1 Os 2 a 0.00000 0.00000 0.00000 1.00000
K1 K 4 d 0.00000 0.50000 0.25000 1.00000
O1 O 4 e 0.00000 0.00000 0.21200 1.00000
Cl1 Cl 8 h 0.23000 0.23000 0.00000 1.00000
```

$K_2OsO_2Cl_4$ ($J1_5$): A4B2C2D_tI18_139_h_d_e_a - POSCAR

```
A4B2C2D_tI18_139_h_d_e_a & a,c/a,z3,x4 --params=6.99,1.25178826896,0.212
      ↪ ,0.23 & I4/mmm D_{4h}^{17} #139 (adeh) & tI18 & $J1_{5}$ &
      ↪ Cl4K2O2Os & Cl4K2O2Os & J. L. Hoard and J. D. Grenko,
      ↪ Zeitschrift f{"u}r Kristallographie - Crystalline Materials 87,
      ↪  100-109 (1934)
   1.00000000000000
  -3.49500000000000    3.49500000000000    4.37500000000000
   3.49500000000000   -3.49500000000000    4.37500000000000
   3.49500000000000    3.49500000000000   -4.37500000000000
    Cl     K     O     Os
     4     2     2      1
Direct
   0.23000000000000    0.23000000000000    0.46000000000000    Cl    (8h)
  -0.23000000000000   -0.23000000000000   -0.46000000000000    Cl    (8h)
   0.23000000000000   -0.23000000000000    0.00000000000000    Cl    (8h)
  -0.23000000000000    0.23000000000000    0.00000000000000    Cl    (8h)
   0.75000000000000    0.25000000000000    0.50000000000000     K    (4d)
   0.25000000000000    0.75000000000000    0.50000000000000     K    (4d)
   0.21200000000000    0.21200000000000    0.00000000000000     O    (4e)
  -0.21200000000000   -0.21200000000000    0.00000000000000     O    (4e)
   0.00000000000000    0.00000000000000    0.00000000000000    Os    (2a)
```

K_2NiF_4: A4B2C_tI14_139_ce_e_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM
_chemical_name_mineral 'F4K2Ni'
_chemical_formula_sum 'F4 K2 Ni'

loop_
_publ_author_name
 'S. N. Ruddlesden'
 'P. Popper'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 10
_journal_year 1957
_journal_page_first 538
_journal_page_last 539
_publ_Section_title
;
 New compounds of the K$_{2}$NiF$_{4}$ type
;

# Found in Ruddlesden-Popper phase, {A$_{2}$BX$_{4}$ series},

_aflow_title 'K$_{2}$NiF$_{4}$ Structure'
_aflow_proto 'A4B2C_tI14_139_ce_e_a'
_aflow_params 'a,c/a,z_{3},z_{4}'
_aflow_params_values '4.0,3.2675,0.151,0.352'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tI14'

_symmetry_space_group_name_H-M "I 4/m 2/m 2/m"
_symmetry_Int_Tables_number 139

_cell_length_a     4.00000
_cell_length_b     4.00000
_cell_length_c     13.07000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 -y,-x,-z
6 -y,x,z
7 y,-x,z
8 y,x,-z
9 -x,-y,-z
10 -x,y,z
11 x,-y,z
12 x,y,-z
13 y,x,z
14 y,-x,-z
15 -y,x,-z
16 -y,-x,z
17 x+1/2,y+1/2,z+1/2
18 x+1/2,-y+1/2,-z+1/2
19 -x+1/2,y+1/2,-z+1/2
20 -x+1/2,-y+1/2,z+1/2
21 -y+1/2,-x+1/2,-z+1/2
22 -y+1/2,x+1/2,z+1/2
23 y+1/2,-x+1/2,z+1/2
24 y+1/2,x+1/2,-z+1/2
25 -x+1/2,-y+1/2,-z+1/2
26 -x+1/2,y+1/2,z+1/2
27 x+1/2,-y+1/2,z+1/2
28 x+1/2,y+1/2,-z+1/2
29 y+1/2,x+1/2,z+1/2
30 y+1/2,-x+1/2,-z+1/2
31 -y+1/2,x+1/2,-z+1/2
32 -y+1/2,-x+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ni1 Ni 2 a 0.00000 0.00000 0.00000 1.00000
F1 F 4 c 0.00000 0.50000 0.00000 1.00000
F2 F 4 e 0.00000 0.00000 0.15100 1.00000
K1 K 4 e 0.00000 0.00000 0.35200 1.00000
```

K_2NiF_4: A4B2C_tI14_139_ce_e_a - POSCAR

```
A4B2C_tI14_139_ce_e_a & a,c/a,z3,z4 --params=4.0,3.2675,0.151,0.352 & I4
      ↪ /mmm D_{4h}^{17} #139 (ace^2) & tI14 & None & F4K2Ni & F4K2Ni &
      ↪  S. N. Ruddlesden and P. Popper, Acta Cryst. 10, 538-539 (1957)
   1.00000000000000
  -2.00000000000000    2.00000000000000    6.53500000000000
   2.00000000000000   -2.00000000000000    6.53500000000000
   2.00000000000000    2.00000000000000   -6.53500000000000
     F     K     Ni
     4     2     1
Direct
   0.50000000000000    0.00000000000000    0.50000000000000     F    (4c)
   0.00000000000000    0.50000000000000    0.50000000000000     F    (4c)
   0.15100000000000    0.15100000000000    0.00000000000000     F    (4e)
  -0.15100000000000   -0.15100000000000    0.00000000000000     F    (4e)
```

```
  0.35200000000000   0.35200000000000   0.00000000000000   K   (4e)
 -0.35200000000000  -0.35200000000000   0.00000000000000   K   (4e)
  0.00000000000000   0.00000000000000   0.00000000000000   Ni  (2a)
```

$K_3TlCl_6{\cdot}2H_2O$ ($J3_1$): A6B2C3D_tI168_139_egikl2m_ejn_bh2n_acf - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Cl6(H2O)2K3Tl'
_chemical_formula_sum 'Cl6 (H2O)2 K3 Tl'

loop_
_publ_author_name
 'J. L. Hoard'
 'L. Goldstein'
_journal_name_full_name
;
 Journal of Chemical Physics
;
_journal_volume 3
_journal_year 1935
_journal_page_first 645
_journal_page_last 649
_publ_Section_title
;
 The Structure of Potassium Hexachlorothalliate Dihydrate
;

_aflow_title 'K$_{3}$TlCl$_{6}$$\cdot$2H$_{2}$O ($J3_{1}$) Structure'
_aflow_proto 'A6B2C3D_tI168_139_egikl2m_ejn_bh2n_acf'
_aflow_params 'a,c/a,z_{4},z_{5},z_{7},x_{8},x_{9},x_{10},x_{11},x_{12},
      ↪ y_{12},x_{13},z_{13},x_{14},z_{14},y_{15},z_{15},y_{16},z_{16},
      ↪ y_{17},z_{17}'
_aflow_params_values '15.841,1.1366075374,0.142,0.347,0.142,0.214,0.161,
      ↪ 0.673,0.364,0.386,0.114,0.181,0.362,0.16,0.163,0.157,0.276,
      ↪ 0.295,0.132,0.293,0.376'
_aflow_Strukturbericht '$J3_{1}$'
_aflow_Pearson 'tI168'

_symmetry_space_group_name_H-M "I 4/m 2/m 2/m"
_symmetry_Int_Tables_number 139

_cell_length_a     15.84100
_cell_length_b     15.84100
_cell_length_c     18.00500
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 -y,-x,-z
6 -y,x,z
7 y,-x,z
8 y,x,-z
9 -x,-y,-z
10 -x,y,z
11 x,-y,z
12 x,y,-z
13 y,x,z
14 y,-x,-z
15 -y,x,-z
16 -y,-x,z
17 x+1/2,y+1/2,z+1/2
18 x+1/2,-y+1/2,-z+1/2
19 -x+1/2,y+1/2,-z+1/2
20 -x+1/2,-y+1/2,z+1/2
21 -y+1/2,-x+1/2,-z+1/2
22 -y+1/2,x+1/2,z+1/2
23 y+1/2,-x+1/2,z+1/2
24 y+1/2,x+1/2,-z+1/2
25 -x+1/2,-y+1/2,-z+1/2
26 -x+1/2,y+1/2,z+1/2
27 x+1/2,-y+1/2,z+1/2
28 x+1/2,y+1/2,-z+1/2
29 y+1/2,x+1/2,z+1/2
30 y+1/2,-x+1/2,-z+1/2
31 -y+1/2,x+1/2,-z+1/2
32 -y+1/2,-x+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Tl1 Tl 2 a 0.00000 0.00000 0.00000 1.00000
K1 K 2 b 0.00000 0.00000 0.50000 1.00000
Tl2 Tl 4 c 0.00000 0.50000 0.00000 1.00000
Cl1 Cl 4 e 0.00000 0.00000 0.14200 1.00000
H2O1 H2O 4 e 0.00000 0.00000 0.34700 1.00000
Tl3 Tl 8 f 0.25000 0.25000 0.25000 1.00000
Cl2 Cl 8 g 0.00000 0.50000 0.14200 1.00000
K2 K 8 h 0.21400 0.21400 0.00000 1.00000
Cl3 Cl 8 i 0.16100 0.00000 0.00000 1.00000
H2O2 H2O 8 j 0.67300 0.50000 0.00000 1.00000
Cl4 Cl 16 k 0.36400 0.86400 0.25000 1.00000
Cl5 Cl 16 l 0.38600 0.11400 0.00000 1.00000
Cl6 Cl 16 m 0.18100 0.18100 0.36200 1.00000
Cl7 Cl 16 m 0.16000 0.16000 0.16300 1.00000
H2O3 H2O 16 n 0.00000 0.15700 0.27600 1.00000
K3 K 16 n 0.00000 0.29500 0.13200 1.00000
K4 K 16 n 0.00000 0.29300 0.37600 1.00000
```

$K_3TlCl_6{\cdot}2H_2O$ ($J3_1$): A6B2C3D_tI168_139_egikl2m_ejn_bh2n_acf - POSCAR

```
A6B2C3D_tI168_139_egikl2m_ejn_bh2n_acf & a,c/a,z4,z5,z7,x8,x9,x10,x11,
      ↪ x12,y12,x13,z13,x14,z14,y15,z15,y16,z16,y17,z17 --params=15.841
      ↪ ,1.1366075374,0.142,0.347,0.142,0.214,0.161,0.673,0.364,0.386,
      ↪ 0.114,0.181,0.362,0.16,0.163,0.157,0.276,0.295,0.132,0.293,
      ↪ 0.376 & I4/mmm D_{4h}^{17} #139 (abce^2fghijklm^2n^3) & tI168 &
      ↪  $J3_{1}$ & Cl6(H2O)2K3Tl & Cl6(H2O)2K3Tl & J. L. Hoard and L.
      ↪ Goldstein, J. Chem. Phys. 3, 645-649 (1935)
   1.00000000000000
  -7.92050000000000   7.92050000000000   9.00250000000000
   7.92050000000000  -7.92050000000000   9.00250000000000
   7.92050000000000   7.92050000000000  -9.00250000000000
    Cl    H2O     K     Tl
    42     14     21      7
Direct
   0.14200000000000   0.14200000000000   0.00000000000000   Cl   (4e)
  -0.14200000000000  -0.14200000000000   0.00000000000000   Cl   (4e)
   0.64200000000000   0.14200000000000   0.50000000000000   Cl   (8g)
   0.14200000000000   0.64200000000000   0.50000000000000   Cl   (8g)
   0.35800000000000  -0.14200000000000   0.50000000000000   Cl   (8g)
  -0.14200000000000   0.35800000000000   0.50000000000000   Cl   (8g)
   0.00000000000000   0.16100000000000   0.16100000000000   Cl   (8i)
   0.00000000000000  -0.16100000000000  -0.16100000000000   Cl   (8i)
   0.16100000000000   0.00000000000000  -0.16100000000000   Cl   (8i)
  -0.16100000000000   0.00000000000000  -0.16100000000000   Cl   (8i)
   1.11400000000000   0.61400000000000   1.22800000000000   Cl   (16k)
   0.38600000000000  -0.11400000000000  -0.22800000000000   Cl   (16k)
   0.61400000000000   0.38600000000000   0.50000000000000   Cl   (16k)
  -0.11400000000000   1.11400000000000   0.50000000000000   Cl   (16k)
  -0.11400000000000   0.38600000000000  -0.22800000000000   Cl   (16k)
   0.61400000000000   1.11400000000000   1.22800000000000   Cl   (16k)
   0.38600000000000   0.61400000000000   0.50000000000000   Cl   (16k)
   1.11400000000000  -0.11400000000000   0.50000000000000   Cl   (16k)
   0.11400000000000   0.38600000000000   0.50000000000000   Cl   (16l)
  -0.11400000000000  -0.38600000000000  -0.50000000000000   Cl   (16l)
   0.38600000000000  -0.11400000000000   0.27200000000000   Cl   (16l)
  -0.38600000000000   0.11400000000000  -0.27200000000000   Cl   (16l)
   0.11400000000000  -0.38600000000000  -0.27200000000000   Cl   (16l)
  -0.11400000000000   0.38600000000000   0.27200000000000   Cl   (16l)
   0.38600000000000   0.11400000000000   0.50000000000000   Cl   (16l)
  -0.38600000000000  -0.11400000000000  -0.50000000000000   Cl   (16l)
   0.54300000000000   0.54300000000000   0.36200000000000   Cl   (16m)
   0.18100000000000   0.18100000000000  -0.36200000000000   Cl   (16m)
   0.54300000000000   0.18100000000000   0.00000000000000   Cl   (16m)
   0.18100000000000   0.54300000000000   0.00000000000000   Cl   (16m)
  -0.18100000000000  -0.54300000000000   0.00000000000000   Cl   (16m)
  -0.54300000000000  -0.18100000000000   0.00000000000000   Cl   (16m)
  -0.18100000000000  -0.18100000000000   0.36200000000000   Cl   (16m)
  -0.54300000000000  -0.54300000000000  -0.36200000000000   Cl   (16m)
   0.32300000000000   0.32300000000000   0.32000000000000   Cl   (16m)
   0.00300000000000   0.00300000000000  -0.32000000000000   Cl   (16m)
   0.32300000000000   0.00300000000000   0.00000000000000   Cl   (16m)
   0.00300000000000   0.32300000000000   0.00000000000000   Cl   (16m)
  -0.00300000000000  -0.32300000000000   0.00000000000000   Cl   (16m)
  -0.32300000000000  -0.00300000000000   0.00000000000000   Cl   (16m)
  -0.00300000000000  -0.00300000000000   0.32000000000000   Cl   (16m)
  -0.32300000000000  -0.32300000000000  -0.32000000000000   Cl   (16m)
   0.34700000000000   0.34700000000000   0.00000000000000   H2O  (4e)
  -0.34700000000000  -0.34700000000000   0.00000000000000   H2O  (4e)
   0.50000000000000   0.67300000000000   1.17300000000000   H2O  (8j)
   0.50000000000000  -0.67300000000000  -0.17300000000000   H2O  (8j)
   0.67300000000000   0.50000000000000   1.17300000000000   H2O  (8j)
  -0.67300000000000   0.50000000000000  -0.17300000000000   H2O  (8j)
   0.43300000000000   0.27600000000000   0.15700000000000   H2O  (16n)
   0.11900000000000   0.27600000000000  -0.15700000000000   H2O  (16n)
   0.27600000000000   0.11900000000000  -0.15700000000000   H2O  (16n)
   0.27600000000000   0.43300000000000   0.15700000000000   H2O  (16n)
  -0.11900000000000  -0.27600000000000   0.15700000000000   H2O  (16n)
  -0.43300000000000  -0.27600000000000  -0.15700000000000   H2O  (16n)
  -0.27600000000000  -0.11900000000000   0.15700000000000   H2O  (16n)
  -0.27600000000000  -0.43300000000000  -0.15700000000000   H2O  (16n)
   0.50000000000000   0.50000000000000   0.00000000000000   K    (2b)
   0.21400000000000   0.21400000000000   0.42800000000000   K    (8h)
  -0.21400000000000  -0.21400000000000  -0.42800000000000   K    (8h)
   0.21400000000000  -0.21400000000000   0.00000000000000   K    (8h)
  -0.21400000000000   0.21400000000000   0.00000000000000   K    (8h)
   0.42700000000000   0.13200000000000   0.29500000000000   K    (16n)
  -0.16300000000000   0.13200000000000  -0.29500000000000   K    (16n)
   0.13200000000000  -0.16300000000000  -0.29500000000000   K    (16n)
   0.13200000000000   0.42700000000000   0.29500000000000   K    (16n)
   0.16300000000000  -0.13200000000000   0.29500000000000   K    (16n)
  -0.42700000000000  -0.13200000000000  -0.29500000000000   K    (16n)
  -0.13200000000000   0.16300000000000   0.29500000000000   K    (16n)
  -0.13200000000000  -0.42700000000000  -0.29500000000000   K    (16n)
   0.66900000000000   0.37600000000000   0.29300000000000   K    (16n)
   0.08300000000000   0.37600000000000  -0.29300000000000   K    (16n)
   0.37600000000000   0.08300000000000  -0.29300000000000   K    (16n)
   0.37600000000000   0.66900000000000   0.29300000000000   K    (16n)
  -0.08300000000000  -0.37600000000000   0.29300000000000   K    (16n)
  -0.66900000000000  -0.37600000000000  -0.29300000000000   K    (16n)
  -0.37600000000000  -0.08300000000000   0.29300000000000   K    (16n)
  -0.37600000000000  -0.66900000000000  -0.29300000000000   K    (16n)
   0.00000000000000   0.00000000000000   0.00000000000000   Tl   (2a)
   0.50000000000000   0.00000000000000   0.50000000000000   Tl   (4c)
   0.00000000000000   0.50000000000000   0.50000000000000   Tl   (4c)
```

```
   0.50000000000000   0.50000000000000   0.50000000000000   Tl   (8f)
   0.00000000000000   0.00000000000000   0.50000000000000   Tl   (8f)
   0.50000000000000   0.00000000000000   0.00000000000000   Tl   (8f)
   0.00000000000000   0.50000000000000   0.00000000000000   Tl   (8f)
```

$Sr_3Ti_2O_7$: A7B3C2_tI24_139_aeg_be_e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'O7Sr3Ti2'
_chemical_formula_sum 'O7 Sr3 Ti2'

loop_
_publ_author_name
 'S. N. Ruddlesden'
 'P. Popper'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 11
_journal_year 1958
_journal_page_first 54
_journal_page_last 55
_publ_Section_title
;
 The compound Sr$_{3}$Ti$_{2}$O$_{7}$ and its structure
;

# Found in Ruddlesden-Popper phase, {A$_{3}$B$_{2}$X$_{7}$ series},

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_aflow_proto 'A7B3C2_tI24_139_aeg_be_e'
_aflow_params 'a,c/a,z_{3},z_{4},z_{5},z_{6}'
_aflow_params_values '3.9,5.22564102564,0.188,0.312,0.094,0.094'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tI24'

_symmetry_space_group_name_H-M "I 4/m 2/m 2/m"
_symmetry_Int_Tables_number 139

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_cell_length_b     3.90000
_cell_length_c     20.38000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 -y,-x,-z
6 -y,x,z
7 y,-x,z
8 y,x,-z
9 -x,-y,-z
10 -x,y,z
11 x,-y,z
12 x,y,-z
13 y,x,z
14 y,-x,-z
15 -y,x,-z
16 -y,-x,z
17 x+1/2,y+1/2,z+1/2
18 x+1/2,-y+1/2,-z+1/2
19 -x+1/2,y+1/2,-z+1/2
20 -x+1/2,-y+1/2,z+1/2
21 -y+1/2,-x+1/2,-z+1/2
22 -y+1/2,x+1/2,z+1/2
23 y+1/2,-x+1/2,z+1/2
24 y+1/2,x+1/2,-z+1/2
25 -x+1/2,-y+1/2,-z+1/2
26 -x+1/2,y+1/2,z+1/2
27 x+1/2,-y+1/2,z+1/2
28 x+1/2,y+1/2,-z+1/2
29 y+1/2,x+1/2,z+1/2
30 y+1/2,-x+1/2,-z+1/2
31 -y+1/2,x+1/2,-z+1/2
32 -y+1/2,-x+1/2,z+1/2

loop_
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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 2 a 0.00000 0.00000 0.00000 1.00000
Sr1 Sr 2 b 0.00000 0.00000 0.50000 1.00000
O2 O 4 e 0.00000 0.00000 0.18800 1.00000
Sr2 Sr 4 e 0.00000 0.00000 0.31200 1.00000
Ti1 Ti 4 e 0.00000 0.00000 0.09400 1.00000
O3 O 8 g 0.00000 0.50000 0.09400 1.00000
```

$Sr_3Ti_2O_7$: A7B3C2_tI24_139_aeg_be_e - POSCAR

```
A7B3C2_tI24_139_aeg_be_e & a,c/a,z3,z4,z5,z6 --params=3.9,5.22564102564,
    ↪ 0.188,0.312,0.094,0.094 & I4/mmm D_{4h}^{17} #139 (abe^3g) &
    ↪ tI24 & None & O7Sr3Ti2 & O7Sr3Ti2 & S. N. Ruddlesden and P.
    ↪ Popper, Acta Cryst. 11, 54-55 (1958)
   1.00000000000000
  -1.95000000000000   1.95000000000000  10.19000000000000
   1.95000000000000  -1.95000000000000  10.19000000000000
   1.95000000000000   1.95000000000000 -10.19000000000000
     O     Sr     Ti
     7     3      2
Direct
   0.00000000000000   0.00000000000000   0.00000000000000   O   (2a)
   0.18800000000000   0.18800000000000   0.00000000000000   O   (4e)
  -0.18800000000000  -0.18800000000000   0.00000000000000   O   (4e)
   0.59400000000000   0.09400000000000   0.50000000000000   O   (8g)
   0.09400000000000   0.59400000000000   0.50000000000000   O   (8g)
   0.40600000000000  -0.09400000000000   0.50000000000000   O   (8g)
  -0.09400000000000   0.40600000000000   0.50000000000000   O   (8g)
   0.50000000000000   0.50000000000000   0.00000000000000   Sr  (2b)
   0.31200000000000   0.31200000000000   0.00000000000000   Sr  (4e)
  -0.31200000000000  -0.31200000000000   0.00000000000000   Sr  (4e)
   0.09400000000000   0.09400000000000   0.00000000000000   Ti  (4e)
  -0.09400000000000  -0.09400000000000   0.00000000000000   Ti  (4e)
```

Fe_8N ($D2_g$): A8B_tI18_139_deh_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Fe8N'
_chemical_formula_sum 'Fe8 N'

loop_
_publ_author_name
 'S. Yamashita'
 'Y. Masubuchi'
 'Y. Nakazawa'
 'T. Okayama'
 'M. Tsuchiya'
 'S. Kikkawa'
_journal_name_full_name
;
 Journal of Solid State Chemistry
;
_journal_volume 194
_journal_year 2012
_journal_page_first 76
_journal_page_last 79
_publ_Section_title
;
 Crystal structure and magnetic properties of ``$\alpha$$\'\'$-Fe$_{16}
    ↪ $N$_{2}$\'\' containing reisdual $\alpha$-Fe prepared by
    ↪ low-temperature ammonia nitridation
;

_aflow_title 'Fe$_{8}$N ($D2_{g}$) Structure'
_aflow_proto 'A8B_tI18_139_deh_a'
_aflow_params 'a,c/a,z_{3},x_{4}'
_aflow_params_values '5.71361,1.10055114017,0.2922,0.24433'
_aflow_Strukturbericht '$D2_{g}$'
_aflow_Pearson 'tI18'

_symmetry_space_group_name_H-M "I 4/m 2/m 2/m"
_symmetry_Int_Tables_number 139

_cell_length_a     5.71361
_cell_length_b     5.71361
_cell_length_c     6.28812
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 -y,-x,-z
6 -y,x,z
7 y,-x,z
8 y,x,-z
9 -x,-y,-z
10 -x,y,z
11 x,-y,z
12 x,y,-z
13 y,x,z
14 y,-x,-z
15 -y,x,-z
16 -y,-x,z
17 x+1/2,y+1/2,z+1/2
18 x+1/2,-y+1/2,-z+1/2
19 -x+1/2,y+1/2,-z+1/2
20 -x+1/2,-y+1/2,z+1/2
21 -y+1/2,-x+1/2,-z+1/2
22 -y+1/2,x+1/2,z+1/2
23 y+1/2,-x+1/2,z+1/2
24 y+1/2,x+1/2,-z+1/2
25 -x+1/2,-y+1/2,-z+1/2
26 -x+1/2,y+1/2,z+1/2
27 x+1/2,-y+1/2,z+1/2
28 x+1/2,y+1/2,-z+1/2
29 y+1/2,x+1/2,z+1/2
```

```
30 y+1/2,-x+1/2,-z+1/2
31 -y+1/2,x+1/2,-z+1/2
32 -y+1/2,-x+1/2,z+1/2

loop_
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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
N1 N 2 a 0.00000 0.00000 0.00000 1.00000
Fe1 Fe 4 d 0.00000 0.50000 0.25000 1.00000
Fe2 Fe 4 e 0.00000 0.00000 0.29220 1.00000
Fe3 Fe 8 h 0.24433 0.24433 0.00000 1.00000
```

Fe_8N ($D2_g$): A8B_tI18_139_deh_a - POSCAR

```
A8B_tI18_139_deh_a & a,c/a,z3,x4 --params=5.71361,1.10055114017,0.2922,
    ↪ 0.24433 & I4/mmm D_{4h}^{17} #139 (adeh) & tI18 & $D2_{g}$ &
    ↪ Fe8N & Fe8N & S. Yamashita et al., J. Solid State Chem. 194,
    ↪ 76-79 (2012)
   1.00000000000000
  -2.85680500000000   2.85680500000000   3.14406000000000
   2.85680500000000  -2.85680500000000   3.14406000000000
   2.85680500000000   2.85680500000000  -3.14406000000000
    Fe    N
     8    1
Direct
   0.75000000000000   0.25000000000000   0.50000000000000  Fe  (4d)
   0.25000000000000   0.75000000000000   0.50000000000000  Fe  (4d)
   0.29220000000000   0.29220000000000   0.00000000000000  Fe  (4e)
  -0.29220000000000  -0.29220000000000   0.00000000000000  Fe  (4e)
   0.24433000000000   0.24433000000000   0.48866000000000  Fe  (8h)
  -0.24433000000000  -0.24433000000000  -0.48866000000000  Fe  (8h)
   0.24433000000000  -0.24433000000000   0.00000000000000  Fe  (8h)
  -0.24433000000000   0.24433000000000   0.00000000000000  Fe  (8h)
   0.00000000000000   0.00000000000000   0.00000000000000   N  (2a)
```

Li_2CN_2: AB2C2_tI10_139_a_d_e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'CLi2N2'
_chemical_formula_sum 'C Li2 N2'

loop_
_publ_author_name
 'M. G. Down'
 'M. J. Haley'
 'P. Hubberstey'
 'R. J. Pulham'
 'A. E. Thunder'
_journal_name_full_name
;
 Journal of the Chemical Society, Dalton Transactions
;
_journal_volume
_journal_year 1978
_journal_page_first 1407
_journal_page_last 1411
_publ_Section_title
;
 Solutions of lithium salts in liquid lithium: preparation and X-ray
     ↪ crystal structure of the dilithium salt of carbodi-imide (
     ↪ cyanamide)
;

# Found in {Commentary: The Materials Project: A materials genome
    ↪ approach to accelerating materials innovation}, 2013

_aflow_title 'Li$_{2}$CN$_{2}$ Structure'
_aflow_proto 'AB2C2_tI10_139_a_d_e'
_aflow_params 'a,c/a,z_{3}'
_aflow_params_values '3.687,2.35096284242,0.1419'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tI10'

_symmetry_space_group_name_H-M "I 4/m 2/m 2/m"
_symmetry_Int_Tables_number 139

_cell_length_a    3.68700
_cell_length_b    3.68700
_cell_length_c    8.66800
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 -y,-x,-z
6 -y,x,z
7 y,-x,z
8 y,x,-z
9 -x,-y,-z
10 -x,y,z
11 x,-y,z
12 x,y,-z
13 y,x,z
14 y,-x,-z
15 -y,x,-z
16 -y,-x,z
17 x+1/2,y+1/2,z+1/2
18 x+1/2,-y+1/2,-z+1/2
19 -x+1/2,y+1/2,-z+1/2
20 -x+1/2,-y+1/2,z+1/2
21 -y+1/2,-x+1/2,-z+1/2
22 -y+1/2,x+1/2,z+1/2
23 y+1/2,-x+1/2,z+1/2
24 y+1/2,x+1/2,-z+1/2
25 -x+1/2,-y+1/2,-z+1/2
26 -x+1/2,y+1/2,z+1/2
27 x+1/2,-y+1/2,z+1/2
28 x+1/2,y+1/2,-z+1/2
29 y+1/2,x+1/2,z+1/2
30 y+1/2,-x+1/2,-z+1/2
31 -y+1/2,x+1/2,-z+1/2
32 -y+1/2,-x+1/2,z+1/2

loop_
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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
C1 C 2 a 0.00000 0.00000 0.00000 1.00000
Li1 Li 4 d 0.00000 0.50000 0.25000 1.00000
N1 N 4 e 0.00000 0.00000 0.14190 1.00000
```

Li_2CN_2: AB2C2_tI10_139_a_d_e - POSCAR

```
AB2C2_tI10_139_a_d_e & a,c/a,z3 --params=3.687,2.35096284242,0.1419 & I4
    ↪ /mmm D_{4h}^{17} #139 (ade) & tI10 & None & CLi2N2 & CLi2N2 &
    ↪ M. G. Down et al., {J. Chem. Soc. Dalton Trans., 1407-1411 (
    ↪ 1978)
   1.00000000000000
  -1.84350000000000   1.84350000000000   4.33400000000000
   1.84350000000000  -1.84350000000000   4.33400000000000
   1.84350000000000   1.84350000000000  -4.33400000000000
     C     Li     N
     1     2     2
Direct
   0.00000000000000   0.00000000000000   0.00000000000000   C   (2a)
   0.75000000000000   0.25000000000000   0.50000000000000   Li  (4d)
   0.25000000000000   0.75000000000000   0.50000000000000   Li  (4d)
   0.14190000000000   0.14190000000000   0.00000000000000   N   (4e)
  -0.14190000000000  -0.14190000000000   0.00000000000000   N   (4e)
```

$H5_9$ [Autunite, $Ca(UO_2)_2(PO_4)_2{\cdot}10\frac{1}{2}H_2O$] (*obsolete*): AB2C2_tI10_139_a_d_e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

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_chemical_formula_sum 'Ca P2 U2'

loop_
_publ_author_name
 'J. Beintema'
_journal_name_full_name
;
 Recueil des Travaux Chimiques des Pays-Bas
;
_journal_volume 57
_journal_year 1938
_journal_page_first 155
_journal_page_last 175
_publ_Section_title
;
 On the composition and the crystallography of autunite and the
     ↪ meta-autunites
;

_aflow_title '$H5_{9}$ [Autunite, Ca(UO$_{2}$)$_{2}$(PO$_{4}$)$_{2}$$\
    ↪ cdot$10$\frac{1}{2}$H$_{2}$O] ({\em{obsolete}}) Structure'
_aflow_proto 'AB2C2_tI10_139_a_d_e'
_aflow_params 'a,c/a,z_{3}'
_aflow_params_values '6.98,2.95558739255,0.208'
_aflow_Strukturbericht '$H5_{9}$'
_aflow_Pearson 'tI10'

_symmetry_space_group_name_H-M "I 4/m 2/m 2/m"
_symmetry_Int_Tables_number 139

_cell_length_a    6.98000
_cell_length_b    6.98000
_cell_length_c    20.63000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
```

```
4 -x,-y,z
5 -y,-x,-z
6 -y,x,z
7 y,-x,z
8 y,x,-z
9 -x,-y,-z
10 -x,y,z
11 x,-y,z
12 x,y,-z
13 y,x,z
14 y,-x,-z
15 -y,x,-z
16 -y,-x,z
17 x+1/2,y+1/2,z+1/2
18 x+1/2,-y+1/2,-z+1/2
19 -x+1/2,y+1/2,-z+1/2
20 -x+1/2,-y+1/2,z+1/2
21 -y+1/2,-x+1/2,-z+1/2
22 -y+1/2,x+1/2,z+1/2
23 y+1/2,-x+1/2,z+1/2
24 y+1/2,x+1/2,-z+1/2
25 -x+1/2,-y+1/2,-z+1/2
26 -x+1/2,y+1/2,z+1/2
27 x+1/2,-y+1/2,z+1/2
28 x+1/2,y+1/2,-z+1/2
29 y+1/2,x+1/2,z+1/2
30 y+1/2,-x+1/2,-z+1/2
31 -y+1/2,x+1/2,-z+1/2
32 -y+1/2,-x+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ca1 Ca 2 a 0.00000 0.00000 0.00000 1.00000
P1 P 4 d 0.00000 0.50000 0.25000 1.00000
U1 U 4 e 0.00000 0.00000 0.20800 1.00000
```

$H5_9$ [Autunite, $Ca(UO_2)_2(PO_4)_2{\cdot}10\frac{1}{2}H_2O$] (*obsolete*): AB2C2_tI10_139_a_d_e - POSCAR

```
AB2C2_tI10_139_a_d_e & a,c/a,z3 --params=6.98,2.95558739255,0.208 & I4/
    ↪ mmm D_{4h}^{17} #139 (ade) & tI10 & $H5_{9}$ & Ca(UO2)2(PO4)2\
    ↪ cdot10\frac{1}{2}H2O & Autunite & J. Beintema, Rec. Trav. Chim.
    ↪ Pays-Bas 57, 155-175 (1938)
   1.00000000000000
  -3.49000000000000   3.49000000000000  10.31500000000000
   3.49000000000000  -3.49000000000000  10.31500000000000
   3.49000000000000   3.49000000000000 -10.31500000000000
    Ca    P    U
     1    2    2
Direct
   0.00000000000000   0.00000000000000   0.00000000000000  Ca  (2a)
   0.75000000000000   0.25000000000000   0.50000000000000  P   (4d)
   0.25000000000000   0.75000000000000   0.50000000000000  P   (4d)
   0.20800000000000   0.20800000000000   0.00000000000000  U   (4e)
  -0.20800000000000  -0.20800000000000   0.00000000000000  U   (4e)
```

$AuCsCl_3$ ($K7_6$): AB3C_tI20_139_ab_eh_d - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'AuCl3Cs'
_chemical_formula_sum 'Au Cl3 Cs'

loop_
_publ_author_name
 'N. Elliott'
 'L. Pauling'
_journal_name_full_name
;
 Journal of the American Chemical Society
;
_journal_volume 60
_journal_year 1938
_journal_page_first 1846
_journal_page_last 1851
_publ_Section_title
;
 The Crystal Structure of Cesium Aurous Auric Chloride, Cs$_{2}$AuAuCl$_
    ↪ {6}$, and Cesium Argentous Auric Chloride, Cs$_{2}$AgAuCl$_{6}
    ↪ $
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'AuCsCl$_{3}$ ($K7_{6}$) Structure'
_aflow_proto 'AB3C_tI20_139_ab_eh_d'
_aflow_params 'a,c/a,z_{4},x_{5}'
_aflow_params_values '7.49,1.45126835781,0.288,0.228'
_aflow_Strukturbericht '$K7_{6}$'
_aflow_Pearson 'tI20'

_symmetry_space_group_name_H-M "I 4/m 2/m 2/m"
_symmetry_Int_Tables_number 139

_cell_length_a    7.49000
_cell_length_b    7.49000
_cell_length_c    10.87000
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 -y,-x,-z
6 -y,x,z
7 y,-x,z
8 y,x,-z
9 -x,-y,-z
10 -x,y,z
11 x,-y,z
12 x,y,-z
13 y,x,z
14 y,-x,-z
15 -y,x,-z
16 -y,-x,z
17 x+1/2,y+1/2,z+1/2
18 x+1/2,-y+1/2,-z+1/2
19 -x+1/2,y+1/2,-z+1/2
20 -x+1/2,-y+1/2,z+1/2
21 -y+1/2,-x+1/2,-z+1/2
22 -y+1/2,x+1/2,z+1/2
23 y+1/2,-x+1/2,z+1/2
24 y+1/2,x+1/2,-z+1/2
25 -x+1/2,-y+1/2,-z+1/2
26 -x+1/2,y+1/2,z+1/2
27 x+1/2,-y+1/2,z+1/2
28 x+1/2,y+1/2,-z+1/2
29 y+1/2,x+1/2,z+1/2
30 y+1/2,-x+1/2,-z+1/2
31 -y+1/2,x+1/2,-z+1/2
32 -y+1/2,-x+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Au1 Au 2 a 0.00000 0.00000 0.00000 1.00000
Au2 Au 2 b 0.00000 0.00000 0.50000 1.00000
Cs1 Cs 4 d 0.00000 0.50000 0.25000 1.00000
Cl1 Cl 4 e 0.00000 0.00000 0.28800 1.00000
Cl2 Cl 8 h 0.22800 0.22800 0.00000 1.00000
```

$AuCsCl_3$ ($K7_6$): AB3C_tI20_139_ab_eh_d - POSCAR

```
AB3C_tI20_139_ab_eh_d & a,c/a,z4,x5 --params=7.49,1.45126835781,0.288,
    ↪ 0.228 & I4/mmm D_{4h}^{17} #139 (abdeh) & tI20 & $K7_{6}$ &
    ↪ AuCl3Cs & AuCl3Cs & N. Elliott and L. Pauling, J. Am. Chem.
    ↪ Soc. 60, 1846-1851 (1938)
   1.00000000000000
  -3.74500000000000   3.74500000000000   5.43500000000000
   3.74500000000000  -3.74500000000000   5.43500000000000
   3.74500000000000   3.74500000000000  -5.43500000000000
    Au   Cl   Cs
     2    6    2
Direct
   0.00000000000000   0.00000000000000   0.00000000000000  Au  (2a)
   0.50000000000000   0.50000000000000   0.00000000000000  Au  (2b)
   0.28800000000000   0.28800000000000   0.00000000000000  Cl  (4e)
  -0.28800000000000  -0.28800000000000   0.00000000000000  Cl  (4e)
   0.22800000000000   0.22800000000000   0.45600000000000  Cl  (8h)
  -0.22800000000000  -0.22800000000000  -0.45600000000000  Cl  (8h)
   0.22800000000000  -0.22800000000000   0.00000000000000  Cl  (8h)
  -0.22800000000000   0.22800000000000   0.00000000000000  Cl  (8h)
   0.75000000000000   0.25000000000000   0.50000000000000  Cs  (4d)
   0.25000000000000   0.75000000000000   0.50000000000000  Cs  (4d)
```

"Martensite Type" FeC_x ($x \leq 0.06$) ($L'2_0$): AB_tI4_139_b_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral '''martensite type'''
_chemical_formula_sum 'C Fe'

loop_
_publ_author_name
 'E. A. Brandes'
 'G. B. Brook'
_journal_year 1992
_publ_Section_title
;
 Smithells Metals Reference Book
;

# Found in Strukturbericht 1913-1928, 1931

_aflow_title '''Martensite Type'' FeC$_{x}$ ($x \le 0.06$) ($L'2_{0}$)
    ↪ Structure'
_aflow_proto 'AB_tI4_139_b_a'
_aflow_params 'a,c/a'
_aflow_params_values '2.84,1.04577464789'
```

```
_aflow_Strukturbericht '$L'2_{0}$'
_aflow_Pearson 'tI4'

_symmetry_space_group_name_H-M "I 4/m 2/m 2/m"
_symmetry_Int_Tables_number 139

_cell_length_a 2.84000
_cell_length_b 2.84000
_cell_length_c 2.97000
_cell_angle_alpha 90.00000
_cell_angle_beta 90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 -y,-x,-z
6 -y,x,z
7 y,-x,z
8 y,x,-z
9 -x,-y,-z
10 -x,y,z
11 x,-y,z
12 x,y,-z
13 y,x,z
14 y,-x,-z
15 -y,x,-z
16 -y,-x,z
17 x+1/2,y+1/2,z+1/2
18 x+1/2,-y+1/2,-z+1/2
19 -x+1/2,y+1/2,-z+1/2
20 -x+1/2,-y+1/2,z+1/2
21 -y+1/2,-x+1/2,-z+1/2
22 -y+1/2,x+1/2,z+1/2
23 y+1/2,-x+1/2,z+1/2
24 y+1/2,x+1/2,-z+1/2
25 -x+1/2,-y+1/2,-z+1/2
26 -x+1/2,y+1/2,z+1/2
27 x+1/2,-y+1/2,z+1/2
28 x+1/2,y+1/2,-z+1/2
29 y+1/2,x+1/2,z+1/2
30 y+1/2,-x+1/2,-z+1/2
31 -y+1/2,x+1/2,-z+1/2
32 -y+1/2,-x+1/2,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Fe1 Fe 2 a 0.00000 0.00000 0.00000 1.00000
C1 C 2 b 0.00000 0.00000 0.50000 0.06000
```

"Martensite Type" FeC_x ($x \leq 0.06$) ($L'2_0$): AB_tI4_139_b_a - POSCAR

```
AB_tI4_139_b_a & a,c/a --params=2.84,1.04577464789 & I4/mmm D_{4h}^{17}
    ↪ #139 (ab) & tI4 & $L'2_{0}$ & CxFe & ``martensite type'' & E.
    ↪ A. Brandes and G. B. Brook, (1992)
   1.00000000000000
  -1.42000000000000   1.42000000000000   1.48500000000000
   1.42000000000000  -1.42000000000000   1.48500000000000
   1.42000000000000   1.42000000000000  -1.48500000000000
     C     Fe
     1      1
Direct
   0.50000000000000   0.50000000000000   0.00000000000000   C   (2b)
   0.00000000000000   0.00000000000000   0.00000000000000   Fe  (2a)
```

V_4SiSb_2: A2BC4_tI28_140_h_a_k - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Sb2SiV4'
_chemical_formula_sum 'Sb2 Si V4'

loop_
_publ_author_name
 'P. Wollesen'
 'W. Jeitschko'
_journal_name_full_name
;
 Journal of Alloys and Compounds
;
_journal_volume 243
_journal_year 1996
_journal_page_first 67
_journal_page_last 69
_publ_Section_title
;
 V${4}$SiSb${2}$, a vanadium silicide antimonide crystallizing with a
    ↪ defect variant of the W${5}$Si${3}$-type structure
;

_aflow_title 'V${4}$SiSb${2}$ Structure'
_aflow_proto 'A2BC4_tI28_140_h_a_k'
_aflow_params 'a,c/a,x_{2},x_{3},y_{3}'
_aflow_params_values '9.87199,0.476803562402,0.14094,0.0852,0.20674'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tI28'

_symmetry_space_group_name_H-M "I 4/m 2/c 2/m"
_symmetry_Int_Tables_number 140

_cell_length_a 9.87199
_cell_length_b 9.87199
_cell_length_c 4.70700
_cell_angle_alpha 90.00000
_cell_angle_beta 90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z+1/2
3 -x,y,-z+1/2
4 -x,-y,z
5 -y,-x,-z+1/2
6 -y,x,z
7 y,-x,z
8 y,x,-z+1/2
9 -x,-y,-z
10 -x,y,z+1/2
11 x,-y,z+1/2
12 x,y,-z
13 y,x,z+1/2
14 y,-x,-z
15 -y,x,-z
16 -y,-x,z+1/2
17 x+1/2,y+1/2,z+1/2
18 x+1/2,-y+1/2,-z
19 -x+1/2,y+1/2,-z
20 -x+1/2,-y+1/2,z+1/2
21 -y+1/2,-x+1/2,-z
22 -y+1/2,x+1/2,z+1/2
23 y+1/2,-x+1/2,z+1/2
24 y+1/2,x+1/2,-z
25 -x+1/2,-y+1/2,-z+1/2
26 -x+1/2,y+1/2,z
27 x+1/2,-y+1/2,z
28 x+1/2,y+1/2,-z+1/2
29 y+1/2,x+1/2,z
30 y+1/2,-x+1/2,-z+1/2
31 -y+1/2,x+1/2,-z+1/2
32 -y+1/2,-x+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Si1 Si 4 a 0.00000 0.00000 0.25000 1.00000
Sb1 Sb 8 h 0.14094 0.64094 0.00000 1.00000
V1 V 16 k 0.08520 0.20674 0.00000 1.00000
```

V_4SiSb_2: A2BC4_tI28_140_h_a_k - POSCAR

```
A2BC4_tI28_140_h_a_k & a,c/a,x2,x3,y3 --params=9.87199,0.476803562402,
    ↪ 0.14094,0.0852,0.20674 & I4/mcm D_{4h}^{18} #140 (ahk) & tI28 &
    ↪ None & Sb2SiV4 & Sb2SiV4 & P. Wollesen and W. Jeitschko, J.
    ↪ Alloys Compd. 243, 67-69 (1996)
   1.00000000000000
  -4.93599500000000   4.93599500000000   2.35350000000000
   4.93599500000000  -4.93599500000000   2.35350000000000
   4.93599500000000   4.93599500000000  -2.35350000000000
     Sb     Si     V
      4      2     8
Direct
   0.64094000000000   0.14094000000000   0.78188000000000   Sb   (8h)
   0.35906000000000  -0.14094000000000   0.21812000000000   Sb   (8h)
   0.14094000000000   0.35906000000000   0.50000000000000   Sb   (8h)
  -0.14094000000000   0.64094000000000   0.50000000000000   Sb   (8h)
   0.25000000000000   0.25000000000000   0.00000000000000   Si   (4a)
   0.75000000000000   0.75000000000000   0.00000000000000   Si   (4a)
   0.20674000000000   0.08520000000000   0.29194000000000   V   (16k)
  -0.20674000000000  -0.08520000000000  -0.29194000000000   V   (16k)
   0.08520000000000  -0.20674000000000  -0.12154000000000   V   (16k)
  -0.08520000000000   0.20674000000000   0.12154000000000   V   (16k)
   0.70674000000000   0.41480000000000   0.12154000000000   V   (16k)
   0.29326000000000   0.58520000000000  -0.12154000000000   V   (16k)
   0.58520000000000   0.70674000000000   0.29194000000000   V   (16k)
   0.41480000000000   0.29326000000000  -0.29194000000000   V   (16k)
```

KHF_2 ($F5_2$): A2BC_tI16_140_h_d_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'F2HK'
_chemical_formula_sum 'F2 H K'

loop_
_publ_author_name
 'J. A. Ibers'
_journal_name_full_name
```

```
;
 Journal of Chemical Physics
;
_journal_volume 40
_journal_year 1964
_journal_page_first 402
_journal_page_last 404
_publ_Section_title
;
 Refinement of Peterson and Levy\'s Neutron Diffraction Data on KHF$_{2}
       ↪ $
;

_aflow_title 'KHF$_{2}$ ($F5_{2}$) Structure'
_aflow_proto 'A2BC_tI16_140_h_d_a'
_aflow_params 'a,c/a,x_{3}'
_aflow_params_values '5.67,1.20105820106,0.142'
_aflow_Strukturbericht '$F5_{2}$'
_aflow_Pearson 'tI16'

_symmetry_space_group_name_H-M "I 4/m 2/c 2/m"
_symmetry_Int_Tables_number 140

_cell_length_a     5.67000
_cell_length_b     5.67000
_cell_length_c     6.81000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z+1/2
3 -x,y,-z+1/2
4 -x,-y,z
5 -y,-x,-z+1/2
6 -y,x,z
7 y,-x,z
8 y,x,-z+1/2
9 -x,-y,-z
10 -x,y,z+1/2
11 x,-y,z+1/2
12 x,y,-z
13 y,x,z+1/2
14 y,-x,-z
15 -y,x,-z
16 -y,-x,z+1/2
17 x+1/2,y+1/2,z+1/2
18 x+1/2,-y+1/2,-z
19 -x+1/2,y+1/2,-z
20 -x+1/2,-y+1/2,z+1/2
21 -y+1/2,-x+1/2,-z
22 -y+1/2,x+1/2,z+1/2
23 y+1/2,-x+1/2,z+1/2
24 y+1/2,x+1/2,-z
25 -x+1/2,-y+1/2,-z+1/2
26 -x+1/2,y+1/2,z
27 x+1/2,-y+1/2,z
28 x+1/2,y+1/2,-z+1/2
29 y+1/2,x+1/2,z
30 y+1/2,-x+1/2,-z+1/2
31 -y+1/2,x+1/2,-z+1/2
32 -y+1/2,-x+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
K1 K 4 a 0.00000 0.00000 0.25000 1.00000
H1 H 4 d 0.00000 0.50000 0.00000 1.00000
F1 F 8 h 0.14200 0.64200 0.00000 1.00000
```

KHF_2 ($F5_2$): A2BC_tI16_140_h_d_a - POSCAR

```
A2BC_tI16_140_h_d_a & a,c/a,x3 --params=5.67,1.20105820106,0.142 & I4/
      ↪ mcm D_{4h}^{18} #140 (adh) & tI16 & $F5_{2}$ & F2HK & F2HK & J.
      ↪  A. Ibers, J. Chem. Phys. 40, 402-404 (1964)
   1.00000000000000
  -2.83500000000000    2.83500000000000    3.40500000000000
   2.83500000000000   -2.83500000000000    3.40500000000000
   2.83500000000000    2.83500000000000   -3.40500000000000
     F     H     K
     4     2     2
Direct
   0.64200000000000    0.14200000000000    0.78400000000000    F    (8h)
   0.35800000000000   -0.14200000000000    0.21600000000000    F    (8h)
   0.14200000000000    0.35800000000000    0.50000000000000    F    (8h)
  -0.14200000000000    0.64200000000000    0.50000000000000    F    (8h)
   0.50000000000000    0.00000000000000    0.50000000000000    H    (4d)
   0.00000000000000    0.50000000000000    0.50000000000000    H    (4d)
   0.25000000000000    0.25000000000000    0.00000000000000    K    (4a)
   0.75000000000000    0.75000000000000    0.00000000000000    K    (4a)
```

$Pu_{31}Rh_{20}$: A31B20_tI204_140_b2gh3m_ac2fh3l - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Pu31Rh20'
_chemical_formula_sum 'Pu31 Rh20'

loop_
_publ_author_name
 'D. T. Cromer'
 'A. C. Larson'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 33
_journal_year 1977
_journal_page_first 2620
_journal_page_last 2627
_publ_Section_title
;
 The Crystal Structure of Pu$_{31}$Pt$_{20}$ and Pu$_{31}$Rh$_{20}$
;

_aflow_title 'Pu$_{31}$Rh$_{20}$ Structure'
_aflow_proto 'A31B20_tI204_140_b2gh3m_ac2fh3l'
_aflow_params 'a,c/a,z_{4},z_{5},z_{6},z_{7},x_{8},x_{9},x_{10},z_{10},
      ↪ x_{11},z_{11},x_{12},z_{12},x_{13},y_{13},z_{13},x_{14},y_{14},
      ↪ z_{14},x_{15},y_{15},z_{15}'
_aflow_params_values '11.076,3.334507945 11,0.094,0.1734,0.0756,0.1656,
      ↪ 0.1586,0.4035,0.1812,0.0726,0.3417,0.1269,0.1536,0.2109,0.2947,
      ↪ 0.4299,0.05,0.2855,0.5774,0.1346,0.2819,0.4125,0.2114'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tI204'

_symmetry_space_group_name_H-M "I 4/m 2/c 2/m"
_symmetry_Int_Tables_number 140

_cell_length_a     11.07600
_cell_length_b     11.07600
_cell_length_c     36.93301
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z+1/2
3 -x,y,-z+1/2
4 -x,-y,z
5 -y,-x,-z+1/2
6 -y,x,z
7 y,-x,z
8 y,x,-z+1/2
9 -x,-y,-z
10 -x,y,z+1/2
11 x,-y,z+1/2
12 x,y,-z
13 y,x,z+1/2
14 y,-x,-z
15 -y,x,-z
16 -y,-x,z+1/2
17 x+1/2,y+1/2,z+1/2
18 x+1/2,-y+1/2,-z
19 -x+1/2,y+1/2,-z
20 -x+1/2,-y+1/2,z+1/2
21 -y+1/2,-x+1/2,-z
22 -y+1/2,x+1/2,z+1/2
23 y+1/2,-x+1/2,z+1/2
24 y+1/2,x+1/2,-z
25 -x+1/2,-y+1/2,-z+1/2
26 -x+1/2,y+1/2,z
27 x+1/2,-y+1/2,z
28 x+1/2,y+1/2,-z+1/2
29 y+1/2,x+1/2,z
30 y+1/2,-x+1/2,-z+1/2
31 -y+1/2,x+1/2,-z+1/2
32 -y+1/2,-x+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Rh1 Rh 4 a 0.00000 0.00000 0.25000 1.00000
Pu1 Pu 4 b 0.00000 0.50000 0.25000 1.00000
Rh2 Rh 4 c 0.00000 0.00000 0.00000 1.00000
Rh3 Rh 8 f 0.00000 0.00000 0.09400 1.00000
Rh4 Rh 8 f 0.00000 0.00000 0.17340 1.00000
Pu2 Pu 8 g 0.00000 0.50000 0.07560 1.00000
Pu3 Pu 8 g 0.00000 0.50000 0.16560 1.00000
Pu4 Pu 8 h 0.15860 0.65860 0.00000 1.00000
Rh5 Rh 8 h 0.40350 0.90350 0.00000 1.00000
Rh6 Rh 16 l 0.18120 0.68120 0.07260 1.00000
Rh7 Rh 16 l 0.34170 0.84170 0.12690 1.00000
Rh8 Rh 16 l 0.15360 0.65360 0.21090 1.00000
Pu5 Pu 32 m 0.29470 0.42990 0.05000 1.00000
Pu6 Pu 32 m 0.28550 0.57740 0.13460 1.00000
Pu7 Pu 32 m 0.28190 0.41250 0.21140 1.00000
```

$Pu_{31}Rh_{20}$: A31B20_tI204_140_b2gh3m_ac2fh3l - POSCAR

```
A31B20_tI204_140_b2gh3m_ac2fh3l & a,c/a,z4,z5,z6,z7,x8,x9,x10,z10,x11,
    ↪ z11,x12,z12,x13,y13,z13,x14,y14,z14,x15,y15,z15 --params=11.076
    ↪ ,3.33450794511,0.094,0.1734,0.0756,0.1656,0.1586,0.4035,0.1812,
    ↪ 0.0726,0.3417,0.1269,0.1536,0.2109,0.2947,0.4299,0.05,0.2855,
    ↪ 0.5774,0.1346,0.2819,0.4125,0.2114 & I4/mcm D_{4h}^{18} #140 (
    ↪ abcf^2g^2h^2l^3m^3) & tI204 & None & Pu31Rh20 & Pu31Rh20 & D.
    ↪ T. Cromer and A. C. Larson, Acta Crystallogr. Sect. B Struct.
    ↪ Sci. 33, 2620-2627 (1977)
  1.00000000000000
 -5.53800000000000   5.53800000000000  18.46650500000000
  5.53800000000000  -5.53800000000000  18.46650500000000
  5.53800000000000   5.53800000000000 -18.46650500000000
   Pu   Rh
   62   40
Direct
  0.75000000000000  0.25000000000000  0.50000000000000  Pu  (4b)
  0.25000000000000  0.75000000000000  0.50000000000000  Pu  (4b)
  0.57560000000000  0.07560000000000  0.50000000000000  Pu  (8g)
  0.07560000000000  0.57560000000000  0.50000000000000  Pu  (8g)
 -0.07560000000000  0.42440000000000  0.50000000000000  Pu  (8g)
  0.42440000000000 -0.07560000000000  0.50000000000000  Pu  (8g)
  0.66560000000000  0.16560000000000  0.50000000000000  Pu  (8g)
  0.16560000000000  0.66560000000000  0.50000000000000  Pu  (8g)
 -0.16560000000000  0.33440000000000  0.50000000000000  Pu  (8g)
  0.33440000000000 -0.16560000000000  0.50000000000000  Pu  (8g)
  0.65860000000000  0.15860000000000  0.81720000000000  Pu  (8h)
  0.34140000000000 -0.15860000000000  0.18280000000000  Pu  (8h)
  0.15860000000000  0.34140000000000  0.50000000000000  Pu  (8h)
 -0.15860000000000  0.65860000000000  0.50000000000000  Pu  (8h)
  0.47990000000000  0.34470000000000  0.72460000000000  Pu (32m)
 -0.37990000000000 -0.24470000000000 -0.72460000000000  Pu (32m)
  0.34470000000000 -0.37990000000000 -0.13520000000000  Pu (32m)
 -0.24470000000000  0.47990000000000  0.13520000000000  Pu (32m)
  0.87990000000000  0.15530000000000  0.13520000000000  Pu (32m)
  0.02010000000000  0.74470000000000 -0.13520000000000  Pu (32m)
  0.74470000000000  0.87990000000000  0.72460000000000  Pu (32m)
  0.15530000000000  0.02010000000000 -0.72460000000000  Pu (32m)
 -0.47990000000000 -0.34470000000000 -0.72460000000000  Pu (32m)
  0.37990000000000  0.24470000000000  0.72460000000000  Pu (32m)
 -0.34470000000000  0.37990000000000  0.13520000000000  Pu (32m)
  0.24470000000000 -0.47990000000000 -0.13520000000000  Pu (32m)
  0.12010000000000  0.84470000000000 -0.13520000000000  Pu (32m)
  0.97990000000000  0.25530000000000  0.13520000000000  Pu (32m)
  0.25530000000000  0.12010000000000 -0.72460000000000  Pu (32m)
  0.84470000000000  0.97990000000000  0.72460000000000  Pu (32m)
  0.71200000000000  0.42010000000000  0.86290000000000  Pu (32m)
 -0.44280000000000 -0.15090000000000 -0.86290000000000  Pu (32m)
  0.42010000000000 -0.44280000000000 -0.29190000000000  Pu (32m)
 -0.15090000000000  0.71200000000000  0.29190000000000  Pu (32m)
  0.94280000000000  0.07990000000000  0.29190000000000  Pu (32m)
 -0.21200000000000  0.65090000000000 -0.29190000000000  Pu (32m)
  0.65090000000000  0.94280000000000  0.86290000000000  Pu (32m)
  0.07990000000000 -0.21200000000000 -0.86290000000000  Pu (32m)
 -0.71200000000000 -0.42010000000000 -0.86290000000000  Pu (32m)
  0.44280000000000  0.15090000000000  0.86290000000000  Pu (32m)
 -0.42010000000000  0.44280000000000  0.29190000000000  Pu (32m)
  0.15090000000000 -0.71200000000000 -0.29190000000000  Pu (32m)
  0.05720000000000  0.92010000000000 -0.29190000000000  Pu (32m)
  1.21200000000000  0.34910000000000  0.29190000000000  Pu (32m)
  0.34910000000000  0.05720000000000 -0.86290000000000  Pu (32m)
  0.92010000000000  1.21200000000000  0.86290000000000  Pu (32m)
  0.62390000000000  0.49330000000000  0.69440000000000  Pu (32m)
 -0.20110000000000 -0.07050000000000 -0.69440000000000  Pu (32m)
  0.49330000000000 -0.20110000000000 -0.13060000000000  Pu (32m)
 -0.07050000000000  0.62390000000000  0.13060000000000  Pu (32m)
  0.70110000000000  0.00670000000000  0.13060000000000  Pu (32m)
 -0.12390000000000  0.57050000000000 -0.13060000000000  Pu (32m)
  0.57050000000000  0.70110000000000  0.69440000000000  Pu (32m)
  0.00670000000000 -0.12390000000000 -0.69440000000000  Pu (32m)
 -0.62390000000000 -0.49330000000000 -0.69440000000000  Pu (32m)
  0.20110000000000  0.07050000000000  0.69440000000000  Pu (32m)
 -0.49330000000000  0.20110000000000  0.13060000000000  Pu (32m)
  0.07050000000000 -0.62390000000000 -0.13060000000000  Pu (32m)
  0.29890000000000  0.99330000000000 -0.13060000000000  Pu (32m)
  1.12390000000000  0.42950000000000  0.13060000000000  Pu (32m)
  0.42950000000000  0.29890000000000 -0.69440000000000  Pu (32m)
  0.99330000000000  1.12390000000000  0.69440000000000  Pu (32m)
  0.25000000000000  0.25000000000000  0.00000000000000  Rh  (4a)
  0.75000000000000  0.75000000000000  0.00000000000000  Rh  (4a)
  0.00000000000000  0.00000000000000  0.00000000000000  Rh  (4c)
  0.50000000000000  0.50000000000000  0.00000000000000  Rh  (4c)
  0.09400000000000  0.09400000000000  0.00000000000000  Rh  (8f)
  0.40600000000000  0.40600000000000  0.00000000000000  Rh  (8f)
 -0.09400000000000 -0.09400000000000  0.00000000000000  Rh  (8f)
  0.59400000000000  0.59400000000000  0.00000000000000  Rh  (8f)
  0.17340000000000  0.17340000000000  0.00000000000000  Rh  (8f)
  0.32660000000000  0.32660000000000  0.00000000000000  Rh  (8f)
 -0.17340000000000 -0.17340000000000  0.00000000000000  Rh  (8f)
  0.67340000000000  0.67340000000000  0.00000000000000  Rh  (8f)
  0.90350000000000  0.40350000000000  1.30700000000000  Rh  (8h)
  0.09650000000000 -0.40350000000000 -0.30700000000000  Rh  (8h)
  0.40350000000000  0.09650000000000  0.50000000000000  Rh  (8h)
 -0.40350000000000  0.90350000000000  0.50000000000000  Rh  (8h)
  0.75380000000000  0.25380000000000  0.86240000000000  Rh (16l)
  0.39140000000000 -0.10860000000000  0.13760000000000  Rh (16l)
  0.25380000000000  0.39140000000000  0.50000000000000  Rh (16l)
 -0.10860000000000  0.75380000000000  0.50000000000000  Rh (16l)
  0.10860000000000  0.24620000000000  0.50000000000000  Rh (16l)
 -0.25380000000000  0.60860000000000  0.50000000000000  Rh (16l)
  0.60860000000000  0.10860000000000  0.86240000000000  Rh (16l)
  0.24620000000000 -0.25380000000000  0.13760000000000  Rh (16l)
  0.96860000000000  0.46860000000000  1.18340000000000  Rh (16l)
  0.28520000000000 -0.21480000000000 -0.18340000000000  Rh (16l)
  0.46860000000000  0.28520000000000  0.50000000000000  Rh (16l)
 -0.21480000000000  0.96860000000000  0.50000000000000  Rh (16l)
  0.21480000000000  0.03140000000000  0.50000000000000  Rh (16l)
 -0.46860000000000  0.71480000000000  0.50000000000000  Rh (16l)
  0.71480000000000  0.21480000000000  1.18340000000000  Rh (16l)
  0.03140000000000 -0.46860000000000 -0.18340000000000  Rh (16l)
  0.86450000000000  0.36450000000000  0.80720000000000  Rh (16l)
  0.55730000000000  0.05730000000000  0.19280000000000  Rh (16l)
  0.36450000000000  0.55730000000000  0.50000000000000  Rh (16l)
  0.05730000000000  0.86450000000000  0.50000000000000  Rh (16l)
 -0.05730000000000  0.13550000000000  0.50000000000000  Rh (16l)
 -0.36450000000000  0.44270000000000  0.50000000000000  Rh (16l)
  0.44270000000000 -0.05730000000000  0.80720000000000  Rh (16l)
  0.13550000000000 -0.36450000000000  0.19280000000000  Rh (16l)
```

$NH_4Pb_2Br_5$ ($K3_4$): A5BC2_tI32_140_bl_a_h - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Br5(NH4)Pb2'
_chemical_formula_sum 'Br5 (NH4) Pb2'

loop_
_publ_author_name
 'H. M. Powell'
 'H. S. Tasker'
_journal_year 1937
_publ_Section_title
;
 The valency angle of bivalent lead: the crystal structure of ammonium,
    ↪ rubidium, and potassium pentabromodiplumbites
;

# Found in Strukturbericht Band V 1937, 1940

_aflow_title 'NH$_{4}$Pb$_{2}$Br$_{5}$ ($K3_{4}$) Structure'
_aflow_proto 'A5BC2_tI32_140_bl_a_h'
_aflow_params 'a,c/a,x_{3},x_{4},z_{4}'
_aflow_params_values '8.39,1.70917759237,0.158,0.163,0.363'
_aflow_Strukturbericht '$K3_{4}$'
_aflow_Pearson 'tI32'

_symmetry_space_group_name_H-M "I 4/m 2/c 2/m"
_symmetry_Int_Tables_number 140

_cell_length_a      8.39000
_cell_length_b      8.39000
_cell_length_c      14.34000
_cell_angle_alpha   90.00000
_cell_angle_beta    90.00000
_cell_angle_gamma   90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z+1/2
3 -x,y,-z+1/2
4 -x,-y,z
5 -y,-x,-z+1/2
6 -y,x,z
7 y,-x,z
8 y,x,-z+1/2
9 -x,-y,-z
10 -x,y,z+1/2
11 x,-y,z+1/2
12 x,y,-z
13 y,x,z+1/2
14 y,-x,-z
15 -y,x,-z
16 -y,-x,z+1/2
17 x+1/2,y+1/2,z+1/2
18 x+1/2,-y+1/2,-z
19 -x+1/2,y+1/2,-z
20 -x+1/2,-y+1/2,z+1/2
21 -y+1/2,-x+1/2,-z
22 -y+1/2,x+1/2,z+1/2
23 y+1/2,-x+1/2,z+1/2
24 y+1/2,x+1/2,-z
25 -x+1/2,-y+1/2,-z+1/2
26 -x+1/2,y+1/2,z
27 x+1/2,-y+1/2,z
28 x+1/2,y+1/2,-z+1/2
29 y+1/2,x+1/2,z
30 y+1/2,-x+1/2,-z+1/2
31 -y+1/2,x+1/2,-z+1/2
32 -y+1/2,-x+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
NH41 NH4 4 a 0.00000 0.00000 0.25000 1.00000
Br1 Br 4 b 0.00000 0.50000 0.25000 1.00000
Pb1 Pb 8 h 0.15800 0.65800 0.00000 1.00000
Br2 Br 16 l 0.16300 0.66300 0.36300 1.00000
```

$NH_4Pb_2Br_5$ ($K3_4$): A5BC2_tI32_140_bl_a_h - POSCAR

```
A5BC2_tI32_140_bl_a_h & a,c/a,x3,x4,z4 --params=8.39,1.70917759237,0.158
    ↪ ,0.163,0.363 & I4/mcm D_{4h}^{18} #140 (abh1) & tI32 & $K3_{4}$
    ↪ & Br5(NH4)Pb2 & Br5(NH4)Pb2 & H. M. Powell and H. S. Tasker, (
    ↪ 1937)
   1.00000000000000
  -4.19500000000000    4.19500000000000    7.17000000000000
   4.19500000000000   -4.19500000000000    7.17000000000000
   4.19500000000000    4.19500000000000   -7.17000000000000
     Br    NH4     Pb
     10      2      4
Direct
   0.75000000000000    0.25000000000000    0.50000000000000    Br    (4b)
   0.25000000000000    0.75000000000000    0.50000000000000    Br    (4b)
   1.02600000000000    0.52600000000000    0.82600000000000    Br    (16l)
   0.70000000000000    0.20000000000000    0.17400000000000    Br    (16l)
   0.52600000000000    0.70000000000000    0.50000000000000    Br    (16l)
   0.20000000000000    1.02600000000000    0.50000000000000    Br    (16l)
  -0.20000000000000   -0.02600000000000    0.50000000000000    Br    (16l)
  -0.52600000000000    0.30000000000000    0.50000000000000    Br    (16l)
   0.30000000000000   -0.20000000000000    0.82600000000000    Br    (16l)
  -0.02600000000000   -0.52600000000000    0.17400000000000    Br    (16l)
   0.25000000000000    0.25000000000000    0.00000000000000    NH4    (4a)
   0.75000000000000    0.75000000000000    0.00000000000000    NH4    (4a)
   0.65800000000000    0.15800000000000    0.81600000000000    Pb    (8h)
   0.34200000000000   -0.15800000000000    0.18400000000000    Pb    (8h)
   0.15800000000000    0.34200000000000    0.50000000000000    Pb    (8h)
  -0.15800000000000    0.65800000000000    0.50000000000000    Pb    (8h)
```

Cs_3CoCl_5 ($K3_1$): A5BC3_tI36_140_cl_b_ah - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Cl5CoCs3'
_chemical_formula_sum 'Cl5 Co Cs3'

loop_
_publ_author_name
 'B. N. Figgis'
 'R. Mason'
 'A. R. P. Smith'
 'G. A. Williams'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 36
_journal_year 1980
_journal_page_first 509
_journal_page_last 512
_publ_Section_title
;
 Neutron Diffraction Structure of Cs$_{3}$CoCl$_{5}$ at 4.2~K
;

_aflow_title 'Cs$_{3}$CoCl$_{5}$ ($K3_{1}$) Structure'
_aflow_proto 'A5BC3_tI36_140_cl_b_ah'
_aflow_params 'a,c/a,x_{4},x_{5},z_{5}'
_aflow_params_values '9.063,1.59439479201,0.66225,0.1421,0.15711'
_aflow_Strukturbericht '$K3_{1}$'
_aflow_Pearson 'tI36'

_symmetry_space_group_name_H-M "I 4/m 2/c 2/m"
_symmetry_Int_Tables_number 140

_cell_length_a     9.06300
_cell_length_b     9.06300
_cell_length_c     14.45000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z+1/2
3 -x,y,-z+1/2
4 -x,-y,z
5 -y,-x,-z+1/2
6 -y,x,z
7 y,-x,z
8 y,x,-z+1/2
9 -x,-y,-z
10 -x,y,z+1/2
11 x,-y,z+1/2
12 x,y,-z
13 y,x,z+1/2
14 y,-x,-z
15 -y,x,-z
16 -y,-x,z+1/2
17 x+1/2,y+1/2,z+1/2
18 x+1/2,-y+1/2,-z
19 -x+1/2,y+1/2,-z
20 -x+1/2,-y+1/2,z+1/2
21 -y+1/2,-x+1/2,-z
22 -y+1/2,x+1/2,z+1/2
23 y+1/2,-x+1/2,z+1/2
24 y+1/2,x+1/2,-z
25 -x+1/2,-y+1/2,-z+1/2
26 -x+1/2,y+1/2,z
27 x+1/2,-y+1/2,z
28 x+1/2,y+1/2,-z+1/2
29 y+1/2,x+1/2,z
30 y+1/2,-x+1/2,-z+1/2
31 -y+1/2,x+1/2,-z+1/2
32 -y+1/2,-x+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cs1 Cs 4 a 0.00000 0.00000 0.25000 1.00000
Co1 Co 4 b 0.00000 0.50000 0.25000 1.00000
Cl1 Cl 4 c 0.00000 0.00000 0.00000 1.00000
Cs2 Cs 8 h 0.66225 0.16225 0.00000 1.00000
Cl2 Cl 16 l 0.14210 0.64210 0.15711 1.00000
```

Cs_3CoCl_5 ($K3_1$): A5BC3_tI36_140_cl_b_ah - POSCAR

```
A5BC3_tI36_140_cl_b_ah & a,c/a,x4,x5,z5 --params=9.063,1.59439479201,
    ↪ 0.66225,0.1421,0.15711 & I4/mcm D_{4h}^{18} #140 (abch1) & tI36
    ↪ & $K3_{1}$ & Cl5CoCs3 & Cl5CoCs3 & B. N. Figgis et al., Acta
    ↪ Crystallogr. Sect. B Struct. Sci. 36, 509-512 (1980)
   1.00000000000000
  -4.53150000000000    4.53150000000000    7.22500000000000
   4.53150000000000   -4.53150000000000    7.22500000000000
   4.53150000000000    4.53150000000000   -7.22500000000000
     Cl     Co     Cs
     10      2      6
Direct
   0.00000000000000    0.00000000000000    0.00000000000000    Cl    (4c)
   0.50000000000000    0.50000000000000    0.00000000000000    Cl    (4c)
   0.79921000000000    0.29921000000000    0.78420000000000    Cl    (16l)
   0.51501000000000    0.01501000000000    0.21580000000000    Cl    (16l)
   0.29921000000000    0.51501000000000    0.50000000000000    Cl    (16l)
   0.01501000000000    0.79921000000000    0.50000000000000    Cl    (16l)
  -0.01501000000000    0.20079000000000    0.50000000000000    Cl    (16l)
  -0.29921000000000    0.48499000000000    0.50000000000000    Cl    (16l)
   0.48499000000000   -0.01501000000000    0.78420000000000    Cl    (16l)
   0.20079000000000   -0.29921000000000    0.21580000000000    Cl    (16l)
   0.75000000000000    0.25000000000000    0.50000000000000    Co     (4b)
   0.25000000000000    0.75000000000000    0.50000000000000    Co     (4b)
   0.25000000000000    0.25000000000000    0.00000000000000    Cs     (4a)
   0.75000000000000    0.75000000000000    0.00000000000000    Cs     (4a)
   1.16225000000000    0.66225000000000    1.82450000000000    Cs     (8h)
  -0.16225000000000   -0.66225000000000   -0.82450000000000    Cs     (8h)
   0.66225000000000   -0.16225000000000    0.50000000000000    Cs     (8h)
  -0.66225000000000    1.16225000000000    0.50000000000000    Cs     (8h)
```

U_6Mn ($D2_c$): AB6_tI28_140_a_hk - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'MnU6'
_chemical_formula_sum 'Mn U6'

loop_
_publ_author_name
 'N. C. Baenziger'
 'R. E. Rundle'
 'A. I. Snow'
 'A. S. Wilson'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 3
_journal_year 1950
_journal_page_first 34
_journal_page_last 40
_publ_Section_title
;
 Compounds of uranium with the transition metals of the first long
    ↪ period
;

# Found in PAULING FILE, {in: Inorganic Solid Phases, SpringerMaterials
    ↪ (online database), Springer, Heidelberg}, 2016

_aflow_title 'U$_{6}$Mn ($D2_{c}$) Structure'
_aflow_proto 'AB6_tI28_140_a_hk'
_aflow_params 'a,c/a,x_{2},x_{3},y_{3}'
_aflow_params_values '10.29,0.509232264334,0.4068,0.2141,0.1021'
_aflow_Strukturbericht '$D2_{c}$'
_aflow_Pearson 'tI28'

_symmetry_space_group_name_H-M "I 4/m 2/c 2/m"
_symmetry_Int_Tables_number 140

_cell_length_a     10.29000
_cell_length_b     10.29000
_cell_length_c     5.24000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z+1/2
```

```
3 -x,y,-z+1/2
4 -x,-y,z
5 -y,-x,-z+1/2
6 -y,x,z
7 y,-x,z
8 y,x,-z+1/2
9 -x,-y,-z
10 -x,y,z+1/2
11 x,-y,z+1/2
12 x,y,-z
13 y,x,z+1/2
14 y,-x,-z
15 -y,x,-z
16 -y,-x,z+1/2
17 x+1/2,y+1/2,z+1/2
18 x+1/2,-y+1/2,-z
19 -x+1/2,y+1/2,-z
20 -x+1/2,-y+1/2,z+1/2
21 -y+1/2,-x+1/2,-z
22 -y+1/2,x+1/2,z+1/2
23 y+1/2,-x+1/2,z+1/2
24 y+1/2,x+1/2,-z
25 -x+1/2,-y+1/2,-z+1/2
26 -x+1/2,y+1/2,z
27 x+1/2,-y+1/2,z
28 x+1/2,y+1/2,-z+1/2
29 y+1/2,x+1/2,z
30 y+1/2,-x+1/2,-z+1/2
31 -y+1/2,x+1/2,-z+1/2
32 -y+1/2,-x+1/2,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Mn1 Mn 4 a 0.00000 0.00000 0.25000 1.00000
U1 U 8 h 0.40680 -0.09320 0.00000 1.00000
U2 U 16 k 0.21410 0.10210 0.00000 1.00000
```

U_6Mn ($D2_c$): AB6_tI28_140_a_hk - POSCAR

```
AB6_tI28_140_a_hk & a,c/a,x2,x3,y3 --params=10.29,0.509232264334,0.4068,
    ↪ 0.2141,0.1021 & I4/mcm D_{4h}^{18} #140 (ahk) & tI28 & $D2_{c}$
    ↪ & MnU6 & MnU6 & N. C. Baenziger et al., Acta Cryst. 3, 34-40 (
    ↪ 1950)
   1.00000000000000
  -5.14500000000000    5.14500000000000    2.62000000000000
   5.14500000000000   -5.14500000000000    2.62000000000000
   5.14500000000000    5.14500000000000   -2.62000000000000
    Mn   U
     2  12
Direct
   0.25000000000000    0.25000000000000    0.00000000000000  Mn   (4a)
   0.75000000000000    0.75000000000000    0.00000000000000  Mn   (4a)
   0.90680000000000    0.40680000000000    1.31360000000000  U    (8h)
   0.09320000000000   -0.40680000000000   -0.31360000000000  U    (8h)
   0.40680000000000    0.09320000000000    0.50000000000000  U    (8h)
  -0.40680000000000    0.90680000000000    0.50000000000000  U    (8h)
   0.10210000000000    0.21410000000000    0.31620000000000  U   (16k)
  -0.10210000000000   -0.21410000000000   -0.31620000000000  U   (16k)
   0.21410000000000   -0.10210000000000    0.11200000000000  U   (16k)
  -0.21410000000000    0.10210000000000   -0.11200000000000  U   (16k)
   0.60210000000000    0.28590000000000   -0.11200000000000  U   (16k)
   0.39790000000000    0.71410000000000    0.11200000000000  U   (16k)
   0.71410000000000    0.60210000000000    0.31620000000000  U   (16k)
   0.28590000000000    0.39790000000000   -0.31620000000000  U   (16k)
```

$BaCd_{11}$: AB11_tI48_141_a_bdi - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'BaCd11'
_chemical_formula_sum 'Ba Cd11'

loop_
_publ_author_name
 'M. J. Sanderson'
 'N. C. Baenziger'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 6
_journal_year 1953
_journal_page_first 627
_journal_page_last 631
_publ_Section_title
;
 The Crystal Structure of BaCd$_{11}$
;

_aflow_title 'BaCd$_{11}$ Structure'
_aflow_proto 'AB11_tI48_141_a_bdi'
_aflow_params 'a,c/a,x_{4},y_{4},z_{4}'
_aflow_params_values '12.02,0.643926788686,0.123,0.455,0.183'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'tI48'
_symmetry_space_group_name_H-M "I 41/a 2/m 2/d (origin choice 2)"
_symmetry_Int_Tables_number 141

_cell_length_a      12.02000
_cell_length_b      12.02000
_cell_length_c      7.74000
_cell_angle_alpha   90.00000
_cell_angle_beta    90.00000
_cell_angle_gamma   90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y+1/2,-z
4 -x,-y+1/2,z
5 -y+1/4,-x+1/4,-z+3/4
6 -y+1/4,x+3/4,z+1/4
7 y+3/4,-x+3/4,z+1/4
8 y+3/4,x+1/4,-z+3/4
9 -x,-y,-z
10 -x,y,z
11 x,-y+1/2,z
12 x,y+1/2,-z
13 y+3/4,x+3/4,z+1/4
14 y+3/4,-x+1/4,-z+3/4
15 -y+1/4,x+1/4,-z+3/4
16 -y+1/4,-x+3/4,z+1/4
17 x+1/2,y+1/2,z+1/2
18 x+1/2,-y+1/2,-z+1/2
19 -x+1/2,y,-z+1/2
20 -x+1/2,-y,z+1/2
21 -y+3/4,-x+3/4,-z+1/4
22 -y+3/4,x+1/4,z+3/4
23 y+1/4,-x+1/4,z+3/4
24 y+1/4,x+3/4,-z+1/4
25 -x+1/2,-y+1/2,-z+1/2
26 -x+1/2,y+1/2,z+1/2
27 x+1/2,-y,z+1/2
28 x+1/2,y,-z+1/2
29 y+1/4,x+1/4,z+3/4
30 y+1/4,-x+3/4,-z+1/4
31 -y+3/4,x+3/4,-z+1/4
32 -y+3/4,-x+1/4,z+3/4

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ba1 Ba 4 a 0.00000 0.75000 0.12500 1.00000
Cd1 Cd 4 b 0.00000 0.25000 0.37500 1.00000
Cd2 Cd 8 d 0.00000 0.00000 0.50000 1.00000
Cd3 Cd 32 i 0.12300 0.45500 0.18300 1.00000
```

$BaCd_{11}$: AB11_tI48_141_a_bdi - POSCAR

```
AB11_tI48_141_a_bdi & a,c/a,x4,y4,z4 --params=12.02,0.643926788686,0.123
    ↪ ,0.455,0.183 & I4_{1}/amd D_{4h}^{19} #141 (abdi) & tI48 & None
    ↪ & BaCd11 & BaCd11 & M. J. Sanderson and N. C. Baenziger, Acta
    ↪ Cryst. 6, 627-631 (1953)
   1.00000000000000
  -6.01000000000000    6.01000000000000    3.87000000000000
   6.01000000000000   -6.01000000000000    3.87000000000000
   6.01000000000000    6.01000000000000   -3.87000000000000
    Ba  Cd
     2  22
Direct
   0.87500000000000    0.12500000000000    0.75000000000000  Ba   (4a)
   0.12500000000000    0.87500000000000    0.25000000000000  Ba   (4a)
   0.62500000000000    0.37500000000000    0.25000000000000  Cd   (4b)
   0.37500000000000    0.62500000000000    0.75000000000000  Cd   (4b)
   0.50000000000000    0.50000000000000    0.00000000000000  Cd   (8d)
   0.00000000000000    0.50000000000000    0.50000000000000  Cd   (8d)
   0.50000000000000    0.00000000000000    0.00000000000000  Cd   (8d)
   0.50000000000000    0.50000000000000    0.50000000000000  Cd   (8d)
   0.63800000000000    0.30600000000000    0.57800000000000  Cd  (32i)
   0.22800000000000    0.06000000000000   -0.07800000000000  Cd  (32i)
   0.30600000000000    0.22800000000000   -0.33200000000000  Cd  (32i)
   0.06000000000000    0.63800000000000    0.83200000000000  Cd  (32i)
   0.77200000000000   -0.30600000000000    0.83200000000000  Cd  (32i)
  -0.63800000000000   -0.06000000000000   -0.33200000000000  Cd  (32i)
  -0.06000000000000    0.77200000000000    0.57800000000000  Cd  (32i)
  -0.30600000000000   -0.63800000000000   -0.07800000000000  Cd  (32i)
  -0.63800000000000   -0.30600000000000   -0.57800000000000  Cd  (32i)
   0.77200000000000   -0.06000000000000    1.07800000000000  Cd  (32i)
  -0.30600000000000    0.77200000000000    0.33200000000000  Cd  (32i)
  -0.06000000000000   -0.63800000000000    0.16800000000000  Cd  (32i)
   0.22800000000000    0.30600000000000    0.16800000000000  Cd  (32i)
   0.63800000000000    0.06000000000000    0.33200000000000  Cd  (32i)
   0.06000000000000    0.22800000000000   -0.57800000000000  Cd  (32i)
   0.30600000000000    0.63800000000000    1.07800000000000  Cd  (32i)
```

Analcime ($NaAlSi_2O_6 \cdot H_2O$, $S6_1$): A2B2C3D12E4_tI184_142_f_f_be_3g_g - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Analcime'
```

```
_chemical_formula_sum 'Al2 (H2O)2 Na3 O12 Si4'

loop_
_publ_author_name
 'F. Mazzi'
 'E. Galli'
_journal_name_full_name
;
 American Mineralogist
;
_journal_volume 63
_journal_year 1978
_journal_page_first 448
_journal_page_last 460
_publ_Section_title
;
 Is each analcime different?
;

# Found in The crystal structure of natural monoclinic analcime (
      ↪ NaAlSi$_{2}$O$_{6}$$\cdot$H$_{2}$O), 1988

_aflow_title 'Analcime (NaAlSi$_{2}$O$_{6}$$\cdot$H$_{2}$O, $S6_{1}$)
      ↪ Structure'
_aflow_proto 'A2B2C3D12E4_tI184_142_f_f_be_3g_g'
_aflow_params 'a,c/a,x_{2},x_{3},x_{4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{
      ↪ 6},x_{7},y_{7},z_{7},x_{8},y_{8},z_{8}'
_aflow_params_values '13.723,0.997303796546,0.1242,0.1631,0.3805,0.105,
      ↪ 0.3705,0.2186,0.2218,0.1028,0.3633,0.3631,0.2182,0.1051,0.1264,
      ↪ 0.1617,0.4118'
_aflow_Strukturbericht '$S6_{1}$'
_aflow_Pearson 'tI184'

_symmetry_space_group_name_H-M "I 41/a 2/c 2/d (origin choice 2)"
_symmetry_Int_Tables_number 142

_cell_length_a     13.72300
_cell_length_b     13.72300
_cell_length_c     13.68600
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z
3 -x+1/2,y,-z
4 -x,-y+1/2,z
5 -y+1/4,-x+1/4,-z+1/4
6 -y+1/4,x+3/4,z+1/4
7 y+3/4,-x+3/4,z+1/4
8 y+3/4,x+1/4,-z+1/4
9 -x,-y,-z
10 -x,y,z+1/2
11 x,-y+1/2,z+1/2
12 x,y+1/2,-z
13 y+1/4,x+1/4,z+1/4
14 y+3/4,-x+1/4,-z+3/4
15 -y+1/4,x+1/4,-z+3/4
16 -y+3/4,-x+1/4,z+1/4
17 x+1/2,y+1/2,z+1/2
18 x,-y,-z+1/2
19 -x,y+1/2,-z+1/2
20 -x+1/2,-y,z+1/2
21 -y+3/4,-x+3/4,-z+3/4
22 -y+3/4,x+1/4,z+3/4
23 y+1/4,-x+1/4,z+3/4
24 y+1/4,x+3/4,-z+3/4
25 -x+1/2,-y+1/2,-z+1/2
26 -x+1/2,y+1/2,z
27 x+1/2,-y,z
28 x+1/2,y,-z+1/2
29 y+3/4,x+3/4,z+3/4
30 y+1/4,-x+3/4,-z+1/4
31 -y+3/4,x+3/4,-z+1/4
32 -y+1/4,-x+3/4,z+3/4

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Na1 Na 8 b 0.00000 0.25000 0.12500 0.23000
Na2 Na 16 e 0.12420 0.00000 0.25000 0.82000
Al1 Al 16 f 0.16310 0.41310 0.12500 1.00000
H2O1 H2O 16 f 0.38050 0.63050 0.12500 1.00000
O1 O 32 g 0.10500 0.37050 0.21860 1.00000
O2 O 32 g 0.22180 0.10280 0.36330 1.00000
O3 O 32 g 0.36310 0.21820 0.10510 1.00000
Si1 Si 32 g 0.12640 0.16170 0.41180 1.00000
```

Analcime ($NaAlSi_2O_6 \cdot H_2O$, $S6_1$): A2B2C3D12E4_tI184_142_f_f_be_3g_g - POSCAR

```
A2B2C3D12E4_tI184_142_f_f_be_3g_g & a,c/a,x2,x3,x4,x5,y5,z5,x6,y6,z6,x7,
      ↪ y7,z7,x8,y8,z8 --params=13.723,0.997303796546,0.1242,0.1631,
      ↪ 0.3805,0.105,0.3705,0.2186,0.2218,0.1028,0.3633,0.3631,0.2182,
      ↪ 0.1051,0.1264,0.1617,0.4118 & I4_{1}/acd D_{4h}^{20} #142 (bef^
      ↪ 2g^4) & tI184 & $S6_{1}$ & Al(H2O)NaO6Si2 & Analcime & F. Mazzi
      ↪  and E. Galli, Am. Mineral. 63, 448-460 (1978)
   1.00000000000000
  -6.86150000000000    6.86150000000000    6.84300000000000
   6.86150000000000   -6.86150000000000    6.84300000000000
   6.86150000000000    6.86150000000000   -6.84300000000000
     Al    H2O     Na      O     Si
      8      8     12     48     16
Direct
   0.53810000000000    0.28810000000000    0.57620000000000   Al   (16f)
   0.21190000000000   -0.03810000000000   -0.07620000000000   Al   (16f)
   0.28810000000000    0.21190000000000    0.75000000000000   Al   (16f)
  -0.03810000000000    0.53810000000000    0.75000000000000   Al   (16f)
   0.46190000000000    0.71190000000000    0.42380000000000   Al   (16f)
   0.78810000000000    1.03810000000000    1.07620000000000   Al   (16f)
   0.71190000000000    0.78810000000000    0.25000000000000   Al   (16f)
   1.03810000000000    0.46190000000000    0.25000000000000   Al   (16f)
   0.75550000000000    0.50550000000000    1.01100000000000   H2O  (16f)
  -0.00550000000000   -0.25550000000000   -0.51100000000000   H2O  (16f)
   0.50550000000000   -0.00550000000000    0.75000000000000   H2O  (16f)
  -0.25550000000000    0.75550000000000    0.75000000000000   H2O  (16f)
   0.24450000000000    0.49450000000000   -0.01100000000000   H2O  (16f)
   1.00550000000000    1.25550000000000    1.51100000000000   H2O  (16f)
   0.49450000000000    1.00550000000000    0.25000000000000   H2O  (16f)
   1.25550000000000    0.24450000000000    0.25000000000000   H2O  (16f)
   0.37500000000000    0.12500000000000    0.25000000000000   Na   (8b)
   0.12500000000000    0.37500000000000    0.75000000000000   Na   (8b)
   0.62500000000000    0.87500000000000    0.75000000000000   Na   (8b)
   0.87500000000000    0.62500000000000    0.25000000000000   Na   (8b)
   0.25000000000000    0.37420000000000    0.12420000000000   Na   (16e)
   0.75000000000000    0.12580000000000    0.37580000000000   Na   (16e)
   0.37420000000000    0.75000000000000    0.12420000000000   Na   (16e)
   0.12580000000000    0.25000000000000    0.37580000000000   Na   (16e)
   0.75000000000000    0.62580000000000   -0.12420000000000   Na   (16e)
   0.25000000000000    0.87420000000000    0.62420000000000   Na   (16e)
   0.62580000000000    0.25000000000000   -0.12420000000000   Na   (16e)
   0.87420000000000    0.75000000000000    0.62420000000000   Na   (16e)
   0.58910000000000    0.32360000000000    0.47550000000000   O    (32g)
   0.34810000000000    0.11360000000000    0.02450000000000   O    (32g)
   0.32360000000000    0.34810000000000   -0.26550000000000   O    (32g)
   0.11360000000000    0.58910000000000    0.76550000000000   O    (32g)
   0.15190000000000    0.17640000000000    0.76550000000000   O    (32g)
  -0.08910000000000    0.38640000000000   -0.26550000000000   O    (32g)
   0.38640000000000    0.15190000000000    0.47550000000000   O    (32g)
   0.17640000000000   -0.08910000000000    0.02450000000000   O    (32g)
  -0.58910000000000   -0.32360000000000   -0.47550000000000   O    (32g)
   0.65190000000000   -0.11360000000000    0.97550000000000   O    (32g)
  -0.32360000000000    0.65190000000000    0.26550000000000   O    (32g)
  -0.11360000000000   -0.58910000000000    0.23450000000000   O    (32g)
  -0.15190000000000    0.82360000000000    0.23450000000000   O    (32g)
   1.08910000000000    0.61360000000000    0.26550000000000   O    (32g)
   0.61360000000000   -0.15190000000000   -0.47550000000000   O    (32g)
   0.82360000000000    1.08910000000000    0.97550000000000   O    (32g)
   0.46610000000000    0.58510000000000    0.32460000000000   O    (32g)
   0.76050000000000    0.14150000000000    0.17540000000000   O    (32g)
   0.58510000000000    0.76050000000000    0.11900000000000   O    (32g)
   0.14150000000000    0.46610000000000    0.38100000000000   O    (32g)
  -0.26050000000000   -0.08510000000000    0.38100000000000   O    (32g)
   0.03390000000000    0.35850000000000    0.11900000000000   O    (32g)
   0.35850000000000   -0.26050000000000    0.32460000000000   O    (32g)
  -0.08510000000000    0.03390000000000    0.17540000000000   O    (32g)
  -0.46610000000000   -0.58510000000000   -0.32460000000000   O    (32g)
   0.23950000000000   -0.14150000000000    0.82460000000000   O    (32g)
  -0.58510000000000    0.23950000000000   -0.11900000000000   O    (32g)
  -0.14150000000000   -0.46610000000000    0.61900000000000   O    (32g)
   0.26050000000000    1.08510000000000    0.61900000000000   O    (32g)
   0.96610000000000    0.64150000000000   -0.11900000000000   O    (32g)
   0.64150000000000    0.26050000000000   -0.32460000000000   O    (32g)
   1.08510000000000    0.96610000000000    0.82460000000000   O    (32g)
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   0.11310000000000    0.03180000000000    0.35510000000000   O    (32g)
   0.17670000000000    0.75800000000000    0.14490000000000   O    (32g)
   0.75800000000000    0.11310000000000    0.58130000000000   O    (32g)
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  -0.11310000000000    0.96820000000000    0.64490000000000   O    (32g)
   0.82330000000000    0.24200000000000   -0.14490000000000   O    (32g)
   0.24200000000000   -0.11310000000000   -0.58130000000000   O    (32g)
   0.96820000000000    0.82330000000000    1.08130000000000   O    (32g)
   0.57350000000000    0.53820000000000    0.28810000000000   Si   (32g)
   0.75010000000000    0.28540000000000    0.21190000000000   Si   (32g)
   0.53820000000000    0.75010000000000   -0.03530000000000   Si   (32g)
   0.28540000000000    0.57350000000000    0.53530000000000   Si   (32g)
  -0.25010000000000   -0.03820000000000    0.53530000000000   Si   (32g)
  -0.07350000000000    0.21460000000000   -0.03530000000000   Si   (32g)
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  -0.53820000000000    0.24990000000000    0.03530000000000   Si   (32g)
  -0.28540000000000   -0.57350000000000    0.46470000000000   Si   (32g)
   0.25010000000000    1.03820000000000    0.46470000000000   Si   (32g)
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   0.78540000000000    0.25010000000000   -0.28810000000000   Si   (32g)
   1.03820000000000    1.07350000000000    0.78810000000000   Si   (32g)
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Cd_3As_2: A2B3_tI160_142_deg_3g - CIF

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_chemical_formula_sum 'As2 Cd3'

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 'M. N. Ali'
 'Q. Gibson'
 'S. Jeon'
 'B. B. Zhou'
 'A. Yazdani'
 'R. J. Cava'
_journal_name_full_name
;
 Inorganic Chemistry
;
_journal_volume 53
_journal_year 2014
_journal_page_first 4062
_journal_page_last 4067
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;
 The Crystal and Electronic Structures of Cd$_{3}$As$_{2}$, the
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;

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_aflow_Strukturbericht 'None'
_aflow_Pearson 'tI160'

_symmetry_space_group_name_H-M "I 41/a 2/c 2/d (origin choice 2)"
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3 -x+1/2,y,-z
4 -x,-y+1/2,z
5 -y+1/4,-x+1/4,-z+1/4
6 -y+1/4,x+3/4,z+1/4
7 y+3/4,-x+3/4,z+1/4
8 y+3/4,x+1/4,-z+1/4
9 -x,-y,-z
10 -x,y,z+1/2
11 x,-y+1/2,z+1/2
12 x,y+1/2,-z
13 y+1/4,x+1/4,z+1/4
14 y+3/4,-x+1/4,-z+3/4
15 -y+1/4,x+1/4,-z+3/4
16 -y+3/4,-x+1/4,z+1/4
17 x+1/2,y+1/2,z+1/2
18 x,-y,-z+1/2
19 -x,y+1/2,-z+1/2
20 -x+1/2,-y,z+1/2
21 -y+3/4,-x+3/4,-z+3/4
22 -y+3/4,x+1/4,z+3/4
23 y+1/4,-x+1/4,z+3/4
24 y+1/4,x+3/4,-z+3/4
25 -x+1/2,-y+1/2,-z+1/2
26 -x+1/2,y+1/2,z
27 x+1/2,-y,z
28 x+1/2,y,-z+1/2
29 y+3/4,x+3/4,z+3/4
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_atom_site_Wyckoff_label
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As2 As 16 e 0.73924 0.00000 0.25000 1.00000
As3 As 32 g 0.24597 0.25789 0.12315 1.00000
Cd1 Cd 32 g 0.13951 0.36959 0.05249 1.00000
Cd2 Cd 32 g 0.11169 0.64224 0.07250 1.00000
Cd3 Cd 32 g 0.11879 0.10610 0.06251 1.00000
```

Cd_3As_2: A2B3_tI160_142_deg_3g - POSCAR

```
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      ↪ 0.12315,0.13951,0.36959,0.05249,0.11169,0.64224,0.0725,0.11879,
      ↪ 0.1061,0.06251 & I4_{1}/acd D_{4h}^{20} #142 (deg^4) & tI160 &
      ↪ None & As2Cd3 & As2Cd3 & M. N. Ali et al., Inorg. Chem. 53,
      ↪ 4062-4067 (2014)
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   6.31650000000000  -6.31650000000000  12.71350000000000
   6.31650000000000   6.31650000000000 -12.71350000000000
    As    Cd
    32    48
Direct
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   0.18290000000000  -0.08702000000000  -0.00910000000000   Cd   (32g)
   0.19200000000000   0.18290000000000  -0.23008000000000   Cd   (32g)
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   0.69200000000000   0.92208000000000   1.00910000000000   Cd   (32g)
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La_3BWO_9 (*P*3): AB3C9D_hP28_143_2a_2d_6d_bc - CIF

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 'J. Han'
 'F. Pan'
 'M. S. Molokeev'
 'J. Dai'
 'M. Peng'
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 'W. Zhou'
 'J. Wang'
_journal_name_full_name
;
 ACS Applied Materials and Interfaces
;
_journal_volume 10
_journal_year 2018
_journal_page_first 13660
_journal_page_last 13668
_publ_Section_title
;
 Redefinition of Crystal Structure and Bi$^{3+}$ Yellow Luminescence
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      ↪ Bi$^{3+}$ Phosphor for White Light-Emitting Diodes
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      ↪ 0.418,0.181,0.498,-0.017,0.728,0.198,0.594,0.134,0.502,0.427,
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_aflow_Strukturbericht 'None'
_aflow_Pearson 'hP28'

_symmetry_space_group_name_H-M "P 3"
_symmetry_Int_Tables_number 143

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2 -y,x-y,z
3 -x+y,-x,z

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B2 B 1 a 0.00000 0.00000 0.84000 1.00000
W1 W 1 b 0.33333 0.66667 0.25000 1.00000
W2 W 1 c 0.66667 0.33333 0.77000 1.00000
La1 La 3 d 0.36300 0.08400 0.22900 1.00000
La2 La 3 d 0.27800 0.36500 0.73700 1.00000
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O2 O 3 d 0.13800 0.16900 0.41800 1.00000
O3 O 3 d 0.18100 0.49800 -0.01700 1.00000
O4 O 3 d 0.72800 0.19800 0.59400 1.00000
O5 O 3 d 0.13400 0.50200 0.42700 1.00000
O6 O 3 d 0.61700 0.14300 -0.00400 1.00000
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La_3BWO_9 ($P3$): AB3C9D_hP28_143_2a_2d_6d_bc - POSCAR

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      ↪ BaLi3O9W & BaLi3O9W & J. Han et al., ACS Appl. Mater.
      ↪ Interfaces 10, 13660-13668 (2018)
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       B     La     O     W
       2      6     18     2
Direct
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     0.00000000000000    0.00000000000000    0.84000000000000   B   (1a)
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$RbNO_3$ (IV): AB3C_hP45_144_3a_9a_3a - CIF

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 'B. W. Lucas'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 38
_journal_year 1982
_journal_page_first 2353
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_publ_Section_title
;
 Structure (neutron) of phase IV rubidium nitrate at 298 and 403~K
;

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      ↪ y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{10},y_{10},
      ↪ z_{10},x_{11},y_{11},z_{11},x_{12},y_{12},z_{12},x_{13},y_{13},
      ↪ z_{13},x_{14},y_{14},z_{14},x_{15},y_{15},z_{15}'
_aflow_params_values '10.55,0.708056872038,0.4388,0.5668,0.1053,0.0962,
      ↪ 0.2042,0.5318,0.742,0.2057,0.1168,0.3371,0.5582,0.0094,0.3964,
      ↪ 0.4704,0.228,0.5646,0.6512,0.0664,-0.0011,0.1089,0.6245,0.2309,
      ↪ 0.2495,0.5637,0.0581,0.2465,0.3949,0.6922,0.1208,0.2503,0.8721,
      ↪ 0.2774,0.0897,0.6541,0.2174,0.014,0.4566,0.5691,0.6236,0.1184,
      ↪ 0.2192,0.0,0.7772,0.2214,0.6381'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hP45'

_symmetry_space_group_name_H-M "P 31"
_symmetry_Int_Tables_number 144

_cell_length_a     10.55000
_cell_length_b     10.55000
_cell_length_c     7.47000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z+1/3
3 -x+y,-x,z+2/3

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
N1 N 3 a 0.43880 0.56680 0.10530 1.00000
N2 N 3 a 0.09620 0.20420 0.53180 1.00000
N3 N 3 a 0.74200 0.20570 0.11680 1.00000
O1 O 3 a 0.33710 0.55820 0.00940 1.00000
O2 O 3 a 0.39640 0.47040 0.22800 1.00000
O3 O 3 a 0.56460 0.65120 0.06640 1.00000
O4 O 3 a -0.00110 0.10890 0.62450 1.00000
O5 O 3 a 0.23090 0.24950 0.56370 1.00000
O6 O 3 a 0.05810 0.24650 0.39490 1.00000
O7 O 3 a 0.69220 0.12080 0.25030 1.00000
O8 O 3 a 0.87210 0.27740 0.08970 1.00000
O9 O 3 a 0.65410 0.21740 0.01400 1.00000
Rb1 Rb 3 a 0.45660 0.56910 0.62360 1.00000
Rb2 Rb 3 a 0.11840 0.21920 0.00000 1.00000
Rb3 Rb 3 a 0.77720 0.22140 0.63810 1.00000
```

$RbNO_3$ (IV): AB3C_hP45_144_3a_9a_3a - POSCAR

```
AB3C_hP45_144_3a_9a_3a & a,c/a,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,z4,x5,y5
      ↪ ,z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10,x11,y11,z11
      ↪ ,x12,y12,z12,x13,y13,z13,x14,y14,z14,x15,y15,z15 --params=10.55
      ↪ ,0.708056872038,0.4388,0.5668,0.1053,0.0962,0.2042,0.5318,0.742
      ↪ ,0.2057,0.1168,0.3371,0.5582,0.0094,0.3964,0.4704,0.228,0.5646,
      ↪ 0.6512,0.0664,-0.0011,0.1089,0.6245,0.2309,0.2495,0.5637,0.0581
      ↪ ,0.2465,0.3949,0.6922,0.1208,0.2503,0.8721,0.2774,0.0897,0.6541
      ↪ ,0.2174,0.014,0.4566,0.5691,0.6236,0.1184,0.2192,0.0,0.7772,
      ↪ 0.2214,0.6381 & P3_{1} C_{3}^{2} #144 (a^15) & hP45 & None &
```

```
      ↪ NO3Rb & NO3Rb & M. Shamsuzzoha and B. W. Lucas, Acta
      ↪ Crystallogr. Sect. B Struct. Sci. 38, 2353-2357 (1982)
   1.00000000000000
   5.27500000000000  -9.13656800992583   0.00000000000000
   5.27500000000000   9.13656800992583   0.00000000000000
   0.00000000000000   0.00000000000000   7.47000000000000
     N     O     Rb
     9    27     9
Direct
   0.43880000000000   0.56680000000000   0.10530000000000   N   (3a)
  -0.56680000000000  -0.12800000000000   0.43863333333333   N   (3a)
   0.12800000000000  -0.43880000000000   0.77196666666667   N   (3a)
   0.09620000000000   0.20420000000000   0.53180000000000   N   (3a)
  -0.20420000000000  -0.10800000000000   0.86513333333333   N   (3a)
   0.10800000000000  -0.09620000000000   1.19846666666667   N   (3a)
   0.74200000000000   0.20570000000000   0.11680000000000   N   (3a)
  -0.20570000000000   0.53630000000000   0.45013333333333   N   (3a)
  -0.53630000000000  -0.74200000000000   0.78346666666667   N   (3a)
   0.33710000000000   0.55820000000000   0.00940000000000   O   (3a)
  -0.55820000000000  -0.22110000000000   0.34273333333333   O   (3a)
   0.22110000000000  -0.33710000000000   0.67606666666667   O   (3a)
   0.39640000000000   0.47040000000000   0.22800000000000   O   (3a)
  -0.47040000000000  -0.07400000000000   0.56133333333333   O   (3a)
   0.07400000000000  -0.39640000000000   0.89466666666667   O   (3a)
   0.56460000000000   0.65120000000000   0.06640000000000   O   (3a)
  -0.65120000000000  -0.08660000000000   0.39973333333333   O   (3a)
   0.08660000000000  -0.56460000000000   0.73306666666667   O   (3a)
  -0.00110000000000   0.10890000000000   0.62450000000000   O   (3a)
  -0.10890000000000  -0.11000000000000   0.95783333333333   O   (3a)
   0.11000000000000   0.00110000000000   1.29116666666667   O   (3a)
   0.23090000000000   0.24950000000000   0.56370000000000   O   (3a)
  -0.24950000000000  -0.01860000000000   0.89703333333333   O   (3a)
   0.01860000000000  -0.23090000000000   1.23036666666667   O   (3a)
   0.05810000000000   0.24650000000000   0.39490000000000   O   (3a)
  -0.24650000000000  -0.18840000000000   0.72823333333333   O   (3a)
   0.18840000000000  -0.05810000000000   1.06156666666667   O   (3a)
   0.69220000000000   0.12080000000000   0.25030000000000   O   (3a)
  -0.12080000000000   0.57140000000000   0.58363333333333   O   (3a)
  -0.57140000000000  -0.69220000000000   0.91696666666667   O   (3a)
   0.87210000000000   0.27740000000000   0.08970000000000   O   (3a)
  -0.27740000000000   0.59470000000000   0.42303333333333   O   (3a)
  -0.59470000000000  -0.87210000000000   0.75636666666667   O   (3a)
   0.65410000000000   0.21740000000000   0.01400000000000   O   (3a)
  -0.21740000000000   0.43670000000000   0.34733333333333   O   (3a)
  -0.43670000000000  -0.65410000000000   0.68066666666667   O   (3a)
   0.45660000000000   0.56910000000000   0.62360000000000   Rb  (3a)
  -0.56910000000000  -0.11250000000000   0.95693333333333   Rb  (3a)
   0.11250000000000  -0.45660000000000   1.29026666666667   Rb  (3a)
   0.11840000000000   0.21920000000000   0.00000000000000   Rb  (3a)
  -0.21920000000000  -0.10080000000000   0.33333333333333   Rb  (3a)
   0.10080000000000  -0.11840000000000   0.66666666666667   Rb  (3a)
   0.77720000000000   0.22140000000000   0.63810000000000   Rb  (3a)
  -0.22140000000000   0.55580000000000   0.97143333333333   Rb  (3a)
  -0.55580000000000  -0.77720000000000   1.30476666666667   Rb  (3a)
```

Na_2SO_3 ($G3_2$): A2B3C_hP12_147_abd_g_d - CIF

```
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_audit_creation_method FINDSYM

_chemical_name_mineral 'Na2O3S'
_chemical_formula_sum 'Na2 O3 S'

loop_
_publ_author_name
 'W. H. Zachariasen'
 'H. E. Buckley'
_journal_name_full_name
;
 Physical Review
;
_journal_volume 37
_journal_year 1931
_journal_page_first 1295
_journal_page_last 1305
_publ_Section_title
;
 The Crystal Lattice of Anhydrous Sodium Sulphite, Na${2}$SO${3}$
;

_aflow_title 'Na${2}$SO${3}$ ($G3_{2}$) Structure'
_aflow_proto 'A2B3C_hP12_147_abd_g_d'
_aflow_params 'a,c/a,z_{3},z_{4},x_{5},y_{5},z_{5}'
_aflow_params_values '5.441,1.12718250322,0.67,0.17,0.14,0.4,0.25'
_aflow_Strukturbericht '$G3_{2}$'
_aflow_Pearson 'hP12'

_symmetry_space_group_name_H-M "P -3"
_symmetry_Int_Tables_number 147

_cell_length_a    5.44100
_cell_length_b    5.44100
_cell_length_c    6.13300
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 -x,-y,-z
5 y,-x+y,-z
6 x-y,x,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Na1 Na 1 a 0.00000 0.00000 0.00000 1.00000
Na2 Na 1 b 0.00000 0.00000 0.50000 1.00000
Na3 Na 2 d 0.33333 0.66667 0.67000 1.00000
S1 S 2 d 0.33333 0.66667 0.17000 1.00000
O1 O 6 g 0.14000 0.40000 0.25000 1.00000
```

Na_2SO_3 ($G3_2$): A2B3C_hP12_147_abd_g_d - POSCAR

```
A2B3C_hP12_147_abd_g_d & a,c/a,z3,z4,x5,y5,z5 --params=5.441,
      ↪ 1.12718250322,0.67,0.17,0.14,0.4,0.25 & P-3 C_{3i}^{1} #147 (
      ↪ abd^2g) & hP12 & $G3_{2}$ & Na2O3S & Na2O3S & W. H. Zachariasen
      ↪ and H. E. Buckley, Phys. Rev. 37, 1295-1305 (1931)
   1.00000000000000
   2.72050000000000  -4.71204422199113   0.00000000000000
   2.72050000000000   4.71204422199113   0.00000000000000
   0.00000000000000   0.00000000000000   6.13300000000000
    Na    O    S
     4    6    2
Direct
   0.00000000000000   0.00000000000000   0.00000000000000   Na   (1a)
   0.00000000000000   0.00000000000000   0.50000000000000   Na   (1b)
   0.33333333333333   0.66666666666667   0.67000000000000   Na   (2d)
   0.66666666666667   0.33333333333333  -0.67000000000000   Na   (2d)
   0.14000000000000   0.40000000000000   0.25000000000000   O    (6g)
  -0.40000000000000  -0.26000000000000   0.25000000000000   O    (6g)
   0.26000000000000  -0.14000000000000   0.25000000000000   O    (6g)
  -0.14000000000000  -0.40000000000000  -0.25000000000000   O    (6g)
   0.40000000000000   0.26000000000000  -0.25000000000000   O    (6g)
  -0.26000000000000   0.14000000000000  -0.25000000000000   O    (6g)
   0.33333333333333   0.66666666666667   0.17000000000000   S    (2d)
   0.66666666666667   0.33333333333333  -0.17000000000000   S    (2d)
```

Dolomite [$MgCa(CO_3)_2$, $G1_1$]: A2BCD6_hR10_148_c_a_b_f - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Dolomite'
_chemical_formula_sum 'C2 Ca Mg O6'

loop_
_publ_author_name
 'R. J. Reeder'
 'S. A. Markgraf'
_journal_name_full_name
;
 American Mineralogist
;
_journal_volume 71
_journal_year 1986
_journal_page_first 795
_journal_page_last 804
_publ_Section_title
;
 High-temperature crystal chemistry of dolomite
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'Dolomite [MgCa(CO${3}$)${2}$, $G1_{1}$] Structure'
_aflow_proto 'A2BCD6_hR10_148_c_a_b_f'
_aflow_params 'a,c/a,x_{3},x_{4},y_{4},z_{4}'
_aflow_params_values '4.8069,3.32896461337,0.24289,0.49193,0.27933,-
      ↪ 0.03947'
_aflow_Strukturbericht '$G1_{1}$'
_aflow_Pearson 'hR10'

_symmetry_space_group_name_H-M "R -3 (hexagonal axes)"
_symmetry_Int_Tables_number 148

_cell_length_a    4.80690
_cell_length_b    4.80690
_cell_length_c    16.00200
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 -x,-y,-z
5 y,-x+y,-z
6 x-y,x,-z
7 x+1/3,y+2/3,z+2/3
8 -y+1/3,x-y+2/3,z+2/3
9 -x+y+1/3,-x+2/3,z+2/3
10 -x+1/3,-y+2/3,-z+2/3
11 y+1/3,-x+y+2/3,-z+2/3
12 x-y+1/3,x+2/3,-z+2/3
```

```
13 x+2/3,y+1/3,z+1/3
14 -y+2/3,x-y+1/3,z+1/3
15 -x+y+2/3,-x+1/3,z+1/3
16 -x+2/3,-y+1/3,-z+1/3
17 y+2/3,-x+y+1/3,-z+1/3
18 x-y+2/3,x+1/3,-z+1/3

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ca1 Ca 3 a 0.00000 0.00000 0.00000 1.00000
Mg1 Mg 3 b 0.00000 0.00000 0.50000 1.00000
C1 C 6 c 0.00000 0.00000 0.24289 1.00000
O1 O 18 f 0.24800 -0.0354 0.24393 1.00000
```

Dolomite [$MgCa(CO_3)_2$, $G1_1$]: A2BCD6_hR10_148_c_a_b_f - POSCAR

```
A2BCD6_hR10_148_c_a_b_f & a,c/a,x3,x4,y4,z4 --params=4.8069,
    ↪ 3.32896461337,0.24289,0.49193,0.27933,-0.03947 & R-3 C_{3i}^{2}
    ↪ #148 (abcf) & hR10 & $G1_{1}$ & C2CaMgO6 & Dolomite & R. J.
    ↪ Reeder and S. A. Markgraf, Am. Mineral. 71, 795-804 (1986)
   1.00000000000000
   2.40345000000000  -1.38763250448381   5.33400000000000
   0.00000000000000   2.77526500896761   5.33400000000000
  -2.40345000000000  -1.38763250448381   5.33400000000000
     C     Ca     Mg     O
     2      1      1     6
Direct
   0.24289000000000   0.24289000000000   0.24289000000000   C    (2c)
  -0.24289000000000  -0.24289000000000  -0.24289000000000   C    (2c)
   0.00000000000000   0.00000000000000   0.00000000000000   Ca   (1a)
   0.50000000000000   0.50000000000000   0.50000000000000   Mg   (1b)
   0.49193000000000   0.27933000000000  -0.03947000000000   O    (6f)
  -0.03947000000000   0.49193000000000   0.27933000000000   O    (6f)
   0.27933000000000  -0.03947000000000   0.49193000000000   O    (6f)
  -0.49193000000000  -0.27933000000000   0.03947000000000   O    (6f)
   0.03947000000000  -0.49193000000000  -0.27933000000000   O    (6f)
  -0.27933000000000   0.03947000000000  -0.49193000000000   O    (6f)
```

$K_2Sn(OH)_6$ ($H6_2$): A6B2C6D_hR15_148_f_c_f_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'H6K2O6Sn'
_chemical_formula_sum 'H6 K2 O6 Sn'

loop_
_publ_author_name
 'H. Jacobs'
 'R. Stahl'
_journal_name_full_name
;
 Zeitschrift fur Anorganische und Allgemeine Chemie
;
_journal_volume 626
_journal_year 2000
_journal_page_first 1863
_journal_page_last 1866
_publ_Section_title
;
 Neubestimmung der Kristallstrukturen der Hexahydroxometallate Na$_{2}
    ↪ $Sn(OH)$_{6}$, K$_{2}$Sn(OH)$_{6}$ und K$_{2}$Pb(OH)$_{6}$
;

# Found in K$_{2}$Sn(OH)$_{6}$ (K$_{2}$Sn[OH]$_{6}$) Crystal Structure,
    ↪ {PAULING FILE in: Inorganic Solid Phases, SpringerMaterials (
    ↪ online database), Springer, Heidelberg (ed.) Springer Materials
    ↪ }, 2016

_aflow_title 'K$_{2}$Sn(OH)$_{6}$ ($H6_{2}$) Structure'
_aflow_proto 'A6B2C6D_hR15_148_f_c_f_a'
_aflow_params 'a,c/a,x_{2},x_{3},y_{3},z_{3},x_{4},y_{4},z_{4}'
_aflow_params_values '6.541,1.95887478979,0.28715,0.248,-0.202,0.326,
    ↪ 0.296,-0.2014,0.1778'
_aflow_Strukturbericht '$H6_{2}$'
_aflow_Pearson 'hR15'

_symmetry_space_group_name_H-M "R -3 (hexagonal axes)"
_symmetry_Int_Tables_number 148

_cell_length_a    6.54100
_cell_length_b    6.54100
_cell_length_c    12.81300
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 -x,-y,-z
5 y,-x+y,-z
6 x-y,x,-z
7 x+1/3,y+2/3,z+2/3
8 -y+1/3,x-y+2/3,z+2/3
9 -x+y+1/3,-x+2/3,z+2/3
10 -x+1/3,-y+2/3,-z+2/3
11 y+1/3,-x+y+2/3,-z+2/3
12 x-y+1/3,x+2/3,-z+2/3
13 x+2/3,y+1/3,z+1/3
14 -y+2/3,x-y+1/3,z+1/3
15 -x+y+2/3,-x+1/3,z+1/3
16 -x+2/3,-y+1/3,-z+1/3
17 y+2/3,-x+y+1/3,-z+1/3
18 x-y+2/3,x+1/3,-z+1/3

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Sn1 Sn 3 a 0.00000 0.00000 0.00000 1.00000
K1 K 6 c 0.00000 0.00000 0.28715 1.00000
H1 H 18 f 0.12400 0.32600 0.12400 1.00000
O1 O 18 f 0.20520 0.29220 0.09080 1.00000
```

$K_2Sn(OH)_6$ ($H6_2$): A6B2C6D_hR15_148_f_c_f_a - POSCAR

```
A6B2C6D_hR15_148_f_c_f_a & a,c/a,x2,x3,y3,z3,x4,y4,z4 --params=6.541,
    ↪ 1.95887478979,0.28715,0.248,-0.202,0.326,0.296,-0.2014,0.1778 &
    ↪ R-3 C_{3i}^{2} #148 (acf^2) & hR15 & $H6_{2}$ & H6K2O6Sn &
    ↪ H6K2O6Sn & H. Jacobs and R. Stahl, Z. Anorg. Allg. Chem. 626,
    ↪ 1863-1866 (2000)
   1.00000000000000
   3.27050000000000  -1.88822405538467   4.27100000000000
   0.00000000000000   3.77644811076934   4.27100000000000
  -3.27050000000000  -1.88822405538467   4.27100000000000
     H     K     O     Sn
     6     2     6      1
Direct
   0.24800000000000  -0.20200000000000   0.32600000000000   H    (6f)
   0.32600000000000   0.24800000000000  -0.20200000000000   H    (6f)
  -0.20200000000000   0.32600000000000   0.24800000000000   H    (6f)
  -0.24800000000000   0.20200000000000  -0.32600000000000   H    (6f)
  -0.32600000000000  -0.24800000000000   0.20200000000000   H    (6f)
   0.20200000000000  -0.32600000000000  -0.24800000000000   H    (6f)
   0.28715000000000   0.28715000000000   0.28715000000000   K    (2c)
  -0.28715000000000  -0.28715000000000  -0.28715000000000   K    (2c)
   0.29600000000000  -0.20140000000000   0.17780000000000   O    (6f)
   0.17780000000000   0.29600000000000  -0.20140000000000   O    (6f)
  -0.20140000000000   0.17780000000000   0.29600000000000   O    (6f)
  -0.29600000000000   0.20140000000000  -0.17780000000000   O    (6f)
  -0.17780000000000  -0.29600000000000   0.20140000000000   O    (6f)
   0.20140000000000  -0.17780000000000  -0.29600000000000   O    (6f)
   0.00000000000000   0.00000000000000   0.00000000000000   Sn   (1a)
```

$Ni(H_2O)_6SnCl_6$ ($I6_1$): A6B6CD_hR14_148_f_f_b_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Cl6(H2O)6NiSn'
_chemical_formula_sum 'Cl6 (H2O)6 Ni Sn'

loop_
_publ_author_name
 'L. Pauling'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 72
_journal_year 1930
_journal_page_first 482
_journal_page_last 492
_publ_Section_title
;
 On the crystal structure of nickel chlorostannate hexahydrate
;

# Found in Strukturbericht Band II 1928-1932, 1937

_aflow_title 'Ni(H$_{2}$O)$_{6}$SnCl$_{6}$ ($I6_{1}$) Structure'
_aflow_proto 'A6B6CD_hR14_148_f_f_b_a'
_aflow_params 'a,c/a,x_{3},y_{3},z_{3},x_{4},y_{4},z_{4}'
_aflow_params_values '10.59967,1.01326739417,0.94,-0.69,0.14,0.44,-0.19,
    ↪ 0.64001'
_aflow_Strukturbericht '$I6_{1}$'
_aflow_Pearson 'hR14'

_symmetry_space_group_name_H-M "R -3 (hexagonal axes)"
_symmetry_Int_Tables_number 148

_cell_length_a    10.59967
_cell_length_b    10.59967
_cell_length_c    10.74030
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
```

```
2 -y,x-y,z
3 -x+y,-x,z
4 -x,-y,-z
5 y,-x+y,-z
6 x-y,x,-z
7 x+1/3,y+2/3,z+2/3
8 -y+1/3,x-y+2/3,z+2/3
9 -x+y+1/3,-x+2/3,z+2/3
10 -x+1/3,-y+2/3,-z+2/3
11 y+1/3,-x+y+2/3,-z+2/3
12 x-y+1/3,x+2/3,-z+2/3
13 x+2/3,y+1/3,z+1/3
14 -y+2/3,x-y+1/3,z+1/3
15 -x+y+2/3,-x+1/3,z+1/3
16 -x+2/3,-y+1/3,-z+1/3
17 y+2/3,-x+y+1/3,-z+1/3
18 x-y+2/3,x+1/3,-z+1/3

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Sn1 Sn 3 a 0.00000 0.00000 0.00000 1.00000
Ni1 Ni 3 b 0.00000 0.00000 0.50000 1.00000
Cl1 Cl 18 f 0.81000 0.82000 0.13000 1.00000
H2O1 H2O 18 f 0.14333 0.48667 0.29667 1.00000
```

$Ni(H_2O)_6SnCl_6$ ($I6_1$): A6B6CD_hR14_148_f_f_b_a - POSCAR

```
A6B6CD_hR14_148_f_f_b_a & a,c/a,x3,y3,z3,x4,y4,z4 --params=10.59967,
      ↪ 1.01326739417,0.94,-0.69,0.14,0.44,-0.19,0.64001 & R-3 C_{3i}^{
      ↪ 2} #148 (abf^2) & hR14 & $I6_{1}$ & Cl6(H2O)6NiSn & Cl6(H2O)
      ↪ 6NiSn & L. Pauling, Zeitschrift f{"u}r Kristallographie -
      ↪ Crystalline Materials 72, 482-492 (1930)
   1.00000000000000
   5.29983500000000  -3.05986116391060   3.58010000000000
   0.00000000000000   6.11972232782120   3.58010000000000
  -5.29983500000000  -3.05986116391060   3.58010000000000
     Cl    H2O     Ni     Sn
      6      6      1      1
Direct
   0.94000000000000  -0.69000000000000   0.14000000000000   Cl    (6f)
   0.14000000000000   0.94000000000000  -0.69000000000000   Cl    (6f)
  -0.69000000000000   0.14000000000000   0.94000000000000   Cl    (6f)
  -0.94000000000000   0.69000000000000  -0.14000000000000   Cl    (6f)
  -0.14000000000000  -0.94000000000000   0.69000000000000   Cl    (6f)
   0.69000000000000  -0.14000000000000  -0.94000000000000   Cl    (6f)
   0.44000000000000  -0.19000000000000   0.64001000000000   H2O   (6f)
   0.64001000000000   0.44000000000000  -0.19000000000000   H2O   (6f)
  -0.19000000000000   0.64001000000000   0.44000000000000   H2O   (6f)
  -0.44000000000000   0.19000000000000  -0.64001000000000   H2O   (6f)
  -0.64001000000000  -0.44000000000000   0.19000000000000   H2O   (6f)
   0.19000000000000  -0.64001000000000  -0.44000000000000   H2O   (6f)
   0.50000000000000   0.50000000000000   0.50000000000000   Ni    (1b)
   0.00000000000000   0.00000000000000   0.00000000000000   Sn    (1a)
```

Li_7TaO_6: A8B6C_hR15_148_cf_f_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Li7O6Ta'
_chemical_formula_sum 'Li8 O6 Ta'

loop_
_publ_author_name
 'G. Wehrum'
 'R. Hoppe'
_journal_name_full_name
;
 Zeitschrift fur Anorganische und Allgemeine Chemie
;
_journal_volume 620
_journal_year 1994
_journal_page_first 659
_journal_page_last 664
_publ_Section_title
;
 Zur Kenntnis 'Kationen-reicher\' Tantalate {\"U}ber Li$_{7}$[TaO$_{6}$]
;

# Found in The solid-state Li-ion conductor Li$_{7}$TaO$_{6}$: A
      ↪ combined computational and experimental study, {arXiv:
      ↪ 1910.11079 [cond-mat.mtrl-sci]}, 2019

_aflow_title 'Li$_{7}$TaO$_{6}$ Structure'
_aflow_proto 'A8B6C_hR15_148_cf_f_a'
_aflow_params 'a,c/a,x_{2},x_{3},y_{3},z_{3},x_{4},y_{4},z_{4}'
_aflow_params_values '5.358,2.81317655842,0.6615,0.5001,-0.2379,0.1053,
      ↪ 0.7714,-0.6257,0.0853'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hR15'

_symmetry_space_group_name_H-M "R -3 (hexagonal axes)"
_symmetry_Int_Tables_number 148

_cell_length_a    5.35800
_cell_length_b    5.35800
_cell_length_c    15.07300
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 -x,-y,-z
5 y,-x+y,-z
6 x-y,x,-z
7 x+1/3,y+2/3,z+2/3
8 -y+1/3,x-y+2/3,z+2/3
9 -x+y+1/3,-x+2/3,z+2/3
10 -x+1/3,-y+2/3,-z+2/3
11 y+1/3,-x+y+2/3,-z+2/3
12 x-y+1/3,x+2/3,-z+2/3
13 x+2/3,y+1/3,z+1/3
14 -y+2/3,x-y+1/3,z+1/3
15 -x+y+2/3,-x+1/3,z+1/3
16 -x+2/3,-y+1/3,-z+1/3
17 y+2/3,-x+y+1/3,-z+1/3
18 x-y+2/3,x+1/3,-z+1/3

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ta1 Ta 3 a 0.00000 0.00000 0.00000 1.00000
Li1 Li 6 c 0.00000 0.00000 0.66150 0.50000
Li2 Li 18 f 0.37760 0.36040 0.12250 1.00000
O1 O 18 f 0.69440 0.70270 0.07700 1.00000
```

Li_7TaO_6: A8B6C_hR15_148_cf_f_a - POSCAR

```
A8B6C_hR15_148_cf_f_a & a,c/a,x2,x3,y3,z3,x4,y4,z4 --params=5.358,
      ↪ 2.81317655842,0.6615,0.5001,-0.2379,0.1053,0.7714,-0.6257,
      ↪ 0.0853 & R-3 C_{3i}^{2} #148 (acf^2) & hR15 & None & Li7O6Ta &
      ↪ Li7O6Ta & G. Wehrum and R. Hoppe, Z. Anorg. Allg. Chem. 620,
      ↪ 659-664 (1994)
   1.00000000000000
   2.67900000000000  -1.54672137115901   5.02433333333333
   0.00000000000000   3.09344274231802   5.02433333333333
  -2.67900000000000  -1.54672137115901   5.02433333333333
     Li     O     Ta
      8     6      1
Direct
   0.66150000000000   0.66150000000000   0.66150000000000   Li    (2c)
  -0.66150000000000  -0.66150000000000  -0.66150000000000   Li    (2c)
   0.50010000000000  -0.23790000000000   0.10530000000000   Li    (6f)
   0.10530000000000   0.50010000000000  -0.23790000000000   Li    (6f)
  -0.23790000000000   0.10530000000000   0.50010000000000   Li    (6f)
  -0.50010000000000   0.23790000000000  -0.10530000000000   Li    (6f)
  -0.10530000000000  -0.50010000000000   0.23790000000000   Li    (6f)
   0.23790000000000  -0.10530000000000  -0.50010000000000   Li    (6f)
   0.77140000000000  -0.62570000000000   0.08530000000000   O     (6f)
   0.08530000000000   0.77140000000000  -0.62570000000000   O     (6f)
  -0.62570000000000   0.08530000000000   0.77140000000000   O     (6f)
  -0.77140000000000   0.62570000000000  -0.08530000000000   O     (6f)
  -0.08530000000000  -0.77140000000000   0.62570000000000   O     (6f)
   0.62570000000000  -0.08530000000000  -0.77140000000000   O     (6f)
   0.00000000000000   0.00000000000000   0.00000000000000   Ta    (1a)
```

$SrCl_2{\cdot}(H_2O)_6$: A2B12C6D_hP21_150_d_2g_ef_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Cl2H12O6Sr'
_chemical_formula_sum 'Cl2 H12 O6 Sr'

loop_
_publ_author_name
 'P. A. Agron'
 'W. R. Busing'
_journal_name_full_name
;
 Acta Crystallographica Section C: Structural Chemistry
;
_journal_volume 42
_journal_year 1986
_journal_page_first 141
_journal_page_last 143
_publ_Section_title
;
 Calcium and strontium dichloride hexahydrates by neutron diffraction
;

# Found in PAULING FILE, 2016 Found in PAULING FILE, {in: Inorganic
      ↪ Solid Phases, SpringerMaterials (online database), Springer,
      ↪ Heidelberg SpringerMaterials},

_aflow_title 'SrCl$_{2}$$\cdot$(H$_{2}$O)$_{6}$ Structure'
_aflow_proto 'A2B12C6D_hP21_150_d_2g_ef_a'
_aflow_params 'a,c/a,z_{2},x_{3},x_{4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{
      ↪ 6}'
```

```
_aflow_params_values '7.9596,0.518154178602,0.429,0.3114,0.7868,0.4326,
    ↪ 0.0988,-0.0926,0.767,0.1113,0.4835'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hP21'

_symmetry_space_group_name_H-M "P 3 2 1"
_symmetry_Int_Tables_number 150

_cell_length_a 7.95960
_cell_length_b 7.95960
_cell_length_c 4.12430
_cell_angle_alpha 90.00000
_cell_angle_beta 90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 x-y,-y,-z
5 y,x,-z
6 -x,-x+y,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Sr1 Sr 1 a 0.00000 0.00000 0.00000 1.00000
Cl1 Cl 2 d 0.33333 0.66667 0.42900 1.00000
O1 O 3 e 0.31140 0.00000 0.00000 1.00000
O2 O 3 f 0.78680 0.00000 0.50000 1.00000
H1 H 6 g 0.43260 0.09880 -0.09260 1.00000
H2 H 6 g 0.76700 0.11130 0.48350 1.00000
```

$SrCl_2{\cdot}(H_2O)_6$: A2B12C6D_hP21_150_d_2g_ef_a - POSCAR

```
A2B12C6D_hP21_150_d_2g_ef_a & a,c/a,z2,x3,x4,x5,y5,z5,x6,y6,z6 --params=
    ↪ 7.9596,0.518154178602,0.429,0.3114,0.7868,0.4326,0.0988,-0.0926
    ↪ ,0.767,0.1113,0.4835 & P321 D_{3}^{2} #150 (adefg^2) & hP21 &
    ↪ None & Cl2H12O6Sr & Cl2H12O6Sr & P. A. Agron and W. R. Busing,
    ↪ Acta Crystallogr. C 42, 141-143 (1986)
   1.00000000000000
   3.97980000000000  -6.89321580396262   0.00000000000000
   3.97980000000000   6.89321580396262   0.00000000000000
   0.00000000000000   0.00000000000000   4.12430000000000
     Cl    H    O   Sr
      2   12    6    1
Direct
   0.33333333333333   0.66666666666667   0.42900000000000  Cl  (2d)
   0.66666666666667   0.33333333333333  -0.42900000000000  Cl  (2d)
   0.43260000000000   0.09880000000000  -0.09260000000000  H  (6g)
  -0.09880000000000   0.33380000000000  -0.09260000000000  H  (6g)
  -0.33380000000000  -0.43260000000000  -0.09260000000000  H  (6g)
   0.09880000000000   0.43260000000000   0.09260000000000  H  (6g)
   0.33380000000000  -0.09880000000000   0.09260000000000  H  (6g)
  -0.43260000000000  -0.33380000000000   0.09260000000000  H  (6g)
   0.76700000000000   0.11130000000000   0.48350000000000  H  (6g)
  -0.11130000000000   0.65570000000000   0.48350000000000  H  (6g)
  -0.65570000000000  -0.76700000000000   0.48350000000000  H  (6g)
   0.11130000000000   0.76700000000000  -0.48350000000000  H  (6g)
   0.65570000000000  -0.11130000000000  -0.48350000000000  H  (6g)
  -0.76700000000000  -0.65570000000000  -0.48350000000000  H  (6g)
   0.31140000000000   0.00000000000000   0.00000000000000  O  (3e)
   0.00000000000000   0.31140000000000   0.00000000000000  O  (3e)
  -0.31140000000000  -0.31140000000000   0.00000000000000  O  (3e)
   0.78680000000000   0.00000000000000   0.50000000000000  O  (3f)
   0.00000000000000   0.78680000000000   0.50000000000000  O  (3f)
  -0.78680000000000  -0.78680000000000   0.50000000000000  O  (3f)
   0.00000000000000   0.00000000000000   0.00000000000000  Sr  (1a)
```

$Cs_3As_2Cl_9$ ($K7_3$): A2B9C3_hP14_150_d_eg_ad - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'As2Cl9Cs3'
_chemical_formula_sum 'As2 Cl9 Cs3'

loop_
_publ_author_name
 'J. L. Hoard'
 'L. Goldstein'
_journal_name_full_name
;
 Journal of Chemical Physics
;
_journal_volume 3
_journal_year 1935
_journal_page_first 117
_journal_page_last 122
_publ_Section_title
;
 The Structure of Caesium Enneachlordiarsenite, Cs$_{3}$As$_{2}$Cl$_{9}$
;

_aflow_title 'Cs$_{3}$As$_{2}$Cl$_{9}$ ($K7_{3}$) Structure'
_aflow_proto 'A2B9C3_hP14_150_d_eg_ad'
_aflow_params 'a,c/a,z_{2},z_{3},x_{4},x_{5},y_{5},z_{5}'
_aflow_params_values '7.37,1.20895522388,0.805,0.32,0.48,0.365,0.2,0.325
    ↪ '
_aflow_Strukturbericht '$K7_{3}$'
_aflow_Pearson 'hP14'

_symmetry_space_group_name_H-M "P 3 2 1"
_symmetry_Int_Tables_number 150

_cell_length_a 7.37000
_cell_length_b 7.37000
_cell_length_c 8.91000
_cell_angle_alpha 90.00000
_cell_angle_beta 90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 x-y,-y,-z
5 y,x,-z
6 -x,-x+y,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cs1 Cs 1 a 0.00000 0.00000 0.00000 1.00000
As1 As 2 d 0.33333 0.66667 0.80500 1.00000
Cs2 Cs 2 d 0.33333 0.66667 0.32000 1.00000
Cl1 Cl 3 e 0.48000 0.00000 0.00000 1.00000
Cl2 Cl 6 g 0.36500 0.20000 0.32500 1.00000
```

$Cs_3As_2Cl_9$ ($K7_3$): A2B9C3_hP14_150_d_eg_ad - POSCAR

```
A2B9C3_hP14_150_d_eg_ad & a,c/a,z2,z3,x4,x5,y5,z5 --params=7.37,
    ↪ 1.20895522388,0.805,0.32,0.48,0.365,0.2,0.325 & P321 D_{3}^{2}
    ↪ #150 (ad^2eg) & hP14 & $K7_{3}$ & As2Cl9Cs3 & As2Cl9Cs3 & J. L.
    ↪ Hoard and L. Goldstein, J. Chem. Phys. 3, 117-122 (1935)
   1.00000000000000
   3.68500000000000  -6.38260722589131   0.00000000000000
   3.68500000000000   6.38260722589131   0.00000000000000
   0.00000000000000   0.00000000000000   8.91000000000000
     As    Cl    Cs
      2     9     3
Direct
   0.33333333333333   0.66666666666667   0.80500000000000  As  (2d)
   0.66666666666667   0.33333333333333  -0.80500000000000  As  (2d)
   0.48000000000000   0.00000000000000   0.00000000000000  Cl  (3e)
   0.00000000000000   0.48000000000000   0.00000000000000  Cl  (3e)
  -0.48000000000000  -0.48000000000000   0.00000000000000  Cl  (3e)
   0.36500000000000   0.20000000000000   0.32500000000000  Cl  (6g)
  -0.20000000000000   0.16500000000000   0.32500000000000  Cl  (6g)
  -0.16500000000000  -0.36500000000000   0.32500000000000  Cl  (6g)
   0.20000000000000   0.36500000000000  -0.32500000000000  Cl  (6g)
   0.16500000000000  -0.20000000000000  -0.32500000000000  Cl  (6g)
  -0.36500000000000  -0.16500000000000  -0.32500000000000  Cl  (6g)
   0.00000000000000   0.00000000000000   0.00000000000000  Cs  (1a)
   0.33333333333333   0.66666666666667   0.32000000000000  Cs  (2d)
   0.66666666666667   0.33333333333333  -0.32000000000000  Cs  (2d)
```

Paralstonite ($BaCa(CO_3)_2$): AB2CD6_hP30_150_e_c2d_f_3g - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Paralstonite'
_chemical_formula_sum 'Ba C2 Ca O6'

loop_
_publ_author_name
 'H. Effenberger'
_journal_name_full_name
;
 Neues Jahrbuch fur Mineralogie, Monatshefte
;
_journal_volume 1980
_journal_year 1980
_journal_page_first 353
_journal_page_last 363
_publ_Section_title
;
 Die Kristallstruktur des Minerals Paralstonite, BaCa(CO$_{3}$)$_{2}$
;

# Found in A new BaCa(CO$_{3}$)$_{2}$ polymorph, 2019

_aflow_title 'Paralstonite (BaCa(CO$_{3}$)$_{2}$) Structure'
_aflow_proto 'AB2CD6_hP30_150_e_c2d_f_3g'
_aflow_params 'a,c/a,z_{1},z_{2},z_{3},x_{4},x_{5},x_{6},y_{6},z_{6},x_{
    ↪ 7},y_{7},z_{7},x_{8},y_{8},z_{8}'
_aflow_params_values '8.692,0.707317073171,0.267,0.356,0.817,0.687,
    ↪ 0.3586,0.191,0.677,0.355,0.173,0.517,0.822,0.15,0.001,0.255'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hP30'
```

```
_symmetry_space_group_name_H-M "P 3 2 1"
_symmetry_Int_Tables_number 150

_cell_length_a     8.69200
_cell_length_b     8.69200
_cell_length_c     6.14800
_cell_angle_alpha 90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 x-y,-y,-z
5 y,x,-z
6 -x,-x+y,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
C1 C 2 c 0.00000 0.00000 0.26700 1.00000
C2 C 2 d 0.33333 0.66667 0.35600 1.00000
C3 C 2 d 0.33333 0.66667 0.81700 1.00000
Ba1 Ba 3 e 0.68700 0.00000 0.00000 1.00000
Ca1 Ca 3 f 0.35860 0.00000 0.50000 1.00000
O1 O 6 g 0.19100 0.67700 0.35500 1.00000
O2 O 6 g 0.17300 0.51700 0.82200 1.00000
O3 O 6 g 0.15000 0.00100 0.25500 1.00000
```

Paralstonite ($BaCa(CO_3)_2$): AB2CD6_hP30_150_e_c2d_f_3g - POSCAR

```
AB2CD6_hP30_150_e_c2d_f_3g & a,c/a,z1,z2,z3,x4,x5,x6,y6,z6,x7,y7,z7,x8,
      ↪ y8,z8 --params=8.692,0.707317073171,0.267,0.356,0.817,0.687,
      ↪ 0.3586,0.191,0.677,0.355,0.173,0.517,0.822,0.15,0.001,0.255 &
      ↪ P321 D_{3}^{2} #150 (cd^2efg^3) & hP30 & None & BaC2CaO6 &
      ↪ Paralstonite & H. Effenberger, Neues Jahrb. Mineral. Monatsh.
      ↪ 1980, 353-363 (1980)
   1.00000000000000
   4.34600000000000  -7.52749280969434   0.00000000000000
   4.34600000000000   7.52749280969434   0.00000000000000
   0.00000000000000   0.00000000000000   6.14800000000000
    Ba    C    Ca    O
     3    6     3    18
Direct
   0.68700000000000   0.00000000000000   0.00000000000000  Ba   (3e)
   0.00000000000000   0.68700000000000   0.00000000000000  Ba   (3e)
  -0.68700000000000  -0.68700000000000   0.00000000000000  Ba   (3e)
   0.00000000000000   0.00000000000000   0.26700000000000  C    (2c)
   0.00000000000000   0.00000000000000  -0.26700000000000  C    (2c)
   0.33333333333333   0.66666666666667   0.35600000000000  C    (2d)
   0.66666666666667   0.33333333333333  -0.35600000000000  C    (2d)
   0.33333333333333   0.66666666666667   0.81700000000000  C    (2d)
   0.66666666666667   0.33333333333333  -0.81700000000000  C    (2d)
   0.35860000000000   0.00000000000000   0.50000000000000  Ca   (3f)
   0.00000000000000   0.35860000000000   0.50000000000000  Ca   (3f)
  -0.35860000000000  -0.35860000000000   0.50000000000000  Ca   (3f)
   0.19100000000000   0.67700000000000   0.35500000000000  O    (6g)
  -0.67700000000000  -0.48600000000000   0.35500000000000  O    (6g)
   0.48600000000000  -0.19100000000000   0.35500000000000  O    (6g)
   0.67700000000000   0.19100000000000  -0.35500000000000  O    (6g)
  -0.48600000000000  -0.67700000000000  -0.35500000000000  O    (6g)
  -0.19100000000000   0.48600000000000  -0.35500000000000  O    (6g)
   0.17300000000000   0.51700000000000   0.82200000000000  O    (6g)
  -0.51700000000000  -0.34400000000000   0.82200000000000  O    (6g)
   0.34400000000000  -0.17300000000000   0.82200000000000  O    (6g)
   0.51700000000000   0.17300000000000  -0.82200000000000  O    (6g)
  -0.34400000000000  -0.51700000000000  -0.82200000000000  O    (6g)
  -0.17300000000000   0.34400000000000  -0.82200000000000  O    (6g)
   0.15000000000000   0.00100000000000   0.25500000000000  O    (6g)
  -0.00100000000000   0.14900000000000   0.25500000000000  O    (6g)
  -0.14900000000000  -0.15000000000000   0.25500000000000  O    (6g)
   0.00100000000000   0.15000000000000  -0.25500000000000  O    (6g)
   0.14900000000000  -0.00100000000000  -0.25500000000000  O    (6g)
  -0.15000000000000  -0.14900000000000  -0.25500000000000  O    (6g)
```

KSO_3 ($K1_1$): AB3C_hP30_150_ef_3g_c2d - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'KO3S'
_chemical_formula_sum 'K O3 S'

loop_
_publ_author_name
 'M. L. Huggins'
 'G. O. Frank'
_journal_name_full_name
;
 American Mineralogist
;
_journal_volume 16
_journal_year 1931
_journal_page_first 580
_journal_page_last 591
_publ_Section_title
;
 The crystal structure of potassium dithionate, K$_{2}$S$_{2}$O$_{6}$
;

# Found in Strukturbericht Band II 1928-1932, 1937

_aflow_title 'KSO$_{3}$ ($K1_{1}$) Structure'
_aflow_proto 'AB3C_hP30_150_ef_3g_c2d'
_aflow_params 'a,c/a,z_{1},z_{2},z_{3},x_{4},x_{5},x_{6},y_{6},z_{6},x_{
      ↪ 7},y_{7},z_{7},x_{8},y_{8},z_{8}'
_aflow_params_values '9.82,0.647657841141,0.16,0.59,0.27,0.3,0.62,0.16,
      ↪ 0.11,0.23,0.61,0.17,0.34,0.5,0.21,0.8'
_aflow_Strukturbericht '$K1_{1}$'
_aflow_Pearson 'hP30'

_symmetry_space_group_name_H-M "P 3 2 1"
_symmetry_Int_Tables_number 150

_cell_length_a     9.82000
_cell_length_b     9.82000
_cell_length_c     6.36000
_cell_angle_alpha 90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 x-y,-y,-z
5 y,x,-z
6 -x,-x+y,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
S1 S 2 c 0.00000 0.00000 0.16000 1.00000
S2 S 2 d 0.33333 0.66667 0.59000 1.00000
S3 S 2 d 0.33333 0.66667 0.27000 1.00000
K1 K 3 e 0.30000 0.00000 0.00000 1.00000
K2 K 3 f 0.62000 0.00000 0.50000 1.00000
O1 O 6 g 0.16000 0.11000 0.23000 1.00000
O2 O 6 g 0.61000 0.17000 0.34000 1.00000
O3 O 6 g 0.50000 0.21000 0.80000 1.00000
```

KSO_3 ($K1_1$): AB3C_hP30_150_ef_3g_c2d - POSCAR

```
AB3C_hP30_150_ef_3g_c2d & a,c/a,z1,z2,z3,x4,x5,x6,y6,z6,x7,y7,z7,x8,y8,
      ↪ z8 --params=9.82,0.647657841141,0.16,0.59,0.27,0.3,0.62,0.16,
      ↪ 0.11,0.23,0.61,0.17,0.34,0.5,0.21,0.8 & P321 D_{3}^{2} #150 (cd
      ↪ ^2efg^3) & hP30 & $K1_{1}$ & KO3S & KO3S & M. L. Huggins and G.
      ↪ O. Frank, Am. Mineral. 16, 580-591 (1931)
   1.00000000000000
   4.91000000000000  -8.50436946516319   0.00000000000000
   4.91000000000000   8.50436946516319   0.00000000000000
   0.00000000000000   0.00000000000000   6.36000000000000
     K     O     S
     6    18     6
Direct
   0.30000000000000   0.00000000000000   0.00000000000000  K    (3e)
   0.00000000000000   0.30000000000000   0.00000000000000  K    (3e)
  -0.30000000000000  -0.30000000000000   0.00000000000000  K    (3e)
   0.62000000000000   0.00000000000000   0.50000000000000  K    (3f)
   0.00000000000000   0.62000000000000   0.50000000000000  K    (3f)
  -0.62000000000000  -0.62000000000000   0.50000000000000  K    (3f)
   0.16000000000000   0.11000000000000   0.23000000000000  O    (6g)
  -0.11000000000000   0.05000000000000   0.23000000000000  O    (6g)
  -0.05000000000000  -0.16000000000000   0.23000000000000  O    (6g)
   0.11000000000000   0.16000000000000  -0.23000000000000  O    (6g)
   0.05000000000000  -0.11000000000000  -0.23000000000000  O    (6g)
  -0.16000000000000  -0.05000000000000  -0.23000000000000  O    (6g)
   0.61000000000000   0.17000000000000   0.34000000000000  O    (6g)
  -0.17000000000000   0.44000000000000   0.34000000000000  O    (6g)
  -0.44000000000000  -0.61000000000000   0.34000000000000  O    (6g)
   0.17000000000000   0.61000000000000  -0.34000000000000  O    (6g)
   0.44000000000000  -0.17000000000000  -0.34000000000000  O    (6g)
  -0.61000000000000  -0.44000000000000  -0.34000000000000  O    (6g)
   0.50000000000000   0.21000000000000   0.80000000000000  O    (6g)
  -0.21000000000000   0.29000000000000   0.80000000000000  O    (6g)
  -0.29000000000000  -0.50000000000000   0.80000000000000  O    (6g)
   0.21000000000000   0.50000000000000  -0.80000000000000  O    (6g)
   0.29000000000000  -0.21000000000000  -0.80000000000000  O    (6g)
  -0.50000000000000  -0.29000000000000  -0.80000000000000  O    (6g)
   0.00000000000000   0.00000000000000   0.16000000000000  S    (2c)
   0.00000000000000   0.00000000000000  -0.16000000000000  S    (2c)
   0.33333333333333   0.66666666666667   0.59000000000000  S    (2d)
   0.66666666666667   0.33333333333333  -0.59000000000000  S    (2d)
   0.33333333333333   0.66666666666667   0.27000000000000  S    (2d)
   0.66666666666667   0.33333333333333  -0.27000000000000  S    (2d)
```

Steklite [$KAl(SO_4)_2$, $H3_2$]: ABC8D2_hP12_150_b_a_dg_d - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM
```

```
_chemical_name_mineral 'Steklite'
_chemical_formula_sum 'Al K O8 S2'

loop_
_publ_author_name
 'L. Vegard'
 'A. Maurstad'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 69
_journal_year 1929
_journal_page_first 519
_journal_page_last 532
_publ_Section_title
;
 Die Kristallstruktur der wasserfreien Alaune $R$\'$R$\'\'(SO$_{4}$)$_{2
    ↪ }$
;

# Found in Strukturbericht 1913-1928, 1931

_aflow_title 'Steklite [KAl(SO$_{4}$)$_{2}$, $H3_{2}$] Structure'
_aflow_proto 'ABC8D2_hP12_150_b_a_dg_d'
_aflow_params 'a,c/a,z_{3},z_{4},x_{5},y_{5},z_{5}'
_aflow_params_values '4.7,1.69361702128,0.016,0.222,0.328,0.344,0.317'
_aflow_Strukturbericht '$H3_{2}$'
_aflow_Pearson 'hP12'

_symmetry_space_group_name_H-M "P 3 2 1"
_symmetry_Int_Tables_number 150

_cell_length_a     4.70000
_cell_length_b     4.70000
_cell_length_c     7.96000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 x-y,-y,-z
5 y,x,-z
6 -x,-x+y,-z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
K1 K 1 a 0.00000 0.00000 0.00000 1.00000
Al1 Al 1 b 0.00000 0.00000 0.50000 1.00000
O1 O 2 d 0.33333 0.66667 0.01600 1.00000
S1 S 2 d 0.33333 0.66667 0.22200 1.00000
O2 O 6 g 0.32800 0.34400 0.31700 1.00000
```

Steklite [$KAl(SO_4)_2$, $H3_2$]: ABC8D2_hP12_150_b_a_dg_d - POSCAR

```
ABC8D2_hP12_150_b_a_dg_d & a,c/a,z3,z4,x5,y5,z5 --params=4.7,
    ↪ 1.69361702128,0.016,0.222,0.328,0.344,0.317 & P321 D_{3}^{2} #
    ↪ 150 (abd^2g) & hP12 & $H3_{2}$ & AlKO8S2 & Steklite & L. Vegard
    ↪  and A. Maurstad, Zeitschrift f{"u}r Kristallographie -
    ↪ Crystalline Materials 69, 519-532 (1929)
   1.00000000000000
   2.35000000000000  -4.07031939778686   0.00000000000000
   2.35000000000000   4.07031939778686   0.00000000000000
   0.00000000000000   0.00000000000000   7.96000000000000
    Al   K   O   S
     1   1   8   2
Direct
   0.00000000000000   0.00000000000000   0.50000000000000   Al   (1b)
   0.00000000000000   0.00000000000000   0.00000000000000   K    (1a)
   0.33333333333333   0.66666666666667   0.01600000000000   O    (2d)
   0.66666666666667   0.33333333333333  -0.01600000000000   O    (2d)
   0.32800000000000   0.34400000000000   0.31700000000000   O    (6g)
  -0.34400000000000  -0.01600000000000   0.31700000000000   O    (6g)
   0.01600000000000  -0.32800000000000   0.31700000000000   O    (6g)
   0.34400000000000   0.32800000000000  -0.31700000000000   O    (6g)
  -0.01600000000000  -0.34400000000000  -0.31700000000000   O    (6g)
  -0.32800000000000   0.01600000000000  -0.31700000000000   O    (6g)
   0.33333333333333   0.66666666666667   0.22200000000000   S    (2d)
   0.66666666666667   0.33333333333333  -0.22200000000000   S    (2d)
```

$KBe_2BO_3F_2$: AB2C2DE3_hR9_155_b_c_c_a_e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'BBe2F2KO3'
_chemical_formula_sum 'B Be2 F2 K O3'

loop_
_publ_author_name
 'L. Mei'
 'X. Huang'
 'Y. Wang'
 'Q. Wu'
 'C. Chen'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 210
_journal_year 1995
_journal_page_first 93
_journal_page_last 95
_publ_Section_title
;
 Crystal Structure of KBe$_{2}$BO$_{3}$F$_{2}$
;

_aflow_title 'KBe$_{2}$BO$_{3}$F$_{2}$ Structure'
_aflow_proto 'AB2C2DE3_hR9_155_b_c_c_a_e'
_aflow_params 'a,c/a,x_{3},x_{4},y_{5}'
_aflow_params_values '4.427,4.2340185227,0.19743,0.27848,0.19063'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hR9'

_symmetry_space_group_name_H-M "R 3 2 (hexagonal axes)"
_symmetry_Int_Tables_number 155

_cell_length_a     4.42700
_cell_length_b     4.42700
_cell_length_c     18.74400
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 y,x,-z
5 -x,-x+y,-z
6 x-y,-y,-z
7 x+1/3,y+2/3,z+2/3
8 -y+1/3,x-y+2/3,z+2/3
9 -x+y+1/3,-x+2/3,z+2/3
10 y+1/3,x+2/3,-z+2/3
11 -x+1/3,-x+y+2/3,-z+2/3
12 x-y+1/3,-y+2/3,-z+2/3
13 x+2/3,y+1/3,z+1/3
14 -y+2/3,x-y+1/3,z+1/3
15 -x+y+2/3,-x+1/3,z+1/3
16 y+2/3,x+1/3,-z+1/3
17 -x+2/3,-x+y+1/3,-z+1/3
18 x-y+2/3,-y+1/3,-z+1/3

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
K1 K 3 a 0.00000 0.00000 0.00000 1.00000
B1 B 3 b 0.00000 0.00000 0.50000 1.00000
Be1 Be 6 c 0.00000 0.00000 0.19743 1.00000
F1 F 6 c 0.00000 0.00000 0.27848 1.00000
O1 O 9 e 0.69063 0.00000 0.50000 1.00000
```

$KBe_2BO_3F_2$: AB2C2DE3_hR9_155_b_c_c_a_e - POSCAR

```
AB2C2DE3_hR9_155_b_c_c_a_e & a,c/a,x3,x4,y5 --params=4.427,4.2340185227,
    ↪ 0.19743,0.27848,0.19063 & R32 D_{3}^{7} #155 (abc^2e) & hR9 &
    ↪ None & BBe2F2KO3 & BBe2F2KO3 & L. Mei et al., Zeitschrift f{"u}
    ↪ r Kristallographie - Crystalline Materials 210, 93-95 (1995)
   1.00000000000000
   2.21350000000000  -1.27796482085124   6.24800000000000
   0.00000000000000   2.55592964170247   6.24800000000000
  -2.21350000000000  -1.27796482085124   6.24800000000000
    B   Be   F   K   O
    1    2   2   1   3
Direct
   0.50000000000000   0.50000000000000   0.50000000000000   B    (1b)
   0.19743000000000   0.19743000000000   0.19743000000000   Be   (2c)
  -0.19743000000000  -0.19743000000000  -0.19743000000000   Be   (2c)
   0.27848000000000   0.27848000000000   0.27848000000000   F    (2c)
  -0.27848000000000  -0.27848000000000  -0.27848000000000   F    (2c)
   0.00000000000000   0.00000000000000   0.00000000000000   K    (1a)
   0.50000000000000   0.19063000000000   0.19063000000000   O    (3e)
  -0.19063000000000   0.50000000000000   0.19063000000000   O    (3e)
   0.19063000000000  -0.19063000000000   0.50000000000000   O    (3e)
```

SbI_3S_{24}: A3B24C_hR28_160_b_2b3c_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'I3S24Sb'
_chemical_formula_sum 'I3 S24 Sb'

loop_
_publ_author_name
 'T. Bjorvatten'
```

```
 'O. Hassel'
 'A. Lindheim'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 17
_journal_year 1963
_journal_page_first 689
_journal_page_last 702
_publ_Section_title
;
 Crystal Structure of the Addition Compound SbI$_{3}$:3S$_{8}$
;

_aflow_title 'SbI$_{3}$S$_{24}$ Structure'
_aflow_proto 'A3B24C_hR28_160_b_2b3c_a'
_aflow_params 'a,c/a,x_{1},x_{2},z_{2},x_{3},z_{3},x_{4},z_{4},x_{5},y_{
      ↪ 5},z_{5},x_{6},y_{6},z_{6},x_{7},y_{7},z_{7}'
_aflow_params_values '24.817,0.178422049402,0.3146,-0.0551,0.1102,0.5891
      ↪ ,-0.00921,0.4515,0.17749,0.8641,-0.2688,0.24779,0.6198,-0.5701,
      ↪ 0.08749,0.7609,-0.3733,0.37359'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hR28'

_symmetry_space_group_name_H-M "R 3 m (hexagonal axes)"
_symmetry_Int_Tables_number 160

_cell_length_a    24.81700
_cell_length_b    24.81700
_cell_length_c    4.42790
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 -y,-x,z
5 x,x-y,z
6 -x+y,y,z
7 x+1/3,y+2/3,z+2/3
8 -y+1/3,x-y+2/3,z+2/3
9 -x+y+1/3,-x+2/3,z+2/3
10 -y+1/3,-x+2/3,z+2/3
11 x+1/3,x-y+2/3,z+2/3
12 -x+y+1/3,y+2/3,z+2/3
13 x+2/3,y+1/3,z+1/3
14 -y+2/3,x-y+1/3,z+1/3
15 -x+y+2/3,-x+1/3,z+1/3
16 -y+2/3,-x+1/3,z+1/3
17 x+2/3,x-y+1/3,z+1/3
18 -x+y+2/3,y+1/3,z+1/3

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Sb1 Sb 3 a 0.00000 0.00000 0.31460 1.00000
I1 I 9 b -0.05510 0.05510 0.00000 1.00000
S1 S 9 b 0.53277 0.46723 0.05633 1.00000
S2 S 9 b 0.42467 0.57533 0.02683 1.00000
S3 S 18 c 0.58307 0.54983 0.28103 1.00000
S4 S 18 c 0.57407 0.61583 0.04573 1.00000
S5 S 18 c 0.50717 0.62703 0.25373 1.00000
```

SbI_3S_{24}: A3B24C_hR28_160_b_2b3c_a - POSCAR

```
A3B24C_hR28_160_b_2b3c_a & a,c/a,x1,x2,z2,x3,z3,x4,z4,x5,y5,z5,x6,y6,z6,
      ↪ x7,y7,z7 --params=24.817,0.178422049402,0.3146,-0.0551,0.1102,
      ↪ 0.5891,-0.00921,0.4515,0.17749,0.8641,-0.2688,0.24779,0.6198,-
      ↪ 0.5701,0.08749,0.7609,-0.3733,0.37359 & R3m C_{3v}^{5} #160 (ab
      ↪ ^3c^3) & hR28 & None & I3S24Sb & I3S24Sb & T. Bjorvatten and O.
      ↪  Hassel and A. Lindheim, Acta Chem. Scand. 17, 689-702 (1963)
   1.00000000000000
  12.40850000000000   -7.16405081523947    1.47596666666667
   0.00000000000000   14.32810163047890    1.47596666666667
 -12.40850000000000   -7.16405081523947    1.47596666666667
     I     S     Sb
     3    24     1
Direct
  -0.05510000000000  -0.05510000000000   0.11020000000000     I   (3b)
   0.11020000000000  -0.05510000000000  -0.05510000000000     I   (3b)
  -0.05510000000000   0.11020000000000  -0.05510000000000     I   (3b)
   0.58910000000000   0.58910000000000  -0.00921000000000     S   (3b)
  -0.00921000000000   0.58910000000000   0.58910000000000     S   (3b)
   0.58910000000000  -0.00921000000000   0.58910000000000     S   (3b)
   0.45150000000000   0.45150000000000   0.17749000000000     S   (3b)
   0.17749000000000   0.45150000000000   0.45150000000000     S   (3b)
   0.45150000000000   0.17749000000000   0.45150000000000     S   (3b)
   0.86410000000000  -0.26880000000000   0.24779000000000     S   (6c)
   0.24779000000000   0.86410000000000  -0.26880000000000     S   (6c)
  -0.26880000000000   0.24779000000000   0.86410000000000     S   (6c)
   0.24779000000000  -0.26880000000000   0.86410000000000     S   (6c)
  -0.26880000000000   0.86410000000000   0.24779000000000     S   (6c)
   0.86410000000000   0.24779000000000  -0.26880000000000     S   (6c)
   0.61980000000000  -0.57010000000000   0.08749000000000     S   (6c)
   0.08749000000000   0.61980000000000  -0.57010000000000     S   (6c)
  -0.57010000000000   0.08749000000000   0.61980000000000     S   (6c)
   0.08749000000000  -0.57010000000000   0.61980000000000     S   (6c)
  -0.57010000000000   0.61980000000000   0.08749000000000     S   (6c)
   0.61980000000000   0.08749000000000  -0.57010000000000     S   (6c)
   0.76090000000000  -0.37330000000000   0.37359000000000     S   (6c)
   0.37359000000000   0.76090000000000  -0.37330000000000     S   (6c)
  -0.37330000000000   0.37359000000000   0.76090000000000     S   (6c)
   0.37359000000000  -0.37330000000000   0.76090000000000     S   (6c)
  -0.37330000000000   0.76090000000000   0.37359000000000     S   (6c)
   0.76090000000000   0.37359000000000  -0.37330000000000     S   (6c)
   0.31460000000000   0.31460000000000   0.31460000000000    Sb   (1a)
```

Fe_3PO_7: A3B7C_hR11_160_b_a2b_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Fe3O7P'
_chemical_formula_sum 'Fe3 O7 P'

loop_
_publ_author_name
 'A. Modaressi'
 'A. Courtois'
 'R. Gerardin'
 'B. Malaman'
 'C. Gleitzer'
_journal_name_full_name
;
 Journal of Solid State Chemistry
;
_journal_volume 47
_journal_year 1983
_journal_page_first 245
_journal_page_last 255
_publ_Section_title
;
 Fe$_{3}$PO$_{7}$, Un cas de coordinence 5 du fer trivalent, etude
      ↪ structurale et magnetique
;

# Found in C. L. Sarkis et al., Partial antiferromagnetic helical order
      ↪ in single-crystal Fe$_{3}$PO$_{4}$O$_{3}$, Phys. Rev. B 101,
      ↪ 184417 (2020)

_aflow_title 'Fe$_{3}$PO$_{7}$ Structure'
_aflow_proto 'A3B7C_hR11_160_b_a2b_a'
_aflow_params 'a,c/a,x_{1},x_{2},x_{3},z_{3},x_{4},z_{4},x_{5},z_{5}'
_aflow_params_values '8.006,0.857232075943,0.2289,0.0,0.5395,0.14849,
      ↪ 0.3895,0.76679,0.833,0.144'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hR11'

_symmetry_space_group_name_H-M "R 3 m (hexagonal axes)"
_symmetry_Int_Tables_number 160

_cell_length_a    8.00600
_cell_length_b    8.00600
_cell_length_c    6.86300
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 -y,-x,z
5 x,x-y,z
6 -x+y,y,z
7 x+1/3,y+2/3,z+2/3
8 -y+1/3,x-y+2/3,z+2/3
9 -x+y+1/3,-x+2/3,z+2/3
10 -y+1/3,-x+2/3,z+2/3
11 x+1/3,x-y+2/3,z+2/3
12 -x+y+1/3,y+2/3,z+2/3
13 x+2/3,y+1/3,z+1/3
14 -y+2/3,x-y+1/3,z+1/3
15 -x+y+2/3,-x+1/3,z+1/3
16 -y+2/3,-x+1/3,z+1/3
17 x+2/3,x-y+1/3,z+1/3
18 -x+y+2/3,y+1/3,z+1/3

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 3 a 0.00000 0.00000 0.22890 1.00000
P1 P 3 a 0.00000 0.00000 0.00000 1.00000
Fe1 Fe 9 b 0.46367 0.53633 0.07583 1.00000
O2 O 9 b 0.20757 0.79243 0.18193 1.00000
O3 O 9 b 0.56300 0.43700 0.27000 1.00000
```

Fe_3PO_7: A3B7C_hR11_160_b_a2b_a - POSCAR

```
A3B7C_hR11_160_b_a2b_a & a,c/a,x1,x2,x3,z3,x4,z4,x5,z5 --params=8.006,
    ↪ 0.857232075943,0.2289,0.0,0.5395,0.14849,0.3895,0.76679,0.833,
    ↪ 0.144 & R3m C_{3v}^{5} #160 (a^2b^3) & hR11 & None & Fe3O7P &
    ↪ Fe3O7P & A. Modaressi et al., J. Solid State Chem. 47, 245-255
    ↪ (1983)
  1.00000000000000
  4.00300000000000  -2.31113312756607   2.28766666666667
  0.00000000000000   4.62226625513214   2.28766666666667
 -4.00300000000000  -2.31113312756607   2.28766666666667
   Fe   O   P
    3   7   1
Direct
  0.53950000000000  0.53950000000000  0.14849000000000  Fe  (3b)
  0.14849000000000  0.53950000000000  0.53950000000000  Fe  (3b)
  0.53950000000000  0.14849000000000  0.53950000000000  Fe  (3b)
  0.22890000000000  0.22890000000000  0.22890000000000  O   (1a)
  0.38950000000000  0.38950000000000  0.76679000000000  O   (3b)
  0.76679000000000  0.38950000000000  0.38950000000000  O   (3b)
  0.38950000000000  0.76679000000000  0.38950000000000  O   (3b)
  0.83300000000000  0.83300000000000  0.14400000000000  O   (3b)
  0.14400000000000  0.83300000000000  0.83300000000000  O   (3b)
  0.83300000000000  0.14400000000000  0.83300000000000  O   (3b)
  0.00000000000000  0.00000000000000  0.00000000000000  P   (1a)
```

Cronstedtite {Fe(Fe,Si)[(OH)$_2$,O]O$_3$, $S5_7$}: AB3C2D_hR7_160_a_b_2a_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Cronstedtite'
_chemical_formula_sum 'Fe O3 (OH)2 Si'

loop_
_publ_author_name
 'S. B. Hendricks'
_journal_name_full_name
;
 American Mineralogist
;
_journal_volume 24
_journal_year 1939
_journal_page_first 529
_journal_page_last 539
_publ_Section_title
;
 Random structures of layer minerals as illustrated by cronstedite (
    ↪ 2FeO$\cdot$Fe$_{2}$O$_{3}$$\cdot$SiO$_{2}$$\cdot$2H$_{2}$O).
    ↪ Possible iron content of kaolin
;

# Found in Strukturbericht Band VII 1939, 1943

_aflow_title 'Cronstedtite \{Fe(Fe,Si)[(OH)$_{2}$O]O$_{3}$OH, $S5_{7}$\}
    ↪ Structure'
_aflow_proto 'AB3C2D_hR7_160_a_b_2a_a'
_aflow_params 'a,c/a,x_{1},x_{2},x_{3},x_{4},x_{5},z_{5}'
_aflow_params_values '5.48,3.87773722628,0.03,0.11,0.54,0.83,0.5,0.0'
_aflow_Strukturbericht '$S5_{7}$'
_aflow_Pearson 'hR7'

_symmetry_space_group_name_H-M "R 3 m (hexagonal axes)"
_symmetry_Int_Tables_number 160

_cell_length_a    5.48000
_cell_length_b    5.48000
_cell_length_c    21.25000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 -y,-x,z
5 x,x-y,z
6 -x+y,y,z
7 x+1/3,y+2/3,z+2/3
8 -y+1/3,x-y+2/3,z+2/3
9 -x+y+1/3,-x+2/3,z+2/3
10 -y+1/3,-x+2/3,z+2/3
11 x+1/3,x-y+2/3,z+2/3
12 -x+y+1/3,y+2/3,z+2/3
13 x+2/3,y+1/3,z+1/3
14 -y+2/3,x-y+1/3,z+1/3
15 -x+y+2/3,-x+1/3,z+1/3
16 -y+2/3,-x+1/3,z+1/3
17 x+2/3,x-y+1/3,z+1/3
18 -x+y+2/3,y+1/3,z+1/3

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Fe1 Fe 3 a 0.00000 0.00000 0.03000 0.90000
OH1 OH 3 a 0.00000 0.00000 0.11000 1.00000
OH2 OH 3 a 0.00000 0.00000 0.54000 1.00000
Si1 Si 3 a 0.00000 0.00000 0.83000 1.00000
O1 O 9 b 0.50000 0.50000 0.00000 1.00000
```

Cronstedtite {Fe(Fe,Si)[(OH)$_2$,O]O$_3$, $S5_7$}: AB3C2D_hR7_160_a_b_2a_a - POSCAR

```
AB3C2D_hR7_160_a_b_2a_a & a,c/a,x1,x2,x3,x4,x5,z5 --params=5.48,
    ↪ 3.87773722628,0.03,0.11,0.54,0.83,0.5,0.0 & R3m C_{3v}^{5} #160
    ↪  (a^4b) & hR7 & $S5_{7}$ & FeH2O5Si & Cronstedtite & S. B.
    ↪ Hendricks, Am. Mineral. 24, 529-539 (1939)
  1.00000000000000
  2.74000000000000  -1.58193973757957   7.08333333333333
  0.00000000000000   3.16387947515915   7.08333333333333
 -2.74000000000000  -1.58193973757957   7.08333333333333
   Fe   O   OH   Si
    1   3    2    1
Direct
  0.03000000000000  0.03000000000000  0.03000000000000  Fe  (1a)
  0.50000000000000  0.50000000000000  0.00000000000000  O   (3b)
  0.00000000000000  0.50000000000000  0.50000000000000  O   (3b)
  0.50000000000000  0.00000000000000  0.50000000000000  O   (3b)
  0.11000000000000  0.11000000000000  0.11000000000000  OH  (1a)
  0.54000000000000  0.54000000000000  0.54000000000000  OH  (1a)
  0.83000000000000  0.83000000000000  0.83000000000000  Si  (1a)
```

Low-Temperature GaMo$_4$S$_8$: AB4C8_hR13_160_a_ab_2a2b - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'GaMo4S8'
_chemical_formula_sum 'Ga Mo4 Si8'

loop_
_publ_author_name
 'M. Fran\c{c}ois'
 'W. Lengauer'
 'K. Yvon'
 'M. Sergent'
 'M. Potel'
 'P. Gougeon'
 'H. {Ben Yaich-Aerrache}'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 196
_journal_year 1991
_journal_page_first 111
_journal_page_last 128
_publ_Section_title
;
 Structural phase transition in GaMo$_{4}$S$_{8}$ by X-ray powder
    ↪ diffraction
;

_aflow_title 'Low-Temperature GaMo$_{4}$S$_{8}$ Structure'
_aflow_proto 'AB4C8_hR13_160_a_ab_2a2b'
_aflow_params 'a,c/a,x_{1},x_{2},x_{3},x_{4},x_{5},z_{5},x_{6},z_{6},x_{
    ↪ 7},z_{7}'
_aflow_params_values '6.90572,2.42010391386,0.0,0.399,0.636,0.135,0.396,
    ↪ 0.814,0.642,0.088,1.138,-0.40801'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hR13'

_symmetry_space_group_name_H-M "R 3 m (hexagonal axes)"
_symmetry_Int_Tables_number 160

_cell_length_a    6.90572
_cell_length_b    6.90572
_cell_length_c    16.71256
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 -y,-x,z
5 x,x-y,z
6 -x+y,y,z
7 x+1/3,y+2/3,z+2/3
8 -y+1/3,x-y+2/3,z+2/3
9 -x+y+1/3,-x+2/3,z+2/3
10 -y+1/3,-x+2/3,z+2/3
11 x+1/3,x-y+2/3,z+2/3
12 -x+y+1/3,y+2/3,z+2/3
13 x+2/3,y+1/3,z+1/3
14 -y+2/3,x-y+1/3,z+1/3
15 -x+y+2/3,-x+1/3,z+1/3
16 -y+2/3,-x+1/3,z+1/3
17 x+2/3,x-y+1/3,z+1/3
18 -x+y+2/3,y+1/3,z+1/3

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
```

```
_atom_site_occupancy
Ga1 Ga 3 a 0.00000 0.00000 0.00000 1.00000
Mo1 Mo 3 a 0.00000 0.00000 0.39900 1.00000
Si1 Si 3 a 0.00000 0.00000 0.63600 1.00000
Si2 Si 3 a 0.00000 0.00000 0.13500 1.00000
Mo2 Mo 9 b 0.19400 0.80600 0.20200 1.00000
Si3 Si 9 b 0.51800 0.48200 0.12400 1.00000
Si4 Si 9 b 0.84867 0.15133 0.28933 1.00000
```

Low-Temperature $GaMo_4S_8$: AB4C8_hR13_160_a_ab_2a2b - POSCAR

```
AB4C8_hR13_160_a_ab_2a2b & a,c/a,x1,x2,x3,x4,x5,z5,x6,z6,x7,z7 --params=
    ↪ 6.90572,2.42010391386,0.0,0.399,0.636,0.135,0.396,0.814,0.642,
    ↪ 0.088,1.138,-0.40801 & R3m C_{3v}^{5} #160 (a^4b^3) & hR13 &
    ↪ None & GaMo4S8 & GaMo4S8 & M. Fran\c{c}ois et al., Zeitschrift
    ↪ f{"u}r Kristallographie - Crystalline Materials 196, 111-128 (
    ↪ 1991)
   1.00000000000000
   3.45286000000000  -1.99350965047409   5.57085333333333
   0.00000000000000   3.98701930094818   5.57085333333333
  -3.45286000000000  -1.99350965047409   5.57085333333333
     Ga    Mo    Si
      1     4     8
Direct
   0.00000000000000   0.00000000000000   0.00000000000000   Ga   (1a)
   0.39900000000000   0.39900000000000   0.39900000000000   Mo   (1a)
   0.39600000000000   0.39600000000000   0.81400000000000   Mo   (3b)
   0.81400000000000   0.39600000000000   0.39600000000000   Mo   (3b)
   0.39600000000000   0.81400000000000   0.39600000000000   Mo   (3b)
   0.63600000000000   0.63600000000000   0.63600000000000   Si   (1a)
   0.13500000000000   0.13500000000000   0.13500000000000   Si   (1a)
   0.64200000000000   0.64200000000000   0.08800000000000   Si   (3b)
   0.08800000000000   0.64200000000000   0.64200000000000   Si   (3b)
   0.64200000000000   0.08800000000000   0.64200000000000   Si   (3b)
   1.13800000000000   1.13800000000000  -0.40801000000000   Si   (3b)
  -0.40801000000000   1.13800000000000   1.13800000000000   Si   (3b)
   1.13800000000000  -0.40801000000000   1.13800000000000   Si   (3b)
```

$KBrO_3$ ($G0_7$): ABC3_hR5_160_a_a_b - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'BrKO3'
_chemical_formula_sum 'Br K O3'

loop_
_publ_author_name
 'D. H. Templeton'
 'L. K. Templeton'
_journal_year 1985
_publ_Section_title
;
 Tensor X-ray optical properties of the bromate ion
;

# Found in Crystal behavior of potassium bromate under compression, 2015

_aflow_title 'KBrO$_{3}$ ($G0_{7}$) Structure'
_aflow_proto 'ABC3_hR5_160_a_a_b'
_aflow_params 'a,c/a,x_{1},x_{2},x_{3},z_{3}'
_aflow_params_values '6.011,1.35618033605,0.4827,0.0,0.54476,1.11108'
_aflow_Strukturbericht '$G0_{7}$'
_aflow_Pearson 'hR5'

_symmetry_space_group_name_H-M "R 3 m (hexagonal axes)"
_symmetry_Int_Tables_number 160

_cell_length_a     6.01100
_cell_length_b     6.01100
_cell_length_c     8.15200
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 -y,-x,z
5 x,x-y,z
6 -x+y,y,z
7 x+1/3,y+2/3,z+2/3
8 -y+1/3,x-y+2/3,z+2/3
9 -x+y+1/3,-x+2/3,z+2/3
10 -y+1/3,-x+2/3,z+2/3
11 x+1/3,x-y+2/3,z+2/3
12 -x+y+1/3,y+2/3,z+2/3
13 x+2/3,y+1/3,z+1/3
14 -y+2/3,x-y+1/3,z+1/3
15 -x+y+2/3,-x+1/3,z+1/3
16 -y+2/3,-x+1/3,z+1/3
17 x+2/3,x-y+1/3,z+1/3
18 -x+y+2/3,y+1/3,z+1/3

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Br1 Br 3 a 0.00000 0.00000 0.48270 1.00000
K1 K 3 a 0.00000 0.00000 0.00000 1.00000
O1 O 9 b 0.14456 0.85544 0.40020 1.00000
```

$KBrO_3$ ($G0_7$): ABC3_hR5_160_a_a_b - POSCAR

```
ABC3_hR5_160_a_a_b & a,c/a,x1,x2,x3,z3 --params=6.011,1.35618033605,
    ↪ 0.4827,0.0,0.54476,1.11108 & R3m C_{3v}^{5} #160 (a^2b) & hR5 &
    ↪ $G0_{7}$ & BrKO3 & BrKO3 & D. H. Templeton and L. K. Templeton
    ↪ , (1985)
   1.00000000000000
   3.00550000000000  -1.73522623404942   2.71733333333333
   0.00000000000000   3.47045246809884   2.71733333333333
  -3.00550000000000  -1.73522623404942   2.71733333333333
     Br    K    O
      1     1     3
Direct
   0.48270000000000   0.48270000000000   0.48270000000000   Br   (1a)
   0.00000000000000   0.00000000000000   0.00000000000000   K   (1a)
   0.54476000000000   0.54476000000000   1.11108000000000   O   (3b)
   1.11108000000000   0.54476000000000   0.54476000000000   O   (3b)
   0.54476000000000   1.11108000000000   0.54476000000000   O   (3b)
```

γ-Potassium Nitrate (KNO_3): ABC3_hR5_160_a_a_b - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'KNO3'
_chemical_formula_sum 'K N O3'

loop_
_publ_author_name
 'J. K. Nimmo'
 'B. W. Lucas'
_journal_year 1976
_publ_Section_title
;
 The crystal structures of $\gamma$-and $\beta$-KNO$_{3}$ and the $\
    ↪ alpha$-$\beta$-$\gamma$ phase transformations
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title '$\gamma$-Potassium Nitrate (KNO$_{3}$) Structure'
_aflow_proto 'ABC3_hR5_160_a_a_b'
_aflow_params 'a,c/a,x_{1},x_{2},x_{3},z_{3}'
_aflow_params_values '5.487,1.66867140514,0.0,0.405,0.565,1.172'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hR5'

_symmetry_space_group_name_H-M "R 3 m (hexagonal axes)"
_symmetry_Int_Tables_number 160

_cell_length_a     5.48700
_cell_length_b     5.48700
_cell_length_c     9.15600
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 -y,-x,z
5 x,x-y,z
6 -x+y,y,z
7 x+1/3,y+2/3,z+2/3
8 -y+1/3,x-y+2/3,z+2/3
9 -x+y+1/3,-x+2/3,z+2/3
10 -y+1/3,-x+2/3,z+2/3
11 x+1/3,x-y+2/3,z+2/3
12 -x+y+1/3,y+2/3,z+2/3
13 x+2/3,y+1/3,z+1/3
14 -y+2/3,x-y+1/3,z+1/3
15 -x+y+2/3,-x+1/3,z+1/3
16 -y+2/3,-x+1/3,z+1/3
17 x+2/3,x-y+1/3,z+1/3
18 -x+y+2/3,y+1/3,z+1/3

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
K1 K 3 a 0.00000 0.00000 0.00000 1.00000
N1 N 3 a 0.00000 0.00000 0.40500 1.00000
O1 O 9 b 0.13100 0.86900 0.43400 1.00000
```

γ-Potassium Nitrate (KNO_3): ABC3_hR5_160_a_a_b - POSCAR

```
ABC3_hR5_160_a_a_b & a,c/a,x1,x2,x3,z3 --params=5.487,1.66867140514,0.0,
    ↪ 0.405,0.565,1.172 & R3m C_{3v}^{5} #160 (a^2b) & hR5 & None &
    ↪ KNO3 & KNO3 & J. K. Nimmo and B. W. Lucas, (1976)
```

```
   1.00000000000000
   2.74350000000000  -1.58396046352174   3.05200000000000
   0.00000000000000   3.16792092704348   3.05200000000000
  -2.74350000000000  -1.58396046352174   3.05200000000000
     K     N     O
     1     1     3
Direct
   0.00000000000000   0.00000000000000   0.00000000000000   K   (1a)
   0.40500000000000   0.40500000000000   0.40500000000000   N   (1a)
   0.56500000000000   0.56500000000000   1.17200000000000   O   (3b)
   1.17200000000000   0.56500000000000   0.56500000000000   O   (3b)
   0.56500000000000   1.17200000000000   0.56500000000000   O   (3b)
```

α-BaB_2O_4 (Low-Temperature): A2BC4_hR42_161_2b_b_4b - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'B2BaO4'
_chemical_formula_sum 'B2 Ba O4'

loop_
_publ_author_name
 'R. Fr\"{o}hlich'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 168
_journal_year 1984
_journal_page_first 109
_journal_page_last 112
_publ_Section_title
;
 Crystal Structure of the low-temperature form of BaB$_{2}$O$_{4}$
;

_aflow_title '$\alpha$-BaB$_{2}$O$_{4}$ (Low-Temperature) Structure'
_aflow_proto 'A2BC4_hR42_161_2b_b_4b'
_aflow_params 'a,c/a,x_{1},y_{1},z_{1},x_{2},y_{2},z_{2},x_{3},y_{3},z_{
    ↪ 3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{6},x_{7},
    ↪ y_{7},z_{7}'
_aflow_params_values '12.519,1.01629523125,0.5427,-0.4113,0.3732,0.2639,
    ↪ 0.3403,0.11859,0.7375,0.01891,-0.62672,0.4582,-0.5851,0.62499,
    ↪ 0.6707,0.2584,-0.4056,0.3591,0.2197,0.14411,0.7788,-0.5047,-
    ↪ 0.07361'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hR42'

_symmetry_space_group_name_H-M "R 3 c (hexagonal axes)"
_symmetry_Int_Tables_number 161

_cell_length_a    12.51900
_cell_length_b    12.51900
_cell_length_c    12.72300
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 -y,-x,z+1/2
5 x,x-y,z+1/2
6 -x+y,y,z+1/2
7 x+1/3,y+2/3,z+2/3
8 -y+1/3,x-y+2/3,z+2/3
9 -x+y+1/3,-x+2/3,z+2/3
10 -y+1/3,-x+2/3,z+1/6
11 x+1/3,x-y+2/3,z+1/6
12 -x+y+1/3,y+2/3,z+1/6
13 x+2/3,y+1/3,z+1/3
14 -y+2/3,x-y+1/3,z+1/3
15 -x+y+2/3,-x+1/3,z+1/3
16 -y+2/3,-x+1/3,z+5/6
17 x+2/3,x-y+1/3,z+5/6
18 -x+y+2/3,y+1/3,z+5/6

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
B1 B 18 b 0.37450 0.57950 0.16820 1.00000
B2 B 18 b 0.02297 -0.09937 0.24093 1.00000
Ba1 Ba 18 b 0.69427 0.02432 0.04323 1.00000
O1 O 18 b 0.29217 0.75113 0.16603 1.00000
O2 O 18 b 0.49620 -0.08390 0.17450 1.00000
O3 O 18 b 0.11813 0.02127 0.24097 1.00000
O4 O 18 b 0.71197 0.57153 0.06683 1.00000
```

α-BaB_2O_4 (Low-Temperature): A2BC4_hR42_161_2b_b_4b - POSCAR

```
A2BC4_hR42_161_2b_b_4b & a,c/a,x1,y1,z1,x2,y2,z2,x3,y3,z3,x4,y4,z4,x5,y5
    ↪ ,z5,x6,y6,z6,x7,y7,z7 --params=12.519,1.01629523125,0.5427,-
    ↪ 0.4113,0.3732,0.2639,0.3403,0.11859,0.7375,0.01891,-0.62672,
    ↪ 0.4582,-0.5851,0.62499,0.6707,0.2584,-0.4056,0.3591,0.2197,
    ↪ 0.14411,0.7788,-0.5047,-0.07361 & R3c C_{3v}^{6} #161 (b^7) &
    ↪ hR42 & None & B2BaO4 & B2BaO4 & R. Fr\"{o}hlich, Zeitschrift f
    ↪ {"u}r Kristallographie - Crystalline Materials 168, 109-112 (
    ↪ 1984)
   1.00000000000000
   6.25950000000000  -3.61392400999246   4.24100000000000
   0.00000000000000   7.22784801998493   4.24100000000000
  -6.25950000000000  -3.61392400999246   4.24100000000000
     B     Ba     O
    12     6     24
Direct
   0.54270000000000  -0.41130000000000   0.37320000000000   B   (6b)
   0.37320000000000   0.54270000000000  -0.41130000000000   B   (6b)
  -0.41130000000000   0.37320000000000   0.54270000000000   B   (6b)
   0.87320000000000   0.08870000000000   1.04270000000000   B   (6b)
   0.08870000000000   1.04270000000000   0.87320000000000   B   (6b)
   1.04270000000000   0.87320000000000   0.08870000000000   B   (6b)
   0.26390000000000   0.34030000000000   0.11859000000000   B   (6b)
   0.11859000000000   0.26390000000000   0.34030000000000   B   (6b)
   0.34030000000000   0.11859000000000   0.26390000000000   B   (6b)
   0.61859000000000   0.84030000000000   0.76390000000000   B   (6b)
   0.84030000000000   0.76390000000000   0.61859000000000   B   (6b)
   0.76390000000000   0.61859000000000   0.84030000000000   B   (6b)
   0.73750000000000   0.01891000000000  -0.62672000000000   Ba  (6b)
  -0.62672000000000   0.73750000000000   0.01891000000000   Ba  (6b)
   0.01891000000000  -0.62672000000000   0.73750000000000   Ba  (6b)
  -0.12672000000000   0.51891000000000   1.23750000000000   Ba  (6b)
   0.51891000000000   1.23750000000000  -0.12672000000000   Ba  (6b)
   1.23750000000000  -0.12672000000000   0.51891000000000   Ba  (6b)
   0.45820000000000  -0.58510000000000   0.62499000000000   O   (6b)
   0.62499000000000   0.45820000000000  -0.58510000000000   O   (6b)
  -0.58510000000000   0.62499000000000   0.45820000000000   O   (6b)
   1.12499000000000  -0.08510000000000   0.95820000000000   O   (6b)
  -0.08510000000000   0.95820000000000   1.12499000000000   O   (6b)
   0.95820000000000   1.12499000000000  -0.08510000000000   O   (6b)
   0.67070000000000   0.25840000000000  -0.40560000000000   O   (6b)
  -0.40560000000000   0.67070000000000   0.25840000000000   O   (6b)
   0.25840000000000  -0.40560000000000   0.67070000000000   O   (6b)
   0.09440000000000   0.75840000000000   1.17070000000000   O   (6b)
   0.75840000000000   1.17070000000000   0.09440000000000   O   (6b)
   1.17070000000000   0.09440000000000   0.75840000000000   O   (6b)
   0.35910000000000   0.21970000000000   0.14411000000000   O   (6b)
   0.14411000000000   0.35910000000000   0.21970000000000   O   (6b)
   0.21970000000000   0.14411000000000   0.35910000000000   O   (6b)
   0.64411000000000   0.71970000000000   0.85910000000000   O   (6b)
   0.71970000000000   0.85910000000000   0.64411000000000   O   (6b)
   0.85910000000000   0.64411000000000   0.71970000000000   O   (6b)
   0.77880000000000  -0.50470000000000  -0.07361000000000   O   (6b)
  -0.07361000000000   0.77880000000000  -0.50470000000000   O   (6b)
  -0.50470000000000  -0.07361000000000   0.77880000000000   O   (6b)
   0.42639000000000  -0.00470000000000   1.27880000000000   O   (6b)
  -0.00470000000000   1.27880000000000   0.42639000000000   O   (6b)
   1.27880000000000   0.42639000000000  -0.00470000000000   O   (6b)
```

$I1_3$ ($SrCl_2 \cdot (H_2O)_6$) (*obsolete*): A2B6C_hP9_162_d_k_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Cl2(H2O)6Sr'
_chemical_formula_sum 'Cl2 (H2O)6 Sr'

loop_
_publ_author_name
 'Z. Herrmann'
_journal_name_full_name
;
 Zeitschrift fur Anorganische und Allgemeine Chemie
;
_journal_volume 187
_journal_year 1930
_journal_page_first 231
_journal_page_last 236
_publ_Section_title
;
 {\"U}ber die Struktur des Strontiumchlorid-Hexahydrats
;

# Found in Strukturbericht Band II 1928-1932, 1937

_aflow_title '$I1_{3}$ ($SrCl$_{2}$$\cdot$(H$_{2}$O)$_{6}$) ({\em{
    ↪ obsolete}}) Structure'
_aflow_proto 'A2B6C_hP9_162_d_k_a'
_aflow_params 'a,c/a,x_{3},z_{3}'
_aflow_params_values '7.906,0.5147988869210,0.387,0.0162'
_aflow_Strukturbericht '$I1_{3}$'
_aflow_Pearson 'hP9'

_symmetry_space_group_name_H-M "P -3 1 2/m"
_symmetry_Int_Tables_number 162

_cell_length_a    7.90600
_cell_length_b    7.90600
_cell_length_c    4.07000
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
```

```
4 x,x-y,-z
5 -x+y,y,-z
6 -y,-x,-z
7 -x,-y,-z
8 y,-x+y,-z
9 x-y,x,-z
10 -x,-x+y,z
11 x-y,-y,z
12 y,x,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Sr1 Sr 1 a 0.00000 0.00000 0.00000 1.00000
Cl1 Cl 2 d 0.33333 0.66667 0.50000 1.00000
H2O1 H2O 6 k 0.38700 0.38700 0.01620 1.00000
```

$I1_3$ ($SrCl_2{\cdot}(H_2O)_6$) (*obsolete*): A2B6C_hP9_162_d_k_a - POSCAR

```
A2B6C_hP9_162_d_k_a & a,c/a,x3,z3 --params=7.906,0.514798886921,0.387,
    ↪ 0.0162 & P-31m D_{3d}^{1} #162 (adk) & hP9 & $I1_{3}$ & Cl2(H2O
    ↪ )6Sr & Cl2(H2O)6Sr & Z. Herrmann, Z. Anorg. Allg. Chem. 187,
    ↪ 231-236 (1930)
   1.00000000000000
   3.95300000000000  -6.84679684231977   0.00000000000000
   3.95300000000000   6.84679684231977   0.00000000000000
   0.00000000000000   0.00000000000000   4.07000000000000
     Cl    H2O     Sr
      2      6      1
Direct
   0.33333333333333   0.66666666666667   0.50000000000000    Cl    (2d)
   0.66666666666667   0.33333333333333   0.50000000000000    Cl    (2d)
   0.38700000000000   0.00000000000000   0.01620000000000   H2O    (6k)
   0.00000000000000   0.38700000000000   0.01620000000000   H2O    (6k)
  -0.38700000000000  -0.38700000000000   0.01620000000000   H2O    (6k)
   0.00000000000000  -0.38700000000000  -0.01620000000000   H2O    (6k)
  -0.38700000000000   0.00000000000000  -0.01620000000000   H2O    (6k)
   0.38700000000000   0.38700000000000  -0.01620000000000   H2O    (6k)
   0.00000000000000   0.00000000000000   0.00000000000000    Sr    (1a)
```

Rosiaite ($PbSb_2O_6$): A6BC2_hP9_162_k_a_d - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Rosiaite'
_chemical_formula_sum 'O6 Pb Sb2'

loop_
_publ_author_name
 'R. Basso'
 'G. Lucchetti'
 'L. Zefiro'
 'A. Palenzona'
_journal_name_full_name
;
 European Journal of Mineralogy
;
_journal_volume 8
_journal_year 1996
_journal_page_first 487
_journal_page_last 492
_publ_Section_title
;
 Rosiaite, PbSb$_{2}$O$_{6}$, a new mineral from the Cetine mine, Siena,
     ↪ Italy
;

_aflow_title 'Rosiaite (PbSb$_{2}$O$_{6}$) Structure'
_aflow_proto 'A6BC2_hP9_162_k_a_d'
_aflow_params 'a,c/a,x_{3},z_{3}'
_aflow_params_values '5.295,1.01454202077,0.377,0.2965'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hP9'

_symmetry_space_group_name_H-M "P -3 1 2/m"
_symmetry_Int_Tables_number 162

_cell_length_a     5.29500
_cell_length_b     5.29500
_cell_length_c     5.37200
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 x,x-y,-z
5 -x+y,y,-z
6 -y,-x,-z
7 -x,-y,-z
8 y,-x+y,-z
9 x-y,x,-z
10 -x,-x+y,z
11 x-y,-y,z
12 y,x,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Pb1 Pb 1 a 0.00000 0.00000 0.00000 1.00000
Sb1 Sb 2 d 0.33333 0.66667 0.50000 1.00000
O1 O 6 k 0.37700 0.00000 0.29650 1.00000
```

Rosiaite ($PbSb_2O_6$): A6BC2_hP9_162_k_a_d - POSCAR

```
A6BC2_hP9_162_k_a_d & a,c/a,x3,z3 --params=5.295,1.01454202077,0.377,
    ↪ 0.2965 & P-31m D_{3d}^{1} #162 (adk) & hP9 & None & O6PbSb2 &
    ↪ Rosiaite & R. Basso et al., Eur. J. Mineral. 8, 487-492 (1996)
   1.00000000000000
   2.64750000000000  -4.58560451303860   0.00000000000000
   2.64750000000000   4.58560451303860   0.00000000000000
   0.00000000000000   0.00000000000000   5.37200000000000
     O     Pb     Sb
     6      1      2
Direct
   0.37700000000000   0.00000000000000   0.29650000000000     O    (6k)
   0.00000000000000   0.37700000000000   0.29650000000000     O    (6k)
  -0.37700000000000  -0.37700000000000   0.29650000000000     O    (6k)
   0.00000000000000  -0.37700000000000  -0.29650000000000     O    (6k)
  -0.37700000000000   0.00000000000000  -0.29650000000000     O    (6k)
   0.37700000000000   0.37700000000000  -0.29650000000000     O    (6k)
   0.00000000000000   0.00000000000000   0.00000000000000    Pb    (1a)
   0.33333333333333   0.66666666666667   0.50000000000000    Sb    (2d)
   0.66666666666667   0.33333333333333   0.50000000000000    Sb    (2d)
```

$NaSbF_4(OH)_2$ ($J1_{12}$): A6BC_hP16_163_i_b_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'F4NaSb'
_chemical_formula_sum 'F6 Na Sb'

loop_
_publ_author_name
 'N. Schrewelius'
_journal_name_full_name
;
 Zeitschrift fur Anorganische und Allgemeine Chemie
;
_journal_volume 238
_journal_year 1938
_journal_page_first 241
_journal_page_last 254
_publ_Section_title
;
 R\"{o}ntgenuntersuchung der Verbindungen NaSb(OH)$_{6}$, NaSbF$_{6}$,
     ↪ NaSbO$_{3}$ und gleichartiger Stoffe
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'NaSbF$_{4}$(OH)$_{2}$ ($J1_{12}$) Structure'
_aflow_proto 'A6BC_hP16_163_i_b_c'
_aflow_params 'a,c/a,x_{3},y_{3},z_{3}'
_aflow_params_values '5.227,1.90931700784,0.33,0.33,0.15'
_aflow_Strukturbericht '$J1_{12}$'
_aflow_Pearson 'hP16'

_symmetry_space_group_name_H-M "P -3 1 2/c"
_symmetry_Int_Tables_number 163

_cell_length_a     5.22700
_cell_length_b     5.22700
_cell_length_c     9.98000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 x,x-y,-z+1/2
5 -x+y,y,-z+1/2
6 -y,-x,-z+1/2
7 -x,-y,-z
8 y,-x+y,-z
9 x-y,x,-z
10 -x,-x+y,z+1/2
11 x-y,-y,z+1/2
12 y,x,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
```

```
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Na1 Na 2 b 0.00000 0.00000 0.00000 1.00000
Sb1 Sb 2 c 0.33333 0.66667 0.25000 1.00000
F1 F 12 i 0.33000 0.33000 0.15000 1.00000
```

$NaSbF_4(OH)_2$ ($J1_{12}$): A6BC_hP16_163_i_b_c - POSCAR

```
A6BC_hP16_163_i_b_c & a,c/a,x3,y3,z3 --params=5.227,1.90931700784,0.33,
    ↪ 0.33,0.15 & P-31c D_{3d}^{2} #163 (bci) & hP16 & $J1_{12}$ &
    ↪ F4NaSb & F4NaSb & N. Schrewelius, Z. Anorg. Allg. Chem. 238,
    ↪ 241-254 (1938)
   1.00000000000000
   2.61350000000000  -4.52671478558126   0.00000000000000
   2.61350000000000   4.52671478558126   0.00000000000000
   0.00000000000000   0.00000000000000   9.98000000000000
     F    Na    Sb
    12     2     2
Direct
   0.33000000000000   0.33000000000000   0.15000000000000   F   (12i)
  -0.33000000000000   0.00000000000000   0.15000000000000   F   (12i)
   0.00000000000000  -0.33000000000000   0.15000000000000   F   (12i)
  -0.33000000000000  -0.33000000000000   0.35000000000000   F   (12i)
   0.00000000000000   0.33000000000000   0.35000000000000   F   (12i)
   0.33000000000000   0.00000000000000   0.35000000000000   F   (12i)
  -0.33000000000000  -0.33000000000000  -0.15000000000000   F   (12i)
   0.33000000000000   0.00000000000000  -0.15000000000000   F   (12i)
   0.00000000000000   0.33000000000000  -0.15000000000000   F   (12i)
   0.33000000000000   0.33000000000000   0.65000000000000   F   (12i)
   0.00000000000000  -0.33000000000000   0.65000000000000   F   (12i)
  -0.33000000000000   0.00000000000000   0.65000000000000   F   (12i)
   0.00000000000000   0.00000000000000   0.00000000000000   Na   (2b)
   0.00000000000000   0.00000000000000   0.50000000000000   Na   (2b)
   0.33333333333333   0.66666666666667   0.25000000000000   Sb   (2c)
   0.66666666666667   0.33333333333333   0.75000000000000   Sb   (2c)
```

Colquiriite ($LiCaAlF_6$): ABC6D_hP18_163_d_b_i_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Colquiriite'
_chemical_formula_sum 'Al Ca F6 Li'

loop_
_publ_author_name
 'S. Kuze'
 'D. du Boulay'
 'N. Ishizawa'
 'N. Kodama'
 'M. Yamaga'
 'B. Henderson'
_journal_name_full_name
;
 Journal of Solid State Chemistry
;
_journal_volume 177
_journal_year 2004
_journal_page_first 3505
_journal_page_last 3513
_publ_Section_title
;
 Structures of LiCaAlF${6}$ and LiSrAlF${6}$ at 120 and 300 K by
    ↪ synchrotron X-ray single-crystal diffraction
;

_aflow_title 'Colquiriite (LiCaAlF${6}$) Structure'
_aflow_proto 'ABC6D_hP18_163_d_b_i_c'
_aflow_params 'a,c/a,x_{4},y_{4},z_{4}'
_aflow_params_values '5.0081,1.92554062419,0.3767,0.031,0.14336'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hP18'

_symmetry_space_group_name_H-M "P -3 1 2/c"
_symmetry_Int_Tables_number 163

_cell_length_a    5.00810
_cell_length_b    5.00810
_cell_length_c    9.64330
_cell_angle_alpha  90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 x,x-y,-z+1/2
5 -x+y,y,-z+1/2
6 -y,-x,-z+1/2
7 -x,-y,-z
8 y,-x+y,-z
9 x-y,x,-z
10 -x,-x+y,z+1/2
11 x-y,-y,z+1/2
12 y,x,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ca1 Ca 2 b 0.00000 0.00000 0.00000 1.00000
Li1 Li 2 c 0.33333 0.66667 0.25000 1.00000
Al1 Al 2 d 0.66667 0.33333 0.25000 1.00000
F1 F 12 i 0.37670 0.03100 0.14336 1.00000
```

Colquiriite ($LiCaAlF_6$): ABC6D_hP18_163_d_b_i_c - POSCAR

```
ABC6D_hP18_163_d_b_i_c & a,c/a,x4,y4,z4 --params=5.0081,1.92554062419,
    ↪ 0.3767,0.031,0.14336 & P-31c D_{3d}^{2} #163 (bcdi) & hP18 &
    ↪ None & AlCaF6Li & Colquiriite & S. Kuze et al., J. Solid State
    ↪ Chem. 177, 3505-3513 (2004)
   1.00000000000000
   2.50405000000000  -4.33714182469285   0.00000000000000
   2.50405000000000   4.33714182469285   0.00000000000000
   0.00000000000000   0.00000000000000   9.64330000000000
    Al    Ca     F    Li
     2     2    12     2
Direct
   0.66666666666667   0.33333333333333   0.25000000000000   Al   (2d)
   0.33333333333333   0.66666666666667   0.75000000000000   Al   (2d)
   0.00000000000000   0.00000000000000   0.00000000000000   Ca   (2b)
   0.00000000000000   0.00000000000000   0.50000000000000   Ca   (2b)
   0.37670000000000   0.03100000000000   0.14336000000000    F   (12i)
  -0.03100000000000   0.34570000000000   0.14336000000000    F   (12i)
  -0.34570000000000  -0.37670000000000   0.14336000000000    F   (12i)
  -0.03100000000000  -0.37670000000000   0.35664000000000    F   (12i)
  -0.34570000000000   0.03100000000000   0.35664000000000    F   (12i)
   0.37670000000000   0.34570000000000   0.35664000000000    F   (12i)
  -0.37670000000000  -0.03100000000000  -0.14336000000000    F   (12i)
   0.03100000000000  -0.34570000000000  -0.14336000000000    F   (12i)
   0.34570000000000   0.37670000000000  -0.14336000000000    F   (12i)
   0.03100000000000   0.37670000000000   0.64336000000000    F   (12i)
   0.34570000000000  -0.03100000000000   0.64336000000000    F   (12i)
  -0.37670000000000  -0.34570000000000   0.64336000000000    F   (12i)
   0.33333333333333   0.66666666666667   0.25000000000000   Li   (2c)
   0.66666666666667   0.33333333333333   0.75000000000000   Li   (2c)
```

Predicted Li_2MgH_{16} 300 GPa: A16B2C_hP19_164_2d2i_d_b - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'H16Li2Mg'
_chemical_formula_sum 'H16 Li2 Mg'

loop_
_publ_author_name
 'Y. Sun'
 'J. Lv'
 'Y. Xie'
 'H. Liu'
 'Y. Ma'
_journal_name_full_name
;
 Physical Review Letters
;
_journal_volume 123
_journal_year 2019
_journal_page_first 097001
_journal_page_last 097001
_publ_Section_title
;
 Route to a Superconducting Phase above Room Temperature in
    ↪ Electron-Doped Hydride Compounds under High Pressure
;

_aflow_title 'Predicted Li${2}$MgH${16}$ 300~GPa Structure'
_aflow_proto 'A16B2C_hP19_164_2d2i_d_b'
_aflow_params 'a,c/a,z_{2},z_{3},z_{4},x_{5},z_{5},x_{6},z_{6}'
_aflow_params_values '2.79596,1.9003812644,0.57451,0.39139,0.83626,
    ↪ 0.17157,0.0797,0.1644,0.23653'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hP19'

_symmetry_space_group_name_H-M "P -3 2/m 1"
_symmetry_Int_Tables_number 164

_cell_length_a    2.79596
_cell_length_b    2.79596
_cell_length_c    5.31339
_cell_angle_alpha  90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 x-y,-y,-z
5 y,x,-z
6 -x,-x+y,-z
7 -x,-y,-z
8 y,-x+y,-z
9 x-y,x,-z
10 -x+y,y,z
11 -y,-x,z
12 x,x-y,z
```

```
loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Mg1 Mg 1 b 0.00000 0.00000 0.50000 1.00000
H1 H 2 d 0.33333 0.66667 0.57451 1.00000
H2 H 2 d 0.33333 0.66667 0.39139 1.00000
Li1 Li 2 d 0.33333 0.66667 0.83626 1.00000
H3 H 6 i 0.17157 0.82843 0.07970 1.00000
H4 H 6 i 0.16440 0.83560 0.23653 1.00000
```

Predicted Li_2MgH_{16} 300 GPa: A16B2C_hP19_164_2d2i_d_b - POSCAR

```
A16B2C_hP19_164_2d2i_d_b & a,c/a,z2,z3,z4,x5,z5,x6,z6 --params=2.79596,
    ↪ 1.9003812644,0.57451,0.39139,0.83626,0.17157,0.0797,0.1644,
    ↪ 0.23653 & P-3m1 D_{3d}^{3} #164 (bd^3i^2) & hP19 & None &
    ↪ H16Li2Mg & H16Li2Mg & Y. Sun et al., Phys. Rev. Lett. 123,
    ↪ 097001(2019)
   1.00000000000000
   1.39798000000000  -2.42137238796514   0.00000000000000
   1.39798000000000   2.42137238796514   0.00000000000000
   0.00000000000000   0.00000000000000   5.31339000000000
     H     Li     Mg
    16     2      1
Direct
   0.33333333333333   0.66666666666667   0.57451000000000   H   (2d)
   0.66666666666667   0.33333333333333  -0.57451000000000   H   (2d)
   0.33333333333333   0.66666666666667   0.39139000000000   H   (2d)
   0.66666666666667   0.33333333333333  -0.39139000000000   H   (2d)
   0.17157000000000  -0.17157000000000   0.07970000000000   H   (6i)
   0.17157000000000   0.34314000000000   0.07970000000000   H   (6i)
  -0.34314000000000  -0.17157000000000   0.07970000000000   H   (6i)
  -0.17157000000000   0.17157000000000  -0.07970000000000   H   (6i)
   0.34314000000000   0.17157000000000  -0.07970000000000   H   (6i)
  -0.17157000000000  -0.34314000000000  -0.07970000000000   H   (6i)
   0.16440000000000  -0.16440000000000   0.23653000000000   H   (6i)
   0.16440000000000   0.32880000000000   0.23653000000000   H   (6i)
  -0.32880000000000  -0.16440000000000   0.23653000000000   H   (6i)
  -0.16440000000000   0.16440000000000  -0.23653000000000   H   (6i)
   0.32880000000000   0.16440000000000  -0.23653000000000   H   (6i)
  -0.16440000000000  -0.32880000000000  -0.23653000000000   H   (6i)
   0.33333333333333   0.66666666666667   0.83626000000000   Li  (2d)
   0.66666666666667   0.33333333333333  -0.83626000000000   Li  (2d)
   0.00000000000000   0.00000000000000   0.50000000000000   Mg  (1b)
```

Ce_2O_2S: A2B2C_hP5_164_d_d_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Ce2O2S'
_chemical_formula_sum 'Ce2 O2 S'

loop_
_publ_author_name
 'W. H. Zachariasen'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 2
_journal_year 1949
_journal_page_first 60
_journal_page_last 62
_publ_Section_title
;
 Crystal chemical studies of the $5f$-series of elements. VII. The
    ↪ crystal structure of Ce$_{2}$O$_{2}$S, La$_{2}$O$_{2}$S and
    ↪ Pu$_{2}$O$_{2}$S
;

_aflow_title 'Ce$_{2}$O$_{2}$S Structure'
_aflow_proto 'A2B2C_hP5_164_d_d_a'
_aflow_params 'a,c/a,z_{2},z_{3}'
_aflow_params_values '4.0,1.705,0.29,0.64'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hP5'

_symmetry_space_group_name_H-M "P -3 2/m 1"
_symmetry_Int_Tables_number 164

_cell_length_a     4.00000
_cell_length_b     4.00000
_cell_length_c     6.82000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 x-y,-y,-z
5 y,x,-z
6 -x,-x+y,-z
7 -x,-y,-z
8 y,-x+y,-z
9 x-y,x,-z
10 -x+y,y,z
11 -y,-x,z
12 x,x-y,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
S1 S 1 a 0.00000 0.00000 0.00000 1.00000
Ce1 Ce 2 d 0.33333 0.66667 0.29000 1.00000
O1 O 2 d 0.33333 0.66667 0.64000 1.00000
```

Ce_2O_2S: A2B2C_hP5_164_d_d_a - POSCAR

```
A2B2C_hP5_164_d_d_a & a,c/a,z2,z3 --params=4.0,1.705,0.29,0.64 & P-3m1
    ↪ D_{3d}^{3} #164 (ad^2) & hP5 & None & Ce2O2S & Ce2O2S & W. H.
    ↪ Zachariasen, Acta Cryst. 2, 60-62 (1949)
   1.00000000000000
   2.00000000000000  -3.46410161513775   0.00000000000000
   2.00000000000000   3.46410161513775   0.00000000000000
   0.00000000000000   0.00000000000000   6.82000000000000
    Ce     O     S
     2     2     1
Direct
   0.33333333333333   0.66666666666667   0.29000000000000   Ce   (2d)
   0.66666666666667   0.33333333333333  -0.29000000000000   Ce   (2d)
   0.33333333333333   0.66666666666667   0.64000000000000   O    (2d)
   0.66666666666667   0.33333333333333  -0.64000000000000   O    (2d)
   0.00000000000000   0.00000000000000   0.00000000000000   S    (1a)
```

Brucite [$Mg(OH)_2$]: A2BC2_hP5_164_d_a_d - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Brucite'
_chemical_formula_sum 'H2 Mg O2'

loop_
_publ_author_name
 'M. Catti'
 'G. Ferraris'
 'S. Hull'
 'A. Pavese'
_journal_name_full_name
;
 Physics and Chemistry of Minerals
;
_journal_volume 22
_journal_year 1995
_journal_page_first 200
_journal_page_last 206
_publ_Section_title
;
 Static compression and H disorder in brucite, Mg(OH)$_{2}$, to 11 GPa:
    ↪ a powder neutron diffraction study
;

_aflow_title 'Brucite [Mg(OH)$_{2}$] Structure'
_aflow_proto 'A2BC2_hP5_164_d_a_d'
_aflow_params 'a,c/a,z_{2},z_{3}'
_aflow_params_values '3.14979,1.51445016969,0.413,0.2203'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hP5'

_symmetry_space_group_name_H-M "P -3 2/m 1"
_symmetry_Int_Tables_number 164

_cell_length_a     3.14979
_cell_length_b     3.14979
_cell_length_c     4.77020
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 x-y,-y,-z
5 y,x,-z
6 -x,-x+y,-z
7 -x,-y,-z
8 y,-x+y,-z
9 x-y,x,-z
10 -x+y,y,z
11 -y,-x,z
12 x,x-y,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
```

```
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Mg1 Mg 1 a 0.00000 0.00000 0.00000 1.00000
H1 H 2 d 0.33333 0.66667 0.41300 1.00000
O1 O 2 d 0.33333 0.66667 0.22030 1.00000
```

Brucite [$Mg(OH)_2$]: A2BC2_hP5_164_d_a_d - POSCAR

```
A2BC2_hP5_164_d_a_d & a,c/a,z2,z3 --params=3.14979,1.51445016969,0.413,
    ↪ 0.2203 & P-3m1 D_{3d}^{3} #164 (ad^2) & hP5 & None & H2MgO2 &
    ↪ Brucite & M. Catti et al., Phys. Chem. Miner. 22, 200-206 (1995
    ↪ )
   1.00000000000000
   1.57489500000000  -2.72779815658619   0.00000000000000
   1.57489500000000   2.72779815658619   0.00000000000000
   0.00000000000000   0.00000000000000   4.77020000000000
     H     Mg     O
     2      1      2
Direct
   0.33333333333333   0.66666666666667   0.41300000000000   H   (2d)
   0.66666666666667   0.33333333333333  -0.41300000000000   H   (2d)
   0.00000000000000   0.00000000000000   0.00000000000000   Mg  (1a)
   0.33333333333333   0.66666666666667   0.22030000000000   O   (2d)
   0.66666666666667   0.33333333333333  -0.22030000000000   O   (2d)
```

$K_2Pt(SCN)_6$ ($H6_3$): A2BC6_hP9_164_d_a_i - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'K2Pt(SCN)6'
_chemical_formula_sum 'K2 Pt S6'

loop_
_publ_author_name
 'S. B. Hendricks'
 'H. E. Merwin'
_journal_name_full_name
;
 American Journal of Science
;
_journal_volume 15
_journal_year 1928
_journal_page_first 487
_journal_page_last 494
_publ_Section_title
;
 The atomic arrangement in crystals of alkali platini-thiocyanates
;

# Found in Strukturbericht 1913-1928, 1931

_aflow_title 'K${2}$Pt(SCN)${6}$ ($H6_{3}$) Structure'
_aflow_proto 'A2BC6_hP9_164_d_a_i'
_aflow_params 'a,c/a,z_{2},x_{3},z_{3}'
_aflow_params_values '6.73,1.52451708767,0.5,0.135,0.1125'
_aflow_Strukturbericht '$H6_{3}$'
_aflow_Pearson 'hP9'

_symmetry_space_group_name_H-M "P -3 2/m 1"
_symmetry_Int_Tables_number 164

_cell_length_a     6.73000
_cell_length_b     6.73000
_cell_length_c     10.26000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 x-y,-y,-z
5 y,x,-z
6 -x,-x+y,-z
7 -x,-y,-z
8 y,-x+y,-z
9 x-y,x,-z
10 -x+y,y,z
11 -y,-x,z
12 x,x-y,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Pt1 Pt 1 a 0.00000 0.00000 0.00000 1.00000
K1 K 2 d 0.33333 0.66667 0.50000 1.00000
S1 S 6 i 0.13500 0.86500 0.11250 1.00000
```

$K_2Pt(SCN)_6$ ($H6_3$): A2BC6_hP9_164_d_a_i - POSCAR

```
A2BC6_hP9_164_d_a_i & a,c/a,z2,x3,z3 --params=6.73,1.52451708767,0.5,
    ↪ 0.135,0.1125 & P-3m1 D_{3d}^{3} #164 (adi) & hP9 & $H6_{3}$ &
    ↪ K2Pt(SCN)6 & K2Pt(SCN)6 & S. B. Hendricks and H. E. Merwin, Am.
    ↪ J. Sci. 15, 487-494 (1928)
   1.00000000000000
   3.36500000000000  -5.82835096746927   0.00000000000000
   3.36500000000000   5.82835096746927   0.00000000000000
   0.00000000000000   0.00000000000000  10.26000000000000
     K     Pt     S
     2      1      6
Direct
   0.33333333333333   0.66666666666667   0.50000000000000   K   (2d)
   0.66666666666667   0.33333333333333  -0.50000000000000   K   (2d)
   0.00000000000000   0.00000000000000   0.00000000000000   Pt  (1a)
   0.13500000000000  -0.13500000000000   0.11250000000000   S   (6i)
   0.13500000000000   0.27000000000000   0.11250000000000   S   (6i)
  -0.27000000000000  -0.13500000000000   0.11250000000000   S   (6i)
  -0.13500000000000   0.13500000000000  -0.11250000000000   S   (6i)
   0.27000000000000   0.13500000000000  -0.11250000000000   S   (6i)
  -0.13500000000000  -0.27000000000000  -0.11250000000000   S   (6i)
```

Bararite (Trigonal $(NH_4)_2SiF_6$, $J1_6$): A6B2C_hP9_164_i_d_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Bararite'
_chemical_formula_sum 'F6 (NH4)2 Si'

loop_
_publ_author_name
 'E. O. Schlemper'
 'W. C. Hamilton'
_journal_name_full_name
;
 Journal of Chemical Physics
;
_journal_volume 45
_journal_year 1966
_journal_page_first 408
_journal_page_last 409
_publ_Section_title
;
 On the Structure of Trigonal Ammonium Flourosilicate
;

# Found in A new modification of diammonium hexafluorosilicate, (NH${4}
    ↪ $)${2}$SiF${6}$, 2001

_aflow_title 'Bararite (Trigonal (NH${4}$)${2}$SiF${6}$, $J1_{6}$)
    ↪ Structure'
_aflow_proto 'A6B2C_hP9_164_i_d_a'
_aflow_params 'a,c/a,z_{2},x_{3},z_{3}'
_aflow_params_values '5.784,0.82918395574,0.33,0.139,0.22'
_aflow_Strukturbericht '$J1_{6}$'
_aflow_Pearson 'hP9'

_symmetry_space_group_name_H-M "P -3 2/m 1"
_symmetry_Int_Tables_number 164

_cell_length_a     5.78400
_cell_length_b     5.78400
_cell_length_c     4.79600
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 x-y,-y,-z
5 y,x,-z
6 -x,-x+y,-z
7 -x,-y,-z
8 y,-x+y,-z
9 x-y,x,-z
10 -x+y,y,z
11 -y,-x,z
12 x,x-y,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Si1 Si 1 a 0.00000 0.00000 0.00000 1.00000
NH41 NH4 2 d 0.33333 0.66667 0.33000 1.00000
F1 F 6 i 0.13900 0.86100 0.22000 1.00000
```

Bararite (Trigonal $(NH_4)_2SiF_6$, $J1_6$): A6B2C_hP9_164_i_d_a - POSCAR

```
A6B2C_hP9_164_i_d_a & a,c/a,z2,x3,z3 --params=5.784,0.82918395574,0.33,
    ↪ 0.139,0.22 & P-3m1 D_{3d}^{3} #164 (adi) & hP9 & $J1_{6}$ & Fe6
    ↪ (NH4)2Si & Bararite & E. O. Schlemper and W. C. Hamilton, J.
    ↪ Chem. Phys. 45, 408-409 (1966)
   1.00000000000000
   2.89200000000000  -5.00909093548919   0.00000000000000
```

```
    2.89200000000000    5.00909093548919    0.00000000000000
    0.00000000000000    0.00000000000000    4.79600000000000
      F    NH4     Si
      6      2      1
Direct
    0.13900000000000   -0.13900000000000    0.22000000000000     F    (6i)
    0.13900000000000    0.27800000000000    0.22000000000000     F    (6i)
   -0.27800000000000   -0.13900000000000    0.22000000000000     F    (6i)
   -0.13900000000000    0.13900000000000   -0.22000000000000     F    (6i)
    0.27800000000000    0.13900000000000   -0.22000000000000     F    (6i)
   -0.13900000000000   -0.27800000000000   -0.22000000000000     F    (6i)
    0.33333333333333    0.66666666666667    0.33000000000000    NH4   (2d)
    0.66666666666667    0.33333333333333   -0.33000000000000    NH4   (2d)
    0.00000000000000    0.00000000000000    0.00000000000000     Si   (1a)
```

K_2GeF_6 ($J1_{13}$): A6BC2_hP9_164_i_a_d - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'F6GeK2'
_chemical_formula_sum 'F6 Ge K2'

loop_
_publ_author_name
 'J. L. Hoard'
 'W. B. Vincent'
_journal_name_full_name
;
 Journal of the American Chemical Society
;
_journal_volume 61
_journal_year 1939
_journal_page_first 2849
_journal_page_last 2852
_publ_Section_title
;
 Structures of Complex Fluorides. Potassium Hexafluogermanate and
      ↪ Ammonium Hexafluogermanate
;

# Found in The crystal structure of the compound Cs$_{2}$CeCl$_{6}$,
      ↪ 1966

_aflow_title 'K$_{2}$GeF$_{6}$ ($J1_{13}$) Structure'
_aflow_proto 'A6BC2_hP9_164_i_a_d'
_aflow_params 'a,c/a,z_{2},x_{3},z_{3}'
_aflow_params_values '5.62,0.827402135231,0.7,0.148,0.22'
_aflow_Strukturbericht '$J1_{13}$'
_aflow_Pearson 'hP9'

_symmetry_space_group_name_H-M "P -3 2/m 1"
_symmetry_Int_Tables_number 164

_cell_length_a     5.62000
_cell_length_b     5.62000
_cell_length_c     4.65000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 x-y,-y,-z
5 y,x,-z
6 -x,-x+y,-z
7 -x,-y,-z
8 y,-x+y,-z
9 x-y,x,-z
10 -x+y,y,z
11 -y,-x,z
12 x,x-y,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ge1 Ge 1 a 0.00000 0.00000 0.00000 1.00000
K1 K 2 d 0.33333 0.66667 0.70000 1.00000
F1 F 6 i 0.14800 0.29600 0.22000 1.00000
```

K_2GeF_6 ($J1_{13}$): A6BC2_hP9_164_i_a_d - POSCAR

```
A6BC2_hP9_164_i_a_d & a,c/a,z2,x3,z3 --params=5.62,0.827402135231,0.7,
      ↪ 0.148,0.22 & P-3m1 D_{3d}^{3} #164 (adi) & hP9 & $J1_{13}$ &
      ↪ F6GeK2 & F6GeK2 & J. L. Hoard and W. B. Vincent, J. Am. Chem.
      ↪ Soc. 61, 2849-2852 (1939)
   1.00000000000000
   2.81000000000000   -4.86706276926854    0.00000000000000
   2.81000000000000    4.86706276926854    0.00000000000000
   0.00000000000000    0.00000000000000    4.65000000000000
     F     Ge     K
     6      1      2
Direct
   0.14800000000000   -0.14800000000000    0.22000000000000     F    (6i)
   0.14800000000000    0.29600000000000    0.22000000000000     F    (6i)
  -0.29600000000000   -0.14800000000000    0.22000000000000     F    (6i)
  -0.14800000000000    0.14800000000000   -0.22000000000000     F    (6i)
   0.29600000000000    0.14800000000000   -0.22000000000000     F    (6i)
  -0.14800000000000   -0.29600000000000   -0.22000000000000     F    (6i)
   0.00000000000000    0.00000000000000    0.00000000000000    Ge    (1a)
   0.33333333333333    0.66666666666667    0.70000000000000     K    (2d)
   0.66666666666667    0.33333333333333   -0.70000000000000     K    (2d)
```

Jacutingaite (Pt_2HgSe_3): AB2C3_hP12_164_d_ae_i - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Jacutingaite'
_chemical_formula_sum 'Hg Pt2 Se3'

loop_
_publ_author_name
 'A. Vymazalov\\'{a}'
 'F. Laufek'
 'M. Dr\\'{a}bek'
 'A. R. Cabral'
 'J. Haloda'
 'T. Sidorinov\\'{a}'
 'B. Lehmann'
 'H. F. Galbiatti'
 'J. Drahokoupil'
_journal_name_full_name
;
 Canadian Mineralogist
;
_journal_volume 50
_journal_year 2012
_journal_page_first 431
_journal_page_last 440
_publ_Section_title
;
 Jacutingaite, Pt$_{2}$HgSe$_{3}$, A New Platinum-Group Mineral Species
      ↪ From the Cau\^{e} Iron-Ore Deposit, Itabira District, Minas
      ↪ Gerais, Brazil
;

# Found in A. Marrazzo, N. Marzari, and M. Gibertini, Emergent dual
      ↪ topology in the three-dimensional Kane-Mele Pt$_{2}$HgSe$_{3}$,
      ↪ Phys. Rev. Research 2, 012063 (2020)

_aflow_title 'Jacutingaite (Pt$_{2}$HgSe$_{3}$) Structure'
_aflow_proto 'AB2C3_hP12_164_d_ae_i'
_aflow_params 'a,c/a,z_{2},x_{4},z_{4}'
_aflow_params_values '7.3477,0.720701716183,0.3507,0.8196,0.2492'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hP12'

_symmetry_space_group_name_H-M "P -3 2/m 1"
_symmetry_Int_Tables_number 164

_cell_length_a     7.34770
_cell_length_b     7.34770
_cell_length_c     5.29550
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 x-y,-y,-z
5 y,x,-z
6 -x,-x+y,-z
7 -x,-y,-z
8 y,-x+y,-z
9 x-y,x,-z
10 -x+y,y,z
11 -y,-x,z
12 x,x-y,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Pt1 Pt 1 a 0.00000 0.00000 0.00000 1.00000
Hg1 Hg 2 d 0.33333 0.66667 0.35070 1.00000
Pt2 Pt 3 e 0.50000 0.00000 0.00000 1.00000
Se1 Se 6 i 0.81960 0.18040 0.24920 1.00000
```

Jacutingaite (Pt_2HgSe_3): AB2C3_hP12_164_d_ae_i - POSCAR

```
AB2C3_hP12_164_d_ae_i & a,c/a,z2,x4,z4 --params=7.3477,0.720701716183,
      ↪ 0.3507,0.8196,0.2492 & P-3m1 D_{3d}^{3} #164 (adei) & hP12 &
      ↪ None & HgPt2Se3 & Jacutingaite & A. Vymazalov\'{a} et al., Can.
      ↪ Mineral. 50, 431-440 (2012)
   1.00000000000000
   3.67385000000000   -6.36329485938692    0.00000000000000
   3.67385000000000    6.36329485938692    0.00000000000000
   0.00000000000000    0.00000000000000    5.29550000000000
```

```
    Hg     Pt     Se
     2      4      6
Direct
   0.33333333333333   0.66666666666667   0.35070000000000   Hg   (2d)
   0.66666666666667   0.33333333333333  -0.35070000000000   Hg   (2d)
   0.00000000000000   0.00000000000000   0.00000000000000   Pt   (1a)
   0.50000000000000   0.00000000000000   0.00000000000000   Pt   (3e)
   0.00000000000000   0.50000000000000   0.00000000000000   Pt   (3e)
   0.50000000000000   0.50000000000000   0.00000000000000   Pt   (3e)
   0.81960000000000  -0.81960000000000   0.24920000000000   Se   (6i)
   0.81960000000000   1.63920000000000   0.24920000000000   Se   (6i)
  -1.63920000000000  -0.81960000000000   0.24920000000000   Se   (6i)
  -0.81960000000000   0.81960000000000  -0.24920000000000   Se   (6i)
   1.63920000000000   0.81960000000000  -0.24920000000000   Se   (6i)
  -0.81960000000000  -1.63920000000000  -0.24920000000000   Se   (6i)
```

$D0_{13}$ ($AlCl_3$) (*obsolete*): AB3_hP4_164_b_ad - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'AlCl3'
_chemical_formula_sum 'Al Cl3'

loop_
_publ_author_name
 'W. E. Laschkarew'
_journal_name_full_name
;
 Zeitschrift fur Anorganische und Allgemeine Chemie
;
_journal_volume 193
_journal_year 1930
_journal_page_first 270
_journal_page_last 276
_publ_Section_title
;
 Zur Struktur AlCl$_{3}$
;

# Found in Strukturbericht Band II 1928-1932, 1937

_aflow_title '$D0_{13}$ (AlCl$_{3}$) ({\em{obsolete}}) Structure'
_aflow_proto 'AB3_hP4_164_b_ad'
_aflow_params 'a,c/a,z_{3}'
_aflow_params_values '3.475,2.45,0.33333'
_aflow_Strukturbericht '$D0_{13}$'
_aflow_Pearson 'hP4'

_symmetry_space_group_name_H-M "P -3 2/m 1"
_symmetry_Int_Tables_number 164

_cell_length_a     3.47500
_cell_length_b     3.47500
_cell_length_c     8.51375
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 x-y,-y,-z
5 y,x,-z
6 -x,-x+y,-z
7 -x,-y,-z
8 y,-x+y,-z
9 x-y,x,-z
10 -x+y,y,z
11 -y,-x,z
12 x,x-y,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cl1 Cl 1 a 0.00000 0.00000 0.00000 1.00000
Al1 Al 1 b 0.00000 0.00000 0.50000 1.00000
Cl2 Cl 2 d 0.33333 0.66667 0.33333 1.00000
```

$D0_{13}$ ($AlCl_3$) (*obsolete*): AB3_hP4_164_b_ad - POSCAR

```
AB3_hP4_164_b_ad & a,c/a,z3 --params=3.475,2.45,0.33333 & P-3m1 D_{3d}^{
    ↪ 3} #164 (abd) & hP4 & $D0_{13}$ & AlCl3 & AlCl3 & W. E.
    ↪ Laschkarew, Z. Anorg. Allg. Chem. 193, 270-276 (1930)
   1.00000000000000
   1.73750000000000  -3.00943827815092   0.00000000000000
   1.73750000000000   3.00943827815092   0.00000000000000
   0.00000000000000   0.00000000000000   8.51375000000000
    Al    Cl
     1     3
Direct
   0.00000000000000   0.00000000000000   0.50000000000000   Al   (1b)
   0.00000000000000   0.00000000000000   0.00000000000000   Cl   (1a)
   0.33333333333333   0.66666666666667   0.33333000000000   Cl   (2d)
   0.66666666666667   0.33333333333333  -0.33333000000000   Cl   (2d)
```

Nevskite (BiSe): AB_hP12_164_c2d_c2d - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Nevskite'
_chemical_formula_sum 'Bi Se'

loop_
_publ_author_name
 'E. Gaudin'
 'S. Jobic'
 'M. Evain'
 'R. Brec'
 'J. Rouxel'
_journal_name_full_name
;
 Materials Research Bulletin
;
_journal_volume 30
_journal_year 1995
_journal_page_first 549
_journal_page_last 561
_publ_Section_title
;
 Charge balance in some Bi$_{x}$Se$_{y}$ phases through atomic structure
    ↪ determination and band structure calculations
;

_aflow_title 'Nevskite (BiSe) Structure'
_aflow_proto 'AB_hP12_164_c2d_c2d'
_aflow_params 'a,c/a,z_{1},z_{2},z_{3},z_{4},z_{5},z_{6}'
_aflow_params_values '4.212,5.44681861349,0.3727,0.1279,0.0419,0.7959,
    ↪ 0.2797,0.5596'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hP12'

_symmetry_space_group_name_H-M "P -3 2/m 1"
_symmetry_Int_Tables_number 164

_cell_length_a     4.21200
_cell_length_b     4.21200
_cell_length_c     22.94200
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 x-y,-y,-z
5 y,x,-z
6 -x,-x+y,-z
7 -x,-y,-z
8 y,-x+y,-z
9 x-y,x,-z
10 -x+y,y,z
11 -y,-x,z
12 x,x-y,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Bi1 Bi 2 c 0.00000 0.00000 0.37270 1.00000
Se1 Se 2 c 0.00000 0.00000 0.12790 1.00000
Bi2 Bi 2 d 0.33333 0.66667 0.04190 1.00000
Bi3 Bi 2 d 0.33333 0.66667 0.79590 1.00000
Se2 Se 2 d 0.33333 0.66667 0.27970 1.00000
Se3 Se 2 d 0.33333 0.66667 0.55960 1.00000
```

Nevskite (BiSe): AB_hP12_164_c2d_c2d - POSCAR

```
AB_hP12_164_c2d_c2d & a,c/a,z1,z2,z3,z4,z5,z6 --params=4.212,
    ↪ 5.44681861349,0.3727,0.1279,0.0419,0.7959,0.2797,0.5596 & P-3m1
    ↪ D_{3d}^{3} #164 (c^2d^4) & hP12 & None & BiSe & Nevskite & E.
    ↪ Gaudin et al., Mater. Res. Bull. 30, 549-561 (1995)
   1.00000000000000
   2.10600000000000  -3.64769900074005   0.00000000000000
   2.10600000000000   3.64769900074005   0.00000000000000
   0.00000000000000   0.00000000000000  22.94200000000000
    Bi    Se
     6     6
Direct
   0.00000000000000   0.00000000000000   0.37270000000000   Bi   (2c)
   0.00000000000000   0.00000000000000  -0.37270000000000   Bi   (2c)
   0.33333333333333   0.66666666666667   0.04190000000000   Bi   (2d)
   0.66666666666667   0.33333333333333  -0.04190000000000   Bi   (2d)
   0.33333333333333   0.66666666666667   0.79590000000000   Bi   (2d)
   0.66666666666667   0.33333333333333  -0.79590000000000   Bi   (2d)
   0.00000000000000   0.00000000000000   0.12790000000000   Se   (2c)
   0.00000000000000   0.00000000000000  -0.12790000000000   Se   (2c)
   0.33333333333333   0.66666666666667   0.27970000000000   Se   (2d)
   0.66666666666667   0.33333333333333  -0.27970000000000   Se   (2d)
   0.33333333333333   0.66666666666667   0.55960000000000   Se   (2d)
   0.66666666666667   0.33333333333333  -0.55960000000000   Se   (2d)
```

$B_{13}C_2$ "B_4C" ($D1_g$): A13B2_hR15_166_b2h_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'B13C2'
_chemical_formula_sum 'B13 C2'

loop_
_publ_author_name
 'G. Will'
 'K. H. Kossobutzki'
_journal_name_full_name
;
 Journal of the Less-Common Metals
;
_journal_volume 44
_journal_year 1976
_journal_page_first 87
_journal_page_last 97
_publ_Section_title
;
 An X-ray structure analysis of boron carbide, B${13}$C${2}$
;

_aflow_title 'B${13}$C${2}$ ''B${4}$C'' (${D1_{g}$) Structure'
_aflow_proto 'A13B2_hR15_166_b2h_c'
_aflow_params 'a,c/a,x_{2},x_{3},z_{3},x_{4},z_{4}'
_aflow_params_values '6.617,1.82847211727,0.3823,0.8044,1.316,0.9936,
      ↪ 1.6714'
_aflow_Strukturbericht '${D1_{g}$'
_aflow_Pearson 'hR15'

_symmetry_space_group_name_H-M "R -3 2/m (hexagonal axes)"
_symmetry_Int_Tables_number 166

_cell_length_a     6.61700
_cell_length_b     6.61700
_cell_length_c     12.09900
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 y,x,-z
5 -x,-x+y,-z
6 x-y,-y,-z
7 -x,-y,-z
8 y,-x+y,-z
9 x-y,x,-z
10 -y,-x,z
11 x,x-y,z
12 -x+y,y,z
13 x+1/3,y+2/3,z+2/3
14 -y+1/3,x-y+2/3,z+2/3
15 -x+y+1/3,-x+2/3,z+2/3
16 y+1/3,x+2/3,-z+2/3
17 -x+1/3,-x+y+2/3,-z+2/3
18 x-y+1/3,-y+2/3,-z+2/3
19 -x+1/3,-y+2/3,-z+2/3
20 y+1/3,-x+y+2/3,-z+2/3
21 x-y+1/3,x+2/3,-z+2/3
22 -y+1/3,-x+2/3,z+2/3
23 x+1/3,x-y+2/3,z+2/3
24 -x+y+1/3,y+2/3,z+2/3
25 x+2/3,y+1/3,z+1/3
26 -y+2/3,x-y+1/3,z+1/3
27 -x+y+2/3,-x+1/3,z+1/3
28 y+2/3,x+1/3,-z+1/3
29 -x+2/3,-x+y+1/3,-z+1/3
30 x-y+2/3,-y+1/3,-z+1/3
31 -x+2/3,-y+1/3,-z+1/3
32 y+2/3,-x+y+1/3,-z+1/3
33 x-y+2/3,x+1/3,-z+1/3
34 -y+2/3,-x+1/3,z+1/3
35 x+2/3,x-y+1/3,z+1/3
36 -x+y+2/3,y+1/3,z+1/3

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
B1 B 3 b 0.00000 0.00000 0.50000 1.00000
C1 C 6 c 0.00000 0.00000 0.38230 1.00000
B2 B 18 h 0.16280 0.83720 0.64160 1.00000
B3 B 18 h 0.10740 0.89260 0.88620 1.00000
```

$B_{13}C_2$ "B_4C" ($D1_g$): A13B2_hR15_166_b2h_c - POSCAR

```
A13B2_hR15_166_b2h_c & a,c/a,x2,x3,z3,x4,z4 --params=6.617,1.82847211727
      ↪ ,0.3823,0.8044,1.316,0.9936,1.6714 & R-3m D_{3d}^{5} #166 (bch^
      ↪ 2) & hR15 & $D1_{g}$ & B13C2 & B13C2 & G. Will and K. H.
      ↪ Kossobutzki, J. Less-Common Met. 44, 87-97 (1976)
   1.00000000000000
    3.30850000000000  -1.91016336561388   4.03300000000000
    0.00000000000000   3.82032673122775   4.03300000000000
   -3.30850000000000  -1.91016336561388   4.03300000000000
      B     C
     13     2
Direct
    0.50000000000000   0.50000000000000   0.50000000000000   B   (1b)
    0.80440000000000   0.80440000000000   1.31600000000000   B   (6h)
    1.31600000000000   0.80440000000000   0.80440000000000   B   (6h)
    0.80440000000000   1.31600000000000   0.80440000000000   B   (6h)
   -1.31600000000000  -0.80440000000000  -0.80440000000000   B   (6h)
   -0.80440000000000  -0.80440000000000  -1.31600000000000   B   (6h)
   -0.80440000000000  -1.31600000000000  -0.80440000000000   B   (6h)
    0.99360000000000   0.99360000000000   1.67140000000000   B   (6h)
    1.67140000000000   0.99360000000000   0.99360000000000   B   (6h)
    0.99360000000000   1.67140000000000   0.99360000000000   B   (6h)
   -1.67140000000000  -0.99360000000000  -0.99360000000000   B   (6h)
   -0.99360000000000  -0.99360000000000  -1.67140000000000   B   (6h)
   -0.99360000000000  -1.67140000000000  -0.99360000000000   B   (6h)
    0.38230000000000   0.38230000000000   0.38230000000000   C   (2c)
   -0.38230000000000  -0.38230000000000  -0.38230000000000   C   (2c)
```

$MnBi_2Te_4$: A2BC4_hR7_166_c_a_2c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Bi2MnTe4'
_chemical_formula_sum 'Bi2 Mn Te4'

loop_
_publ_author_name
 'J.-Q. Yan'
 'Q. Zhang'
 'T. Heitmann'
 'Z. Huang'
 'K. Y. Chen'
 'J.-G. Cheng'
 'W. Wu'
 'D. Vaknin'
 'B. C. Sales'
 'R. J. McQueeney'
_journal_name_full_name
;
 Physical Review Materials
;
_journal_volume 3
_journal_year 2019
_journal_page_first 064202
_journal_page_last 064202
_publ_Section_title
;
 Crystal growth and magnetic structure of MnBi${2}$Te${4}$
;

_aflow_title 'MnBi${2}$Te${4}$ Structure'
_aflow_proto 'A2BC4_hR7_166_c_a_2c'
_aflow_params 'a,c/a,x_{2},x_{3},x_{4}'
_aflow_params_values '4.309,9.44047342771,0.4247,0.1332,0.294'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hR7'

_symmetry_space_group_name_H-M "R -3 2/m (hexagonal axes)"
_symmetry_Int_Tables_number 166

_cell_length_a     4.30900
_cell_length_b     4.30900
_cell_length_c     40.67900
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 y,x,-z
5 -x,-x+y,-z
6 x-y,-y,-z
7 -x,-y,-z
8 y,-x+y,-z
9 x-y,x,-z
10 -y,-x,z
11 x,x-y,z
12 -x+y,y,z
13 x+1/3,y+2/3,z+2/3
14 -y+1/3,x-y+2/3,z+2/3
15 -x+y+1/3,-x+2/3,z+2/3
16 y+1/3,x+2/3,-z+2/3
17 -x+1/3,-x+y+2/3,-z+2/3
18 x-y+1/3,-y+2/3,-z+2/3
19 -x+1/3,-y+2/3,-z+2/3
20 y+1/3,-x+y+2/3,-z+2/3
21 x-y+1/3,x+2/3,-z+2/3
22 -y+1/3,-x+2/3,z+2/3
23 x+1/3,x-y+2/3,z+2/3
24 -x+y+1/3,y+2/3,z+2/3
25 x+2/3,y+1/3,z+1/3
26 -y+2/3,x-y+1/3,z+1/3
27 -x+y+2/3,-x+1/3,z+1/3
28 y+2/3,x+1/3,-z+1/3
29 -x+2/3,-x+y+1/3,-z+1/3
```

```
30 x-y+2/3,-y+1/3,-z+1/3
31 -x+2/3,-y+1/3,-z+1/3
32 y+2/3,-x+y+1/3,-z+1/3
33 x-y+2/3,x+1/3,-z+1/3
34 -y+2/3,-x+1/3,z+1/3
35 x+2/3,x-y+1/3,z+1/3
36 -x+y+2/3,y+1/3,z+1/3

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Mn1 Mn 3 a 0.00000 0.00000 0.00000 1.00000
Bi1 Bi 6 c 0.00000 0.00000 0.42470 1.00000
Te1 Te 6 c 0.00000 0.00000 0.13320 1.00000
Te2 Te 6 c 0.00000 0.00000 0.29400 1.00000
```

$MnBi_2Te_4$: A2BC4_hR7_166_c_a_2c - POSCAR

```
A2BC4_hR7_166_c_a_2c & a,c/a,x2,x3,x4 --params=4.309,9.44047342771,
    ↪ 0.4247,0.1332,0.294 & R-3m D_{3d}^{5} #166 (ac^3) & hR7 & None
    ↪ & Bi2MnTe4 & Bi2MnTe4 & J.-Q. Yan et al., Phys. Rev. Mater. 3,
    ↪ 064202(2019)
   1.00000000000000
   2.15450000000000  -1.24390115496905  13.55966666666670
   0.00000000000000   2.48780230993810  13.55966666666670
  -2.15450000000000  -1.24390115496905  13.55966666666670
     Bi     Mn     Te
      2      1      4
Direct
   0.42470000000000   0.42470000000000   0.42470000000000   Bi   (2c)
  -0.42470000000000  -0.42470000000000  -0.42470000000000   Bi   (2c)
   0.00000000000000   0.00000000000000   0.00000000000000   Mn   (1a)
   0.13320000000000   0.13320000000000   0.13320000000000   Te   (2c)
  -0.13320000000000  -0.13320000000000  -0.13320000000000   Te   (2c)
   0.29400000000000   0.29400000000000   0.29400000000000   Te   (2c)
  -0.29400000000000  -0.29400000000000  -0.29400000000000   Te   (2c)
```

Shandite ($Ni_3Pb_2S_2$): A3B2C2_hR7_166_d_ab_c - CIF

```
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_audit_creation_method FINDSYM

_chemical_name_mineral 'Shandite'
_chemical_formula_sum 'Ni3 Pb2 S2'

loop_
_publ_author_name
 'R. Weihrich'
 'S. F. Matar'
 'V. Eyert'
 'F. Rau'
 'M. Zabel'
 'M. Andratschke'
 'I. Anusca'
 'T. Bernert'
_journal_name_full_name
;
 Progress in Solid State Chemistry
;
_journal_volume 35
_journal_year 2007
_journal_page_first 309
_journal_page_last 322
_publ_Section_title
;
 Structure, ordering, and bonding of half antiperovskites: PbNi$_{3/2}$S
    ↪ and BiPd$_{3/2}$S
;

_aflow_title 'Shandite (Ni$_{3}$Pb$_{2}$S$_{2}$) Structure'
_aflow_proto 'A3B2C2_hR7_166_d_ab_c'
_aflow_params 'a,c/a,x_{3}'
_aflow_params_values '5.595,2.43538873995,0.275'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hR7'

_symmetry_space_group_name_H-M "R -3 2/m (hexagonal axes)"
_symmetry_Int_Tables_number 166

_cell_length_a     5.59500
_cell_length_b     5.59500
_cell_length_c     13.62600
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
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2 -y,x-y,z
3 -x+y,-x,z
4 y,x,-z
5 -x,-x+y,-z
6 x-y,-y,-z
7 -x,-y,-z
8 y,-x+y,-z
9 x-y,x,-z
10 -y,-x,z
11 x,x-y,z
12 -x+y,y,z
13 x+1/3,y+2/3,z+2/3
14 -y+1/3,x-y+2/3,z+2/3
15 -x+y+1/3,-x+2/3,z+2/3
16 y+1/3,x+2/3,-z+2/3
17 -x+1/3,-x+y+2/3,-z+2/3
18 x-y+1/3,-y+2/3,-z+2/3
19 -x+1/3,-y+2/3,-z+2/3
20 y+1/3,-x+y+2/3,-z+2/3
21 x-y+1/3,x+2/3,-z+2/3
22 -y+1/3,-x+2/3,z+2/3
23 x+1/3,x-y+2/3,z+2/3
24 -x+y+1/3,y+2/3,z+2/3
25 x+2/3,y+1/3,z+1/3
26 -y+2/3,x-y+1/3,z+1/3
27 -x+y+2/3,-x+1/3,z+1/3
28 y+2/3,x+1/3,-z+1/3
29 -x+2/3,-x+y+1/3,-z+1/3
30 x-y+2/3,-y+1/3,-z+1/3
31 -x+2/3,-y+1/3,-z+1/3
32 y+2/3,-x+y+1/3,-z+1/3
33 x-y+2/3,x+1/3,-z+1/3
34 -y+2/3,-x+1/3,z+1/3
35 x+2/3,x-y+1/3,z+1/3
36 -x+y+2/3,y+1/3,z+1/3

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Pb1 Pb 3 a 0.00000 0.00000 0.00000 1.00000
Pb2 Pb 3 b 0.00000 0.00000 0.50000 1.00000
S1 S 6 c 0.00000 0.00000 0.27500 1.00000
Ni1 Ni 9 d 0.50000 0.00000 0.50000 1.00000
```

Shandite ($Ni_3Pb_2S_2$): A3B2C2_hR7_166_d_ab_c - POSCAR

```
A3B2C2_hR7_166_d_ab_c & a,c/a,x3 --params=5.595,2.43538873995,0.275 &
    ↪ R-3m D_{3d}^{5} #166 (abcd) & hR7 & None & Ni3Pb2S2 & Shandite
    ↪ & R. Weihrich et al., Prog. Solid State Chem. 35, 309-322 (2007
    ↪ )
   1.00000000000000
   2.79750000000000  -1.61513737805798   4.54200000000000
   0.00000000000000   3.23027475611596   4.54200000000000
  -2.79750000000000  -1.61513737805798   4.54200000000000
     Ni     Pb     S
      3      2      2
Direct
   0.50000000000000   0.00000000000000   0.00000000000000   Ni   (3d)
   0.00000000000000   0.50000000000000   0.00000000000000   Ni   (3d)
   0.00000000000000   0.00000000000000   0.50000000000000   Ni   (3d)
   0.00000000000000   0.00000000000000   0.00000000000000   Pb   (1a)
   0.50000000000000   0.50000000000000   0.50000000000000   Pb   (1b)
   0.27500000000000   0.27500000000000   0.27500000000000   S    (2c)
  -0.27500000000000  -0.27500000000000  -0.27500000000000   S    (2c)
```

Chabazite ($Ca_{1.4}Sr_{0.3}Al_{3.8}Si_{8.3}O_{24}{\cdot}13H_2O$, $S3_4$ (I)): A5B21C24D12_hR62_166_a2c_ehi_fg2h_i - CIF

```
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_audit_creation_method FINDSYM

_chemical_name_mineral 'Chabazite'
_chemical_formula_sum 'Ca5 (H2O)21 O24 Si12'

loop_
_publ_author_name
 'M. Calligaris'
 'G. Nardin'
 'L. Randaccio'
 'P. C. Chiaramonti'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 38
_journal_year 1982
_journal_page_first 602
_journal_page_last 605
_publ_Section_title
;
 Cation-site location in a natural chabazite
;

_aflow_title 'Chabazite (Ca$_{1.4}$Sr$_{0.3}$Al$_{3.8}$Si$_{8.3}$O$_{24}
    ↪ $\cdot$13H$_{2}$O, $S3_{4}$ (I)) Structure'
_aflow_proto 'A5B21C24D12_hR62_166_a2c_ehi_fg2h_i'
_aflow_params 'a,c/a,x_{2},x_{3},x_{5},x_{6},x_{7},z_{7},x_{8},z_{8},x_{
    ↪ 9},z_{9},x_{10},y_{10},z_{10},x_{11},y_{11},z_{11}'
_aflow_params_values '13.80257,1.09220819021,0.2038,0.4065,0.2638,1.1548
    ↪ ,0.418,0.75989,0.2515,0.89459,1.0248,-0.67229,0.2024,-0.5087,
    ↪ 0.31011,0.1044,-0.1251,0.33381'
_aflow_Strukturbericht '$S3_{4}$ (I)'
_aflow_Pearson 'hR62'

_symmetry_space_group_name_H-M "R -3 2/m (hexagonal axes)"
_symmetry_Int_Tables_number 166
```

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_cell_length_a      13.80257
_cell_length_b      13.80257
_cell_length_c      15.07528
_cell_angle_alpha   90.00000
_cell_angle_beta    90.00000
_cell_angle_gamma   120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 y,x,-z
5 -x,-x+y,-z
6 x-y,-y,-z
7 -x,-y,-z
8 y,-x+y,-z
9 x-y,x,-z
10 -y,-x,z
11 x,x-y,z
12 -x+y,y,z
13 x+1/3,y+2/3,z+2/3
14 -y+1/3,x-y+2/3,z+2/3
15 -x+y+1/3,-x+2/3,z+2/3
16 y+1/3,x+2/3,-z+2/3
17 -x+1/3,-x+y+2/3,-z+2/3
18 x-y+1/3,-y+2/3,-z+2/3
19 -x+1/3,-y+2/3,-z+2/3
20 y+1/3,-x+y+2/3,-z+2/3
21 x-y+1/3,x+2/3,-z+2/3
22 -y+1/3,-x+2/3,z+2/3
23 x+1/3,x-y+2/3,z+2/3
24 -x+y+1/3,y+2/3,z+2/3
25 x+2/3,y+1/3,z+1/3
26 -y+2/3,x-y+1/3,z+1/3
27 -x+y+2/3,-x+1/3,z+1/3
28 y+2/3,x+1/3,-z+1/3
29 -x+2/3,-x+y+1/3,-z+1/3
30 x-y+2/3,-y+1/3,-z+1/3
31 -x+2/3,-y+1/3,-z+1/3
32 y+2/3,-x+y+1/3,-z+1/3
33 x-y+2/3,x+1/3,-z+1/3
34 -y+2/3,-x+1/3,z+1/3
35 x+2/3,x-y+1/3,z+1/3
36 -x+y+2/3,y+1/3,z+1/3

loop_
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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ca1 Ca 3 a 0.00000 0.00000 0.00000 0.11000
Ca2 Ca 6 c 0.00000 0.00000 0.20380 0.53000
Ca3 Ca 6 c 0.00000 0.00000 0.40650 0.24000
H2O1 H2O 9 e 0.50000 0.00000 0.00000 0.50000
O1 O 18 f 0.26380 0.00000 0.00000 1.00000
O2 O 18 g 0.65480 0.00000 0.50000 1.00000
H2O2 H2O 18 h 0.21937 0.78063 0.19863 0.57000
O3 O 18 h 0.11897 0.88103 0.13253 1.00000
O4 O 18 h 0.89903 0.10097 0.12577 1.00000
H2O3 H2O 36 i 0.20113 0.50997 0.00127 0.23000
Si1 Si 36 i 0.00003 0.22947 0.10437 1.00000
```

Chabazite ($Ca_{1.4}Sr_{0.3}Al_{3.8}Si_{8.3}O_{24}{\cdot}13H_2O$, $S3_4$ (I)): A5B21C24D12_hR62_166_a2c_ehi_fg2h_i - POSCAR

```
A5B21C24D12_hR62_166_a2c_ehi_fg2h_i & a,c/a,x2,x3,x5,x6,x7,z7,x8,z8,x9,
    ↪ z9,x10,y10,z10,x11,y11,z11 --params=13.80257,1.09220819021,
    ↪ 0.2038,0.4065,0.2638,1.1548,0.418,0.75989,0.2515,0.89459,1.0248
    ↪ ,-0.67229,0.2024,-0.5087,0.31011,0.1044,-0.1251,0.33381 & R-3m
    ↪ D_{3d}^{5} #166 (ac^2efgh^3i^2) & hR62 & $S3_{4}$ (I) &
    ↪ Al3.8Ca1.4H26O37Si8.3Sr0.3 & Chabazite & M. Calligaris et al.,
    ↪ Acta Crystallogr. Sect. B Struct. Sci. 38, 602-605 (1982)
   1.00000000000000
   6.90128500000000  -3.98445875250433   5.02509333333333
   0.00000000000000   7.96891750500865   5.02509333333333
  -6.90128500000000  -3.98445875250433   5.02509333333333
     Ca    H2O     O    Si
      5     21     24    12
Direct
   0.00000000000000   0.00000000000000   0.00000000000000   Ca   (1a)
   0.20380000000000   0.20380000000000   0.20380000000000   Ca   (2c)
  -0.20380000000000  -0.20380000000000  -0.20380000000000   Ca   (2c)
   0.40650000000000   0.40650000000000   0.40650000000000   Ca   (2c)
  -0.40650000000000  -0.40650000000000  -0.40650000000000   Ca   (2c)
   0.00000000000000   0.50000000000000   0.50000000000000   H2O  (3e)
   0.50000000000000   0.00000000000000   0.50000000000000   H2O  (3e)
   0.50000000000000   0.50000000000000   0.00000000000000   H2O  (3e)
   0.41800000000000   0.41800000000000   0.75989000000000   H2O  (6h)
   0.75989000000000   0.41800000000000   0.41800000000000   H2O  (6h)
   0.41800000000000   0.75989000000000   0.41800000000000   H2O  (6h)
  -0.75989000000000  -0.41800000000000  -0.41800000000000   H2O  (6h)
  -0.41800000000000  -0.41800000000000  -0.75989000000000   H2O  (6h)
  -0.41800000000000  -0.75989000000000  -0.41800000000000   H2O  (6h)
   0.20240000000000  -0.50870000000000   0.31011000000000   H2O  (12i)
   0.31011000000000   0.20240000000000  -0.50870000000000   H2O  (12i)
  -0.50870000000000   0.31011000000000   0.20240000000000   H2O  (12i)
  -0.31011000000000   0.50870000000000  -0.20240000000000   H2O  (12i)
   0.50870000000000  -0.20240000000000  -0.31011000000000   H2O  (12i)
  -0.20240000000000  -0.31011000000000   0.50870000000000   H2O  (12i)
  -0.20240000000000   0.50870000000000  -0.31011000000000   H2O  (12i)
  -0.31011000000000  -0.20240000000000   0.50870000000000   H2O  (12i)
   0.50870000000000  -0.31011000000000  -0.20240000000000   H2O  (12i)
   0.31011000000000  -0.50870000000000   0.20240000000000   H2O  (12i)
  -0.50870000000000   0.20240000000000   0.31011000000000   H2O  (12i)
   0.20240000000000   0.31011000000000  -0.50870000000000   H2O  (12i)
   0.26380000000000  -0.26380000000000   0.00000000000000   O    (6f)
   0.00000000000000   0.26380000000000  -0.26380000000000   O    (6f)
  -0.26380000000000   0.00000000000000   0.26380000000000   O    (6f)
  -0.26380000000000   0.26380000000000   0.00000000000000   O    (6f)
   0.00000000000000  -0.26380000000000   0.26380000000000   O    (6f)
   0.26380000000000   0.00000000000000  -0.26380000000000   O    (6f)
   1.15480000000000  -1.15480000000000   0.50000000000000   O    (6g)
   0.50000000000000   1.15480000000000  -1.15480000000000   O    (6g)
  -1.15480000000000   0.50000000000000   1.15480000000000   O    (6g)
  -1.15480000000000   1.15480000000000   0.50000000000000   O    (6g)
   0.50000000000000  -1.15480000000000   1.15480000000000   O    (6g)
   1.15480000000000   0.50000000000000  -1.15480000000000   O    (6g)
   0.25150000000000   0.25150000000000   0.89459000000000   O    (6h)
   0.89459000000000   0.25150000000000   0.25150000000000   O    (6h)
   0.25150000000000   0.89459000000000   0.25150000000000   O    (6h)
  -0.89459000000000  -0.25150000000000  -0.25150000000000   O    (6h)
  -0.25150000000000  -0.25150000000000  -0.89459000000000   O    (6h)
  -0.25150000000000  -0.89459000000000  -0.25150000000000   O    (6h)
   1.02480000000000   1.02480000000000  -0.67229000000000   O    (6h)
  -0.67229000000000   1.02480000000000   1.02480000000000   O    (6h)
   1.02480000000000  -0.67229000000000   1.02480000000000   O    (6h)
   0.67229000000000  -1.02480000000000  -1.02480000000000   O    (6h)
  -1.02480000000000  -1.02480000000000   0.67229000000000   O    (6h)
  -1.02480000000000   0.67229000000000  -1.02480000000000   O    (6h)
   0.10440000000000  -0.12510000000000   0.33381000000000   Si   (12i)
   0.33381000000000   0.10440000000000  -0.12510000000000   Si   (12i)
  -0.12510000000000   0.33381000000000   0.10440000000000   Si   (12i)
  -0.33381000000000   0.12510000000000  -0.10440000000000   Si   (12i)
   0.12510000000000  -0.10440000000000  -0.33381000000000   Si   (12i)
  -0.10440000000000  -0.33381000000000   0.12510000000000   Si   (12i)
  -0.10440000000000   0.12510000000000  -0.33381000000000   Si   (12i)
  -0.33381000000000  -0.10440000000000   0.12510000000000   Si   (12i)
   0.12510000000000  -0.33381000000000  -0.10440000000000   Si   (12i)
   0.33381000000000  -0.12510000000000   0.10440000000000   Si   (12i)
  -0.12510000000000   0.10440000000000   0.33381000000000   Si   (12i)
   0.10440000000000   0.33381000000000  -0.12510000000000   Si   (12i)
```

$CaSi_2$ (*C12*): AB2_hR6_166_c_2c - CIF

```
# CIF file
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_audit_creation_method FINDSYM

_chemical_name_mineral 'CaSi2'
_chemical_formula_sum 'Ca Si2'

loop_
_publ_author_name
 'S. M. Castillo'
 'Z. Tang'
 'A. P. Litvinchuk'
 'A. M. Guloy'
_journal_name_full_name
;
 Inorganic Chemistry
;
_journal_volume 55
_journal_year 2016
_journal_page_first 10203
_journal_page_last 10207
_publ_Section_title
;
 Lattice Dynamics of the Rhombohedral Polymorphs of CaSi$_{2}$
;

_aflow_title 'CaSi$_{2}$ ($C12$) Structure'
_aflow_proto 'AB2_hR6_166_c_2c'
_aflow_params 'a,c/a,x_{1},x_{2},x_{3}'
_aflow_params_values '3.8548,7.95397945419,0.083,0.183,0.35'
_aflow_Strukturbericht '$C12$'
_aflow_Pearson 'hR6'

_symmetry_space_group_name_H-M "R -3 2/m (hexagonal axes)"
_symmetry_Int_Tables_number 166

_cell_length_a      3.85480
_cell_length_b      3.85480
_cell_length_c      30.66100
_cell_angle_alpha   90.00000
_cell_angle_beta    90.00000
_cell_angle_gamma   120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 y,x,-z
5 -x,-x+y,-z
6 x-y,-y,-z
7 -x,-y,-z
8 y,-x+y,-z
9 x-y,x,-z
10 -y,-x,z
11 x,x-y,z
12 -x+y,y,z
13 x+1/3,y+2/3,z+2/3
14 -y+1/3,x-y+2/3,z+2/3
```

```
15 -x+y+1/3,-x+2/3,z+2/3
16 y+1/3,x+2/3,-z+2/3
17 -x+1/3,-x+y+2/3,-z+2/3
18 x-y+1/3,-y+2/3,-z+2/3
19 -x+1/3,-y+2/3,-z+2/3
20 y+1/3,-x+y+2/3,-z+2/3
21 x-y+1/3,x+2/3,-z+2/3
22 -y+1/3,-x+2/3,z+2/3
23 x+1/3,x-y+2/3,z+2/3
24 -x+y+1/3,y+2/3,z+2/3
25 x+2/3,y+1/3,z+1/3
26 -y+2/3,x-y+1/3,z+1/3
27 -x+y+2/3,-x+1/3,z+1/3
28 y+2/3,x+1/3,-z+1/3
29 -x+2/3,-x+y+1/3,-z+1/3
30 x-y+2/3,-y+1/3,-z+1/3
31 -x+2/3,-y+1/3,-z+1/3
32 y+2/3,-x+y+1/3,-z+1/3
33 x-y+2/3,x+1/3,-z+1/3
34 -y+2/3,-x+1/3,z+1/3
35 x+2/3,x-y+1/3,z+1/3
36 -x+y+2/3,y+1/3,z+1/3

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ca1 Ca 6 c 0.00000 0.00000 0.08300 1.00000
Si1 Si 6 c 0.00000 0.00000 0.18300 1.00000
Si2 Si 6 c 0.00000 0.00000 0.35000 1.00000
```

$CaSi_2$ (*C12*): AB2_hR6_166_c_2c - POSCAR

```
AB2_hR6_166_c_2c & a,c/a,x1,x2,x3 --params=3.8548,7.95397945419,0.083,
    ↪ 0.183,0.35 & R-3m D_{3d}^{5} #166 (c^3) & hR6 & $C12$ & CaSi2 &
    ↪  CaSi2 & S. M. Castillo et al., Inorg. Chem. 55, 10203-10207 (
    ↪ 2016)
   1.00000000000000
   1.92740000000000  -1.11278490883608  10.22033333333330
   0.00000000000000   2.22556981767217  10.22033333333330
  -1.92740000000000  -1.11278490883608  10.22033333333330
     Ca      Si
      2       4
Direct
   0.08300000000000   0.08300000000000   0.08300000000000   Ca   (2c)
  -0.08300000000000  -0.08300000000000  -0.08300000000000   Ca   (2c)
   0.18300000000000   0.18300000000000   0.18300000000000   Si   (2c)
  -0.18300000000000  -0.18300000000000  -0.18300000000000   Si   (2c)
   0.35000000000000   0.35000000000000   0.35000000000000   Si   (2c)
  -0.35000000000000  -0.35000000000000  -0.35000000000000   Si   (2c)
```

Rhombohedral $CuTi_2S_4$: AB4C2_hR28_166_2c_2c2h_abh - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'CuS4Ti2'
_chemical_formula_sum 'Cu S4 Ti2'

loop_
_publ_author_name
 'N. Soheilnia'
 'K. M. Kleinke'
 'E. Dashjav'
 'H. L. Cuthbert'
 'J. E. Greedan'
 'H. Kleinke'
_journal_name_full_name
;
 Inorganic Chemistry
;
_journal_volume 43
_journal_year 2004
_journal_page_first 6473
_journal_page_last 6478
_publ_Section_title
;
 Crystal Structure and Physical Properties of a New CuTi$_{2}$S$_{4}$
      ↪ Modification in Comparison to the Thiospinel
;

_aflow_title 'Rhombohedral CuTi$_{2}$S$_{4}$ Structure'
_aflow_proto 'AB4C2_hR28_166_2c_2c2h_abh'
_aflow_params 'a,c/a,x_{3},x_{4},x_{5},x_{6},x_{7},z_{7},x_{8},z_{8},x_{
    ↪ 9},z_{9}'
_aflow_params_values '7.0242,4.95914125452,0.18759,0.35449,0.1232,
    ↪ 0.28892,0.19936,0.7283,0.61457,0.13398,1.08544,-0.42793'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hR28'

_symmetry_space_group_name_H-M "R -3 2/m (hexagonal axes)"
_symmetry_Int_Tables_number 166

_cell_length_a    7.02420
_cell_length_b    7.02420
_cell_length_c    34.83400
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 y,x,-z
5 -x,-x+y,-z
6 x-y,-y,-z
7 -x,-y,-z
8 y,-x+y,-z
9 x-y,x,-z
10 -y,-x,z
11 x,x-y,z
12 -x+y,y,z
13 x+1/3,y+2/3,z+2/3
14 -y+1/3,x-y+2/3,z+2/3
15 -x+y+1/3,-x+2/3,z+2/3
16 y+1/3,x+2/3,-z+2/3
17 -x+1/3,-x+y+2/3,-z+2/3
18 x-y+1/3,-y+2/3,-z+2/3
19 -x+1/3,-y+2/3,-z+2/3
20 y+1/3,-x+y+2/3,-z+2/3
21 x-y+1/3,x+2/3,-z+2/3
22 -y+1/3,-x+2/3,z+2/3
23 x+1/3,x-y+2/3,z+2/3
24 -x+y+1/3,y+2/3,z+2/3
25 x+2/3,y+1/3,z+1/3
26 -y+2/3,x-y+1/3,z+1/3
27 -x+y+2/3,-x+1/3,z+1/3
28 y+2/3,x+1/3,-z+1/3
29 -x+2/3,-x+y+1/3,-z+1/3
30 x-y+2/3,-y+1/3,-z+1/3
31 -x+2/3,-y+1/3,-z+1/3
32 y+2/3,-x+y+1/3,-z+1/3
33 x-y+2/3,x+1/3,-z+1/3
34 -y+2/3,-x+1/3,z+1/3
35 x+2/3,x-y+1/3,z+1/3
36 -x+y+2/3,y+1/3,z+1/3

loop_
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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ti1 Ti 3 a 0.00000 0.00000 0.00000 1.00000
Ti2 Ti 3 b 0.00000 0.00000 0.50000 1.00000
Cu1 Cu 6 c 0.00000 0.00000 0.18759 1.00000
Cu2 Cu 6 c 0.00000 0.00000 0.35449 1.00000
S1 S 6 c 0.00000 0.00000 0.12320 1.00000
S2 S 6 c 0.00000 0.00000 0.28892 1.00000
S3 S 18 h 0.15702 0.84298 0.04234 1.00000
S4 S 18 h 0.49353 0.50647 0.12104 1.00000
Ti3 Ti 18 h 0.83779 0.16221 0.24765 1.00000
```

Rhombohedral $CuTi_2S_4$: AB4C2_hR28_166_2c_2c2h_abh - POSCAR

```
AB4C2_hR28_166_2c_2c2h_abh & a,c/a,x3,x4,x5,x6,x7,z7,x8,z8,x9,z9 --
    ↪ params=7.0242,4.95914125452,0.18759,0.35449,0.1232,0.28892,
    ↪ 0.19936,0.7283,0.61457,0.13398,1.08544,-0.42793 & R-3m D_{3d}^{
    ↪ 5} #166 (abc^4h^3) & hR28 & None & CuS4Ti2 & CuS4Ti2 & N.
    ↪ Soheilnia et al., Inorg. Chem. 43, 6473-6478 (2004)
   1.00000000000000
   3.51210000000000  -2.02771188042089  11.61133333333330
   0.00000000000000   4.05542376084177  11.61133333333330
  -3.51210000000000  -2.02771188042089  11.61133333333330
     Cu      S      Ti
      4      16      8
Direct
   0.18759000000000   0.18759000000000   0.18759000000000   Cu   (2c)
  -0.18759000000000  -0.18759000000000  -0.18759000000000   Cu   (2c)
   0.35449000000000   0.35449000000000   0.35449000000000   Cu   (2c)
  -0.35449000000000  -0.35449000000000  -0.35449000000000   Cu   (2c)
   0.12320000000000   0.12320000000000   0.12320000000000   S    (2c)
  -0.12320000000000  -0.12320000000000  -0.12320000000000   S    (2c)
   0.28892000000000   0.28892000000000   0.28892000000000   S    (2c)
  -0.28892000000000  -0.28892000000000  -0.28892000000000   S    (2c)
   0.19936000000000   0.19936000000000   0.72830000000000   S    (6h)
   0.72830000000000   0.19936000000000   0.19936000000000   S    (6h)
   0.19936000000000   0.72830000000000   0.19936000000000   S    (6h)
  -0.72830000000000  -0.19936000000000  -0.19936000000000   S    (6h)
  -0.19936000000000  -0.19936000000000  -0.72830000000000   S    (6h)
  -0.19936000000000  -0.72830000000000  -0.19936000000000   S    (6h)
   0.61457000000000   0.61457000000000   0.13398000000000   S    (6h)
   0.13398000000000   0.61457000000000   0.61457000000000   S    (6h)
   0.61457000000000   0.13398000000000   0.61457000000000   S    (6h)
  -0.13398000000000  -0.61457000000000  -0.61457000000000   S    (6h)
  -0.61457000000000  -0.61457000000000  -0.13398000000000   S    (6h)
  -0.61457000000000  -0.13398000000000  -0.61457000000000   S    (6h)
   0.00000000000000   0.00000000000000   0.00000000000000   Ti   (1a)
   0.50000000000000   0.50000000000000   0.50000000000000   Ti   (1b)
   1.08544000000000   1.08544000000000  -0.42793000000000   Ti   (6h)
  -0.42793000000000   1.08544000000000   1.08544000000000   Ti   (6h)
   1.08544000000000  -0.42793000000000   1.08544000000000   Ti   (6h)
   0.42793000000000  -1.08544000000000  -1.08544000000000   Ti   (6h)
  -1.08544000000000  -1.08544000000000   0.42793000000000   Ti   (6h)
  -1.08544000000000   0.42793000000000  -1.08544000000000   Ti   (6h)
```

$CaCu_4P_2$: AB4C2_hR7_166_a_2c_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'CaCu4P2'
_chemical_formula_sum 'Ca Cu4 P2'

loop_
_publ_author_name
 'A. Mewis'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Naturforschung B
;
_journal_volume 35
_journal_year 1980
_journal_page_first 942
_journal_page_last 945
_publ_Section_title
;
 Darstellung und Struktur der Verbindung CaCu$_{4}$P$_{2}$
;

_aflow_title 'CaCu$_{4}$P$_{2}$ Structure'
_aflow_proto 'AB4C2_hR7_166_a_2c_c'
_aflow_params 'a,c/a,x_{2},x_{3},x_{4}'
_aflow_params_values '4.036,5.51833498513,0.1453,0.4311,0.2503'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hR7'

_symmetry_space_group_name_H-M "R -3 2/m (hexagonal axes)"
_symmetry_Int_Tables_number 166

_cell_length_a    4.03600
_cell_length_b    4.03600
_cell_length_c    22.27200
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
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2 -y,x-y,z
3 -x+y,-x,z
4 y,x,-z
5 -x,-x+y,-z
6 x-y,-y,-z
7 -x,-y,-z
8 y,-x+y,-z
9 x-y,x,-z
10 -y,-x,z
11 x,x-y,z
12 -x+y,y,z
13 x+1/3,y+2/3,z+2/3
14 -y+1/3,x-y+2/3,z+2/3
15 -x+y+1/3,-x+2/3,z+2/3
16 y+1/3,x+2/3,-z+2/3
17 -x+1/3,-x+y+2/3,-z+2/3
18 x-y+1/3,-y+2/3,-z+2/3
19 -x+1/3,-y+2/3,-z+2/3
20 y+1/3,-x+y+2/3,-z+2/3
21 x-y+1/3,x+2/3,-z+2/3
22 -y+1/3,-x+2/3,z+2/3
23 x+1/3,x-y+2/3,z+2/3
24 -x+y+1/3,y+2/3,z+2/3
25 x+2/3,y+1/3,z+1/3
26 -y+2/3,x-y+1/3,z+1/3
27 -x+y+2/3,-x+1/3,z+1/3
28 y+2/3,x+1/3,-z+1/3
29 -x+2/3,-x+y+1/3,-z+1/3
30 x-y+2/3,-y+1/3,-z+1/3
31 -x+2/3,-y+1/3,-z+1/3
32 y+2/3,-x+y+1/3,-z+1/3
33 x-y+2/3,x+1/3,-z+1/3
34 -y+2/3,-x+1/3,z+1/3
35 x+2/3,x-y+1/3,z+1/3
36 -x+y+2/3,y+1/3,z+1/3

loop_
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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ca1 Ca 3 a 0.00000 0.00000 0.00000 1.00000
Cu1 Cu 6 c 0.00000 0.00000 0.14530 1.00000
Cu2 Cu 6 c 0.00000 0.00000 0.43110 1.00000
P1 P 6 c 0.00000 0.00000 0.25030 1.00000
```

$CaCu_4P_2$: AB4C2_hR7_166_a_2c_c - POSCAR

```
AB4C2_hR7_166_a_2c_c & a,c/a,x2,x3,x4 --params=4.036,5.51833498513,
    ↪ 0.1453,0.4311,0.2503 & R-3m D_{3d}^{5} #166 (ac^3) & hR7 & None
    ↪ & CaCu4P2 & CaCu4P2 & A. Mewis, Z. Naturforsch. B 35, 942-945
    ↪ (1980)
   1.00000000000000
   2.01800000000000  -1.16509284322466   7.42400000000000
   0.00000000000000   2.33018568644933   7.42400000000000
  -2.01800000000000  -1.16509284322466   7.42400000000000
     Ca     Cu     P
     1     4     2
Direct
   0.00000000000000   0.00000000000000   0.00000000000000  Ca  (1a)
   0.14530000000000   0.14530000000000   0.14530000000000  Cu  (2c)
  -0.14530000000000  -0.14530000000000  -0.14530000000000  Cu  (2c)
   0.43110000000000   0.43110000000000   0.43110000000000  Cu  (2c)
  -0.43110000000000  -0.43110000000000  -0.43110000000000  Cu  (2c)
   0.25030000000000   0.25030000000000   0.25030000000000  P  (2c)
  -0.25030000000000  -0.25030000000000  -0.25030000000000  P  (2c)
```

$CaUO_4$: AB4C_hR6_166_b_2c_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'CaO4U'
_chemical_formula_sum 'Ca O4 U'

loop_
_publ_author_name
 'B. O. Loopstra'
 'H. M. Rietveld'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 25
_journal_year 1969
_journal_page_first 787
_journal_page_last 791
_publ_Section_title
;
 The structure of some alkaline-earth metal uranates
;

_aflow_title 'CaUO$_{4}$ Structure'
_aflow_proto 'AB4C_hR6_166_b_2c_a'
_aflow_params 'a,c/a,x_{3},x_{4}'
_aflow_params_values '3.87818,4.52899555977,0.1118,0.3627'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hR6'

_symmetry_space_group_name_H-M "R -3 2/m (hexagonal axes)"
_symmetry_Int_Tables_number 166

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_cell_length_b    3.87818
_cell_length_c    17.56426
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_cell_angle_beta  90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 y,x,-z
5 -x,-x+y,-z
6 x-y,-y,-z
7 -x,-y,-z
8 y,-x+y,-z
9 x-y,x,-z
10 -y,-x,z
11 x,x-y,z
12 -x+y,y,z
13 x+1/3,y+2/3,z+2/3
14 -y+1/3,x-y+2/3,z+2/3
15 -x+y+1/3,-x+2/3,z+2/3
16 y+1/3,x+2/3,-z+2/3
17 -x+1/3,-x+y+2/3,-z+2/3
18 x-y+1/3,-y+2/3,-z+2/3
19 -x+1/3,-y+2/3,-z+2/3
20 y+1/3,-x+y+2/3,-z+2/3
21 x-y+1/3,x+2/3,-z+2/3
22 -y+1/3,-x+2/3,z+2/3
23 x+1/3,x-y+2/3,z+2/3
24 -x+y+1/3,y+2/3,z+2/3
25 x+2/3,y+1/3,z+1/3
26 -y+2/3,x-y+1/3,z+1/3
27 -x+y+2/3,-x+1/3,z+1/3
28 y+2/3,x+1/3,-z+1/3
29 -x+2/3,-x+y+1/3,-z+1/3
30 x-y+2/3,-y+1/3,-z+1/3
31 -x+2/3,-y+1/3,-z+1/3
32 y+2/3,-x+y+1/3,-z+1/3
33 x-y+2/3,x+1/3,-z+1/3
34 -y+2/3,-x+1/3,z+1/3
35 x+2/3,x-y+1/3,z+1/3
36 -x+y+2/3,y+1/3,z+1/3

loop_
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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
U1 U 3 a 0.00000 0.00000 0.00000 1.00000
Ca1 Ca 3 b 0.00000 0.00000 0.50000 1.00000
O1 O 6 c 0.00000 0.00000 0.11180 1.00000
```

```
O2 O 6 c 0.00000 0.00000 0.36270 1.00000
```

$CaUO_4$: AB4C_hR6_166_b_2c_a - POSCAR

```
AB4C_hR6_166_b_2c_a & a,c/a,x3,x4 --params=3.87818,4.52899555977,0.1118,
    ↪ 0.3627 & R-3m D_{3d}^{5} #166 (abc^2) & hR6 & None & CaO4U &
    ↪ CaO4U & B. O. Loopstra and H. M. Rietveld, Acta Crystallogr.
    ↪ Sect. B Struct. Sci. 25, 787-791 (1969)
   1.00000000000000
   1.93909000000000  -1.11953413348291   5.85475333333333
   0.00000000000000   2.23906826696582   5.85475333333333
  -1.93909000000000  -1.11953413348291   5.85475333333333
     Ca     O     U
      1     4     1
Direct
   0.50000000000000   0.50000000000000   0.50000000000000   Ca   (1b)
   0.11180000000000   0.11180000000000   0.11180000000000   O    (2c)
  -0.11180000000000  -0.11180000000000  -0.11180000000000   O    (2c)
   0.36270000000000   0.36270000000000   0.36270000000000   O    (2c)
  -0.36270000000000  -0.36270000000000  -0.36270000000000   O    (2c)
   0.00000000000000   0.00000000000000   0.00000000000000   U    (1a)
```

$TaTi_7$ (BCC SQS-16): AB7_hR16_166_c_c2h - CIF

```
# AFLOW.org Repositories
# TaTi/AB7_hR16_166_c_c2h-001.AB params=9.2519726786,1.22474487138,0.875
    ↪ ,0.625,1.625,0.1250000001,1.125,1.6249999999 SG=166 [ANRL doi:
    ↪ 10.1016/j.commatsci.2017.01.017 (part 1), doi: 10.1016/
    ↪ j.commatsci.2018.10.043 (part 2)]
data_TaTi
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_chemical_formula_sum 'Ta Ti7'

loop_
_publ_author_name
 'T. Chakraborty'
 'J. Rogal'
 'R. Drautz'
_journal_name_full_name
;
 Physical Review B
;
_journal_volume 94
_journal_year 2016
_journal_page_first 224104
_journal_page_last 224104
_publ_Section_title
;
 Unraveling the composition dependence of the martensitic transformation
    ↪ temperature: A first-principles study of Ti-Ta alloys
;

_aflow_title 'TaTi$_{7}$ (BCC SQS-16) Structure'
_aflow_proto 'AB7_hR16_166_c_c2h'
_aflow_params 'a,c/a,x_{1},x_{2},x_{3},z_{3},x_{4},z_{4}'
_aflow_params_values '9.2519726786,1.22474487138,0.375,0.125,1.125,
    ↪ 1.6250000002,1.625,0.1249999998'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hR16'

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_cell_length_b    9.2519726786
_cell_length_c    11.3313060883
_cell_angle_alpha    90.0000000000
_cell_angle_beta    90.0000000000
_cell_angle_gamma    120.0000000000
_symmetry_space_group_name_H-M    'R-3m'
_symmetry_Int_Tables_Number    166
loop_
_symmetry_equiv_pos_site_id
_symmetry_equiv_pos_as_xyz
   1   x,y,z
   2   -y,x-y,z
   3   -x+y,-x,z
   4   y,x,-z
   5   x-y,-y,-z
   6   -x,-x+y,-z
   7   -x,-y,-z
   8   y,-x+y,-z
   9   x-y,x,-z
   10   -y,-x,z
   11   -x+y,y,z
   12   x,x-y,z
   13   x+2/3,y+1/3,z+1/3
   14   -y+2/3,x-y+1/3,z+1/3
   15   -x+y+2/3,-x+1/3,z+1/3
   16   y+2/3,x+1/3,-z+1/3
   17   x-y+2/3,-y+1/3,-z+1/3
   18   -x+2/3,-x+y+1/3,-z+1/3
   19   -x+2/3,-y+1/3,-z+1/3
   20   y+2/3,-x+y+1/3,-z+1/3
   21   x-y+2/3,x+1/3,-z+1/3
   22   -y+2/3,-x+1/3,z+1/3
   23   -x+y+2/3,y+1/3,z+1/3
   24   x+2/3,x-y+1/3,z+1/3
   25   x+1/3,y+2/3,z+2/3
   26   -y+1/3,x-y+2/3,z+2/3
   27   -x+y+1/3,-x+2/3,z+2/3
   28   y+1/3,x+2/3,-z+2/3
   29   x-y+1/3,-y+2/3,-z+2/3
   30   -x+1/3,-x+y+2/3,-z+2/3
   31   -x+1/3,-y+2/3,-z+2/3
   32   y+1/3,-x+y+2/3,-z+2/3
   33   x-y+1/3,x+2/3,-z+2/3
   34   -y+1/3,-x+2/3,z+2/3
   35   -x+y+1/3,y+2/3,z+2/3
   36   x+1/3,x-y+2/3,z+2/3
loop_
 _atom_site_label
 _atom_site_occupancy
 _atom_site_fract_x
 _atom_site_fract_y
 _atom_site_fract_z
 _atom_site_thermal_displace_type
 _atom_site_B_iso_or_equiv
 _atom_site_type_symbol
 _atom_site_symmetry_multiplicity
 _atom_site_Wyckoff_label
Ta1  1.0000000000  0.0000000000  -0.0000000000  0.3750000000  Biso  1.0  Ta  6  c
Ti1  1.0000000000  0.0000000000  -0.0000000000  0.1250000000  Biso  1.0  Ti  6  c
Ti2  1.0000000000  0.1666666666  0.8333333334  0.9583333334  Biso  1.0  Ti  18  h
Ti3  1.0000000000  0.8333333334  0.1666666666  0.7916666666  Biso  1.0  Ti  18  h
```

$TaTi_7$ (BCC SQS-16): AB7_hR16_166_c_c2h - POSCAR

```
AB7_hR16_166_c_c2h & a,c/a,x1,x2,x3,z3,x4,z4 --params=9.2519726786,
    ↪ 1.22474487138,0.375,0.125,1.125,1.6250000002,1.625,0.1249999998
    ↪ & R-3m D_{3d}^{5} #166 (c^2h^2) & hR16 & None & TaTi7 & TaTi7
    ↪ & T. Chakraborty and J. Rogal and R. Drautz, Phys. Rev. B 94,
    ↪ 224104(2016)
   1.00000000000000
   4.62598633930000  -2.67081445826239   3.77710202943333
   0.00000000000000   5.34162891652477   3.77710202943333
  -4.62598633930000  -2.67081445826239   3.77710202943333
     Ta    Ti
      2    14
Direct
   0.37500000000000   0.37500000000000   0.37500000000000   Ta   (2c)
  -0.37500000000000  -0.37500000000000  -0.37500000000000   Ta   (2c)
   0.12500000000000   0.12500000000000   0.12500000000000   Ti   (2c)
  -0.12500000000000  -0.12500000000000  -0.12500000000000   Ti   (2c)
   1.12500000000000   1.12500000000000   1.62500000020000   Ti   (6h)
   1.62500000020000   1.12500000000000   1.12500000000000   Ti   (6h)
   1.12500000000000   1.62500000020000   1.12500000000000   Ti   (6h)
  -1.62500000020000  -1.12500000000000  -1.12500000000000   Ti   (6h)
  -1.12500000000000  -1.12500000000000  -1.62500000020000   Ti   (6h)
  -1.12500000000000  -1.62500000020000  -1.12500000000000   Ti   (6h)
   1.62500000000000   1.62500000000000   0.12499999980000   Ti   (6h)
   0.12499999980000   1.62500000000000   1.62500000000000   Ti   (6h)
   1.62500000000000   0.12499999980000   1.62500000000000   Ti   (6h)
  -0.12499999980000  -1.62500000000000  -1.62500000000000   Ti   (6h)
  -1.62500000000000  -1.62500000000000  -0.12499999980000   Ti   (6h)
  -1.62500000000000  -0.12499999980000  -1.62500000000000   Ti   (6h)
```

Rhombohedral Delafossite ($CuFeO_2$): ABC2_hR4_166_a_b_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Delafossite'
_chemical_formula_sum 'Cu Fe O2'

loop_
_publ_author_name
 'C. T. Prewitt'
 'R. D. Shannon'
 'D. B. Rogers'
_journal_name_full_name
;
 Inorganic Chemistry
;
_journal_volume 10
_journal_year 1971
_journal_page_first 719
_journal_page_last 723
_publ_Section_title
;
 Chemistry of noble metal oxides. II. Crystal structures of platinum
    ↪ cobalt dioxide, palladium cobalt dioxide, coppper iron dioxide
    ↪ , and silver iron dioxide
;

# Found in Crystal chemistry and electrical properties of the
    ↪ delafossite structure, 2006

_aflow_title 'Rhombohedral Delafossite (CuFeO$_{2}$) Structure'
_aflow_proto 'ABC2_hR4_166_a_b_c'
_aflow_params 'a,c/a,x_{3}'
_aflow_params_values '3.0351,5.65582682613,0.1066'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hR4'

_symmetry_space_group_name_H-M "R -3 2/m (hexagonal axes)"
_symmetry_Int_Tables_number 166

_cell_length_a     3.03510
_cell_length_b     3.03510
_cell_length_c     17.16600
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
```

```
3 -x+y,-x,z
4 y,x,-z
5 -x,-x+y,-z
6 x-y,-y,-z
7 -x,-y,-z
8 y,-x+y,-z
9 x-y,x,-z
10 -y,-x,z
11 x,x-y,z
12 -x+y,y,z
13 x+1/3,y+2/3,z+2/3
14 -y+1/3,x-y+2/3,z+2/3
15 -x+y+1/3,-x+2/3,z+2/3
16 y+1/3,x+2/3,-z+2/3
17 -x+1/3,-x+y+2/3,-z+2/3
18 x-y+1/3,-y+2/3,-z+2/3
19 -x+1/3,-y+2/3,-z+2/3
20 y+1/3,-x+y+2/3,-z+2/3
21 x-y+1/3,x+2/3,-z+2/3
22 -y+1/3,-x+2/3,z+2/3
23 x+1/3,x-y+2/3,z+2/3
24 -x+y+1/3,y+2/3,z+2/3
25 x+2/3,y+1/3,z+1/3
26 -y+2/3,x-y+1/3,z+1/3
27 -x+y+2/3,-x+1/3,z+1/3
28 y+2/3,x+1/3,-z+1/3
29 -x+2/3,-x+y+1/3,-z+1/3
30 x-y+2/3,-y+1/3,-z+1/3
31 -x+2/3,-y+1/3,-z+1/3
32 y+2/3,-x+y+1/3,-z+1/3
33 x-y+2/3,x+1/3,-z+1/3
34 -y+2/3,-x+1/3,z+1/3
35 x+2/3,x-y+1/3,z+1/3
36 -x+y+2/3,y+1/3,z+1/3

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cu1 Cu 3 a 0.00000 0.00000 0.00000 1.00000
Fe1 Fe 3 b 0.00000 0.00000 0.50000 1.00000
O1 O 6 c 0.00000 0.00000 0.10660 1.00000
```

Rhombohedral Delafossite ($CuFeO_2$): ABC2_hR4_166_a_b_c - POSCAR

```
ABC2_hR4_166_a_b_c & a,c/a,x3 --params=3.0351,5.65582682613,0.1066 &
    ↪ R-3m D_{3d}^{5} #166 (abc) & hR4 & None & CuFeO2 & Delafossite
    ↪ & C. T. Prewitt and R. D. Shannon and D. B. Rogers, Inorg.
    ↪ Chem. 10, 719-723 (1971)
   1.00000000000000
   1.51755000000000  -0.87615790100872   5.72200000000000
   0.00000000000000   1.75231580201743   5.72200000000000
  -1.51755000000000  -0.87615790100872   5.72200000000000
     Cu    Fe    O
      1     1    2
Direct
   0.00000000000000   0.00000000000000   0.00000000000000   Cu   (1a)
   0.50000000000000   0.50000000000000   0.50000000000000   Fe   (1b)
   0.10660000000000   0.10660000000000   0.10660000000000   O    (2c)
  -0.10660000000000  -0.10660000000000  -0.10660000000000   O    (2c)
```

β-Potassium Nitrate (KNO_3): ABC6_hR8_166_a_b_h - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'KNO3'
_chemical_formula_sum 'K N O6'

loop_
_publ_author_name
 'J. K. Nimmo'
 'B. W. Lucas'
_journal_year 1976
_publ_Section_title
;
 The crystal structures of $\gamma$-and $\beta$-KNO$_{3}$ and the $\
    ↪ alpha$-$\beta$-$\gamma$ phase transformations
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title '$\beta$-Potassium Nitrate (KNO$_{3}$) Structure'
_aflow_proto 'ABC6_hR8_166_a_b_h'
_aflow_params 'a,c/a,x_{3},z_{3}'
_aflow_params_values '5.425,1.8130875576,0.605,1.215'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hR8'

_symmetry_space_group_name_H-M "R -3 2/m (hexagonal axes)"
_symmetry_Int_Tables_number 166

_cell_length_a      5.42500
_cell_length_b      5.42500
_cell_length_c      9.83600
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma 120.00000
loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 y,x,-z
5 -x,-x+y,-z
6 x-y,-y,-z
7 -x,-y,-z
8 y,-x+y,-z
9 x-y,x,-z
10 -y,-x,z
11 x,x-y,z
12 -x+y,y,z
13 x+1/3,y+2/3,z+2/3
14 -y+1/3,x-y+2/3,z+2/3
15 -x+y+1/3,-x+2/3,z+2/3
16 y+1/3,x+2/3,-z+2/3
17 -x+1/3,-x+y+2/3,-z+2/3
18 x-y+1/3,-y+2/3,-z+2/3
19 -x+1/3,-y+2/3,-z+2/3
20 y+1/3,-x+y+2/3,-z+2/3
21 x-y+1/3,x+2/3,-z+2/3
22 -y+1/3,-x+2/3,z+2/3
23 x+1/3,x-y+2/3,z+2/3
24 -x+y+1/3,y+2/3,z+2/3
25 x+2/3,y+1/3,z+1/3
26 -y+2/3,x-y+1/3,z+1/3
27 -x+y+2/3,-x+1/3,z+1/3
28 y+2/3,x+1/3,-z+1/3
29 -x+2/3,-x+y+1/3,-z+1/3
30 x-y+2/3,-y+1/3,-z+1/3
31 -x+2/3,-y+1/3,-z+1/3
32 y+2/3,-x+y+1/3,-z+1/3
33 x-y+2/3,x+1/3,-z+1/3
34 -y+2/3,-x+1/3,z+1/3
35 x+2/3,x-y+1/3,z+1/3
36 -x+y+2/3,y+1/3,z+1/3

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
K1 K 3 a 0.00000 0.00000 0.00000 1.00000
N1 N 3 b 0.00000 0.00000 0.50000 1.00000
O1 O 18 h 0.13000 0.87000 0.47500 0.50000
```

β-Potassium Nitrate (KNO_3): ABC6_hR8_166_a_b_h - POSCAR

```
ABC6_hR8_166_a_b_h & a,c/a,x3,z3 --params=5.425,1.8130875576,0.605,1.215
    ↪ & R-3m D_{3d}^{5} #166 (abh) & hR8 & None & KNO3 & KNO3 & J.
    ↪ K. Nimmo and B. W. Lucas, (1976)
   1.00000000000000
   2.71250000000000  -1.56606260517686   3.27866666666667
   0.00000000000000   3.13212521035372   3.27866666666667
  -2.71250000000000  -1.56606260517686   3.27866666666667
     K     N     O
     1     1     6
Direct
   0.00000000000000   0.00000000000000   0.00000000000000   K   (1a)
   0.50000000000000   0.50000000000000   0.50000000000000   N   (1b)
   0.60500000000000   0.60500000000000   1.21500000000000   O   (6h)
   1.21500000000000   0.60500000000000   0.60500000000000   O   (6h)
   0.60500000000000   1.21500000000000   0.60500000000000   O   (6h)
  -1.21500000000000  -0.60500000000000  -0.60500000000000   O   (6h)
  -0.60500000000000  -0.60500000000000  -1.21500000000000   O   (6h)
  -0.60500000000000  -1.21500000000000  -0.60500000000000   O   (6h)
```

K(SH) ($B22$): AB_hR2_166_a_b - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'K(SH)'
_chemical_formula_sum 'K S'

loop_
_publ_author_name
 'C. D. West'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 88
_journal_year 1934
_journal_page_first 97
_journal_page_last 115
_publ_Section_title
;
 The Crystal Structures of Some Alkali Hydrosulfides and Monosulfides
;

# Found in Strukturbericht Band III 1933-1935, 1937

_aflow_title 'K(SH) ($B22$) Structure'
_aflow_proto 'AB_hR2_166_a_b'
_aflow_params 'a,c/a'
_aflow_params_values '9.88622,1.00556532224'
```

```
_aflow_Strukturbericht '$B22$'
_aflow_Pearson 'hR2'

_symmetry_space_group_name_H-M "R -3 2/m (hexagonal axes)"
_symmetry_Int_Tables_number 166

_cell_length_a      9.88622
_cell_length_b      9.88622
_cell_length_c      9.94124
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 y,x,-z
5 -x,-x+y,-z
6 x-y,-y,-z
7 -x,-y,-z
8 y,-x+y,-z
9 x-y,x,-z
10 -y,-x,z
11 x,x-y,z
12 -x+y,y,z
13 x+1/3,y+2/3,z+2/3
14 -y+1/3,x-y+2/3,z+2/3
15 -x+y+1/3,-x+2/3,z+2/3
16 y+1/3,x+2/3,-z+2/3
17 -x+1/3,-x+y+2/3,-z+2/3
18 x-y+1/3,-y+2/3,-z+2/3
19 -x+1/3,-y+2/3,-z+2/3
20 y+1/3,-x+y+2/3,-z+2/3
21 x-y+1/3,x+2/3,-z+2/3
22 -y+1/3,-x+2/3,z+2/3
23 x+1/3,x-y+2/3,z+2/3
24 -x+y+1/3,y+2/3,z+2/3
25 x+2/3,y+1/3,z+1/3
26 -y+2/3,x-y+1/3,z+1/3
27 -x+y+2/3,-x+1/3,z+1/3
28 y+2/3,x+1/3,-z+1/3
29 -x+2/3,-x+y+1/3,-z+1/3
30 x-y+2/3,-y+1/3,-z+1/3
31 -x+2/3,-y+1/3,-z+1/3
32 y+2/3,-x+y+1/3,-z+1/3
33 x-y+2/3,x+1/3,-z+1/3
34 -y+2/3,-x+1/3,z+1/3
35 x+2/3,x-y+1/3,z+1/3
36 -x+y+2/3,y+1/3,z+1/3

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
K1 K 3 a 0.00000 0.00000 0.00000 1.00000
S1 S 3 b 0.00000 0.00000 0.50000 1.00000
```

K(SH) (*B22*): AB_hR2_166_a_b - POSCAR

```
AB_hR2_166_a_b & a,c/a --params=9.88622,1.00556532224 & R-3m D_{3d}^{5}
    ↪ #166 (ab) & hR2 & $B22$ & K(SH) & K(SH) & C. D. West,
    ↪ Zeitschrift f{"u}r Kristallographie - Crystalline Materials 88,
    ↪ 97-115 (1934)
   1.00000000000000
   4.94311000000000  -2.85390588913393   3.31374666666667
   0.00000000000000   5.70781177826786   3.31374666666667
  -4.94311000000000  -2.85390588913393   3.31374666666667
     K     S
     1     1
Direct
   0.00000000000000   0.00000000000000   0.00000000000000   K   (1a)
   0.50000000000000   0.50000000000000   0.50000000000000   S   (1b)
```

$Zr_{21}Re_{25}$: A25B21_hR92_167_b2e3f_e3f - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Re25Zr21'
_chemical_formula_sum 'Re25 Zr21'

loop_
_publ_author_name
 'K. Cenzual'
 'E. Parth\\'{e}'
 'R. M. Waterstrat'
_journal_name_full_name
;
 Acta Crystallographica Section C: Structural Chemistry
;
_journal_volume 42
_journal_year 1986
_journal_page_first 261
_journal_page_last 266
_publ_Section_title
;
 Zr$_{21}$Re$_{25}$, a new rhombohedral structure type containing 12~{\
    ↪ AA}-thick infinite MgZn$_{2}$(Laves)-type columns
;

# Found in The intermetallic compound Mg$_{21}$Zn$_{25}$, 2002

_aflow_title 'Zr$_{21}$Re$_{25}$ Structure'
_aflow_proto 'A25B21_hR92_167_b2e3f_e3f'
_aflow_params 'a,c/a,x_{2},x_{3},x_{4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{
    ↪ 6},x_{7},y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{10
    ↪ },y_{10},z_{10}'
_aflow_params_values '25.847,0.339343057221,0.189,0.8054,0.5557,0.3062,
    ↪ 0.0702,0.3676,0.4819,0.0723,0.1889,0.235,-0.1163,-0.1187,0.0445
    ↪ ,-0.1882,0.2709,0.182,-0.0505,0.0662,0.561,0.1397,0.0436'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hR92'

_symmetry_space_group_name_H-M "R -3 2/c (hexagonal axes)"
_symmetry_Int_Tables_number 167

_cell_length_a      25.84700
_cell_length_b      25.84700
_cell_length_c      8.77100
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 y,x,-z+1/2
5 -x,-x+y,-z+1/2
6 x-y,-y,-z+1/2
7 -x,-y,-z
8 y,-x+y,-z
9 x-y,x,-z
10 -y,-x,z+1/2
11 x,x-y,z+1/2
12 -x+y,y,z+1/2
13 x+1/3,y+2/3,z+2/3
14 -y+1/3,x-y+2/3,z+2/3
15 -x+y+1/3,-x+2/3,z+2/3
16 y+1/3,x+2/3,-z+1/6
17 -x+1/3,-x+y+2/3,-z+1/6
18 x-y+1/3,-y+2/3,-z+1/6
19 -x+1/3,-y+2/3,-z+2/3
20 y+1/3,-x+y+2/3,-z+2/3
21 x-y+1/3,x+2/3,-z+2/3
22 -y+1/3,-x+2/3,z+1/6
23 x+1/3,x-y+2/3,z+1/6
24 -x+y+1/3,y+2/3,z+1/6
25 x+2/3,y+1/3,z+1/3
26 -y+2/3,x-y+1/3,z+1/3
27 -x+y+2/3,-x+1/3,z+1/3
28 y+2/3,x+1/3,-z+5/6
29 -x+2/3,-x+y+1/3,-z+5/6
30 x-y+2/3,-y+1/3,-z+5/6
31 -x+2/3,-y+1/3,-z+1/3
32 y+2/3,-x+y+1/3,-z+1/3
33 x-y+2/3,x+1/3,-z+1/3
34 -y+2/3,-x+1/3,z+5/6
35 x+2/3,x-y+1/3,z+5/6
36 -x+y+2/3,y+1/3,z+5/6

loop_
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_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Re1 Re 6 b 0.00000 0.00000 0.00000 1.00000
Re2 Re 18 e 0.06100 0.00000 0.25000 1.00000
Re3 Re 18 e 0.44460 0.00000 0.25000 1.00000
Zr1 Zr 18 e 0.69430 0.00000 0.25000 1.00000
Re4 Re 36 f 0.05820 0.17780 0.24800 1.00000
Re5 Re 36 f 0.23420 0.17540 0.24770 1.00000
Re6 Re 36 f 0.23500 0.11630 0.00000 1.00000
Zr2 Zr 36 f 0.00210 0.23060 0.04240 1.00000
Zr3 Zr 36 f 0.11610 0.11640 0.06590 1.00000
Zr4 Zr 36 f 0.31290 0.10840 0.24810 1.00000
```

$Zr_{21}Re_{25}$: A25B21_hR92_167_b2e3f_e3f - POSCAR

```
A25B21_hR92_167_b2e3f_e3f & a,c/a,x2,x3,x4,x5,y5,z5,x6,y6,z6,x7,y7,z7,x8
    ↪ ,y8,z8,x9,y9,z9,x10,y10,z10 --params=25.847,0.339343057221,
    ↪ 0.189,0.8054,0.5557,0.3062,0.0702,0.3676,0.4819,0.0723,0.1889,
    ↪ 0.235,-0.1163,-0.1187,0.0445,-0.1882,0.2709,0.182,-0.0505,
    ↪ 0.0662,0.561,0.1397,0.0436 & R-3c D_{3d}^{6} #167 (be^3f^6) &
    ↪ hR92 & None & Re25Zr21 & Re25Zr21 & K. Cenzual and E. Parth\'{e
    ↪ } and R. M. Waterstrat, Acta Crystallogr. C 42, 261-266 (1986)
   1.00000000000000
  12.92350000000000  -7.46138620387213   2.92366666666667
   0.00000000000000  14.92277240774430   2.92366666666667
 -12.92350000000000  -7.46138620387213   2.92366666666667
     Re    Zr
     50    42
Direct
   0.00000000000000   0.00000000000000   0.00000000000000   Re   (2b)
   0.50000000000000   0.50000000000000   0.50000000000000   Re   (2b)
```

```
  0.18900000000000   0.31100000000000   0.25000000000000   Re   (6e)
  0.25000000000000   0.18900000000000   0.31100000000000   Re   (6e)
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  0.75000000000000  -0.18900000000000   0.68900000000000   Re   (6e)
  0.68900000000000   0.75000000000000  -0.18900000000000   Re   (6e)
  0.80540000000000  -0.30540000000000   0.25000000000000   Re   (6e)
  0.25000000000000   0.80540000000000  -0.30540000000000   Re   (6e)
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 -0.80540000000000   0.30540000000000   0.75000000000000   Re   (6e)
  0.75000000000000  -0.80540000000000   0.30540000000000   Re   (6e)
  0.30540000000000   0.75000000000000  -0.80540000000000   Re   (6e)
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  0.42980000000000   0.19380000000000   0.13240000000000   Re   (12f)
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  0.42770000000000   0.01810000000000   0.31110000000000   Re   (12f)
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 -0.07230000000000  -0.18890000000000  -0.48190000000000   Re   (12f)
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  0.57230000000000   0.98190000000000   0.68890000000000   Re   (12f)
  0.98190000000000   0.68890000000000   0.57230000000000   Re   (12f)
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  0.26500000000000   0.61870000000000   0.61630000000000   Re   (12f)
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  0.11870000000000  -0.23500000000000   0.11630000000000   Re   (12f)
  0.11630000000000   0.11870000000000  -0.23500000000000   Re   (12f)
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  0.75000000000000  -0.55570000000000   0.05570000000000   Zr   (6e)
  0.05570000000000   0.75000000000000  -0.55570000000000   Zr   (6e)
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  0.27090000000000   0.04450000000000  -0.18820000000000   Zr   (12f)
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  0.18820000000000  -0.27090000000000  -0.04450000000000   Zr   (12f)
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  0.31180000000000   0.54450000000000   0.77090000000000   Zr   (12f)
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  0.55050000000000   0.31800000000000   0.43380000000000   Zr   (12f)
  0.31800000000000   0.43380000000000   0.55050000000000   Zr   (12f)
 -0.18200000000000   0.05050000000000  -0.06620000000000   Zr   (12f)
 -0.06620000000000  -0.18200000000000   0.05050000000000   Zr   (12f)
  0.05050000000000  -0.06620000000000  -0.18200000000000   Zr   (12f)
  0.56620000000000   0.44950000000000   0.68200000000000   Zr   (12f)
  0.44950000000000   0.68200000000000   0.56620000000000   Zr   (12f)
  0.68200000000000   0.56620000000000   0.44950000000000   Zr   (12f)
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  0.04360000000000   0.56100000000000   0.13970000000000   Zr   (12f)
  0.13970000000000   0.04360000000000   0.56100000000000   Zr   (12f)
  0.45640000000000   0.36030000000000  -0.06100000000000   Zr   (12f)
  0.36030000000000  -0.06100000000000   0.45640000000000   Zr   (12f)
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 -0.56100000000000  -0.13970000000000  -0.04360000000000   Zr   (12f)
 -0.04360000000000  -0.56100000000000  -0.13970000000000   Zr   (12f)
 -0.13970000000000  -0.04360000000000  -0.56100000000000   Zr   (12f)
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  0.63970000000000   1.06100000000000   0.54360000000000   Zr   (12f)
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β-BaB_2O_4 (High-Temperature): A2BC4_hR42_167_f_ac_2f - CIF

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_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 20
_journal_year 1966
_journal_page_first 819
_journal_page_last 823
_publ_Section_title
;
 The crystal structure of the high temperature form of barium borate,
      ↪ BaO$\cdot$B$_{2}$O$_{3}$
;

# Found in Crystal Structure of the low-temperature form of BaB$_{2}$O$_
      ↪ {4}$, 1984

_aflow_title '$\beta-BaB$_{2}$O$_{4}$ (High-Temperature) Structure'
_aflow_proto 'A2BC4_hR42_167_f_ac_2f'
_aflow_params 'a,c/a,x_{2},x_{3},y_{3},z_{3},x_{4},y_{4},z_{4},x_{5},y_{
      ↪ 5},z_{5}'
_aflow_params_values '7.235,5.41700069109,0.84983,0.35996,-0.25104,
      ↪ 0.50497,0.335,-0.429,0.71401,0.4775,0.1905,-0.05849'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hR42'

_symmetry_space_group_name_H-M "R -3 2/c (hexagonal axes)"
_symmetry_Int_Tables_number 167

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_cell_length_c      39.19200
_cell_angle_alpha   90.00000
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_space_group_symop_operation_xyz
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2 -y,x-y,z
3 -x+y,-x,z
4 y,x,-z+1/2
5 -x,-x+y,-z+1/2
6 x-y,-y,-z+1/2
7 -x,-y,-z
8 y,-x+y,-z
9 x-y,x,-z
10 -y,-x,z+1/2
11 x,x-y,z+1/2
12 -x+y,y,z+1/2
13 x+1/3,y+2/3,z+2/3
14 -y+1/3,x-y+2/3,z+2/3
15 -x+y+1/3,-x+2/3,z+2/3
16 y+1/3,x+2/3,-z+1/6
17 -x+1/3,-x+y+2/3,-z+1/6
18 x-y+1/3,-y+2/3,-z+1/6
19 -x+1/3,-y+2/3,-z+2/3
20 y+1/3,-x+y+2/3,-z+2/3
21 x-y+1/3,x+2/3,-z+2/3
22 -y+1/3,-x+2/3,z+1/6
23 x+1/3,x-y+2/3,z+1/6
24 -x+y+1/3,y+2/3,z+1/6
25 x+2/3,y+1/3,z+1/3
26 -y+2/3,x-y+1/3,z+1/3
27 -x+y+2/3,-x+1/3,z+1/3
28 y+2/3,x+1/3,-z+5/6
29 -x+2/3,-x+y+1/3,-z+5/6
30 x-y+2/3,-y+1/3,-z+5/6
31 -x+2/3,-y+1/3,-z+1/3
32 y+2/3,-x+y+1/3,-z+1/3
33 x-y+2/3,x+1/3,-z+1/3
34 -y+2/3,-x+1/3,z+5/6
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_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
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Ba2 Ba 12 c 0.00000 0.00000 0.84983 1.00000
B1 B 36 f 0.15533 0.45567 0.20463 1.00000
O1 O 36 f 0.12833 0.63567 0.20667 1.00000
O2 O 36 f 0.27433 0.01267 0.20317 1.00000
```

β-BaB_2O_4 (High-Temperature): A2BC4_hR42_167_f_ac_2f - POSCAR

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_audit_creation_method FINDSYM

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_chemical_formula_sum 'B2 Ba O4'

loop_
_publ_author_name
 'A. D. Mighell'
 'A. Perloff'
 'S. Block'
```

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A2BC4_hR42_167_f_ac_2f & a,c/a,x2,x3,y3,z3,x4,y4,z4,x5,y5,z5 --params=
      ↪ 7.235,5.41700069109,0.84983,0.35996,-0.25104,0.50497,0.335,-
      ↪ 0.429,0.71401,0.4775,0.1905,-0.05849 & R-3c D_{3d}^{6} #167 (
      ↪ acf^3) & hR42 & None & B2BaO4 & B2BaO4 & A. D. Mighell and A.
      ↪ Perloff and S. Block, Acta Cryst. 20, 819-823 (1966)
   1.00000000000000
   3.61750000000000  -2.08856459879347  13.06400000000000
   0.00000000000000   4.17712919758694  13.06400000000000
  -3.61750000000000  -2.08856459879347  13.06400000000000
     B     Ba     O
    12      6     24
Direct
```

```
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   0.50497000000000   0.35996000000000  −0.25104000000000   B   (12f)
  −0.25104000000000   0.50497000000000   0.35996000000000   B   (12f)
  −0.00497000000000   0.75104000000000   0.14004000000000   B   (12f)
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   0.14004000000000  −0.00497000000000   0.75104000000000   B   (12f)
  −0.35996000000000   0.25104000000000  −0.50497000000000   B   (12f)
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   0.75000000000000   0.75000000000000   0.75000000000000   Ba   (2a)
   0.84983000000000   0.84983000000000   0.84983000000000   Ba   (4c)
  −0.34983000000000  −0.34983000000000  −0.34983000000000   Ba   (4c)
  −0.84983000000000  −0.84983000000000  −0.84983000000000   Ba   (4c)
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  −0.42900000000000   0.71401000000000   0.33500000000000   O   (12f)
  −0.21401000000000   0.92900000000000   0.16500000000000   O   (12f)
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   0.97750000000000   0.44151000000000   0.69050000000000   O   (12f)
```

$CrCl_3(H_2O)_6$ ($J2_2$): A3BC6_hR20_167_e_b_f - CIF

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_audit_creation_method FINDSYM

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_chemical_formula_sum 'Cl3 Cr (H2O)6'

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_publ_author_name
 'K. R. Andress'
 'C. Carpenter'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 87
_journal_year 1934
_journal_page_first 446
_journal_page_last 463
_publ_Section_title
;
 Kristallhydrate. II. Die Struktur von Chromchlorid- und
      ↪ Aluminiumchloridhexahydrat
;

# Found in Strukturbericht Band III 1933-1935, 1937

_aflow_title 'CrCl$_{3}$(H$_{2}$O)$_{6}$ ($J2_{2}$) Structure'
_aflow_proto 'A3BC6_hR20_167_e_b_f'
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_aflow_params_values '11.8335,1.00556555541,0.99,0.22,-0.04,0.12'
_aflow_Strukturbericht '$J2_{2}$'
_aflow_Pearson 'hR20'

_symmetry_space_group_name_H-M "R -3 2/c (hexagonal axes)"
_symmetry_Int_Tables_number 167

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_cell_angle_gamma 120.00000

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_space_group_symop_operation_xyz
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3 -x+y,-x,z
4 y,x,-z+1/2
5 -x,-x+y,-z+1/2
6 x-y,-y,-z+1/2
7 -x,-y,-z
8 y,-x+y,-z
9 x-y,x,-z
10 -y,-x,z+1/2
11 x,x-y,z+1/2
12 -x+y,y,z+1/2
13 x+1/3,y+2/3,z+2/3
14 -y+1/3,x-y+2/3,z+2/3
15 -x+y+1/3,-x+2/3,z+2/3
16 y+1/3,x+2/3,-z+1/6
17 -x+1/3,-x+y+2/3,-z+1/6
18 x-y+1/3,-y+2/3,-z+1/6
19 -x+1/3,-y+2/3,-z+2/3
20 y+1/3,-x+y+2/3,-z+2/3
21 x-y+1/3,x+2/3,-z+2/3
22 -y+1/3,-x+2/3,z+1/6
23 x+1/3,x-y+2/3,z+1/6
24 -x+y+1/3,y+2/3,z+1/6
25 x+2/3,y+1/3,z+1/3
26 -y+2/3,x-y+1/3,z+1/3
27 -x+y+2/3,-x+1/3,z+1/3
28 y+2/3,x+1/3,-z+5/6
29 -x+2/3,-x+y+1/3,-z+5/6
30 x-y+2/3,-y+1/3,-z+5/6
31 -x+2/3,-y+1/3,-z+1/3
32 y+2/3,-x+y+1/3,-z+1/3
33 x-y+2/3,x+1/3,-z+1/3
34 -y+2/3,-x+1/3,z+5/6
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36 -x+y+2/3,y+1/3,z+5/6

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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cr1 Cr 6 b 0.00000 0.00000 0.00000 1.00000
Cl1 Cl 18 e 0.26000 0.00000 0.25000 1.00000
H2O1 H2O 36 f 0.12000 0.14000 0.10000 1.00000
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$CrCl_3(H_2O)_6$ ($J2_2$): A3BC6_hR20_167_e_b_f - POSCAR

```
A3BC6_hR20_167_e_b_f & a,c/a,x2,x3,y3,z3 --params=11.8335,1.00556555541,
      ↪ 0.99,0.22,-0.04,0.12 & R-3c D_{3d}^{6} #167 (bef) & hR20 & $J2_
      ↪ {2}$ & Cl3Cr(H2O)6 & Cl3Cr(H2O)6 & K. R. Andress and C.
      ↪ Carpenter, Zeitschrift f{"u}r Kristallographie - Crystalline
      ↪ Materials 87, 446-463 (1934)
   1.00000000000000
   5.91675000000000  -3.41603720522772   3.96645333333333
   0.00000000000000   6.83207441045544   3.96645333333333
  -5.91675000000000  -3.41603720522772   3.96645333333333
     Cl    Cr   H2O
      6     2    12
Direct
   0.99000000000000  -0.49000000000000   0.25000000000000   Cl   (6e)
   0.25000000000000   0.99000000000000  -0.49000000000000   Cl   (6e)
  -0.49000000000000   0.25000000000000   0.99000000000000   Cl   (6e)
  -0.99000000000000   0.49000000000000   0.75000000000000   Cl   (6e)
   0.75000000000000  -0.99000000000000   0.49000000000000   Cl   (6e)
   0.49000000000000   0.75000000000000  -0.99000000000000   Cl   (6e)
   0.00000000000000   0.00000000000000   0.00000000000000   Cr   (2b)
   0.50000000000000   0.50000000000000   0.50000000000000   Cr   (2b)
   0.22000000000000  -0.04000000000000   0.12000000000000   H2O   (12f)
   0.12000000000000   0.22000000000000  -0.04000000000000   H2O   (12f)
  -0.04000000000000   0.12000000000000   0.22000000000000   H2O   (12f)
   0.38000000000000   0.54000000000000   0.28000000000000   H2O   (12f)
   0.54000000000000   0.28000000000000   0.38000000000000   H2O   (12f)
   0.28000000000000   0.38000000000000   0.54000000000000   H2O   (12f)
  -0.22000000000000   0.04000000000000  -0.12000000000000   H2O   (12f)
  -0.12000000000000  -0.22000000000000   0.04000000000000   H2O   (12f)
   0.04000000000000  -0.12000000000000  -0.22000000000000   H2O   (12f)
   0.62000000000000   0.46000000000000   0.72000000000000   H2O   (12f)
   0.46000000000000   0.72000000000000   0.62000000000000   H2O   (12f)
   0.72000000000000   0.62000000000000   0.46000000000000   H2O   (12f)
```

FeF_3 ($D0_{12}$): A3B_hR8_167_e_b - CIF

```
# CIF file
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_chemical_name_mineral 'F3Fe'
_chemical_formula_sum 'F3 Fe'

loop_
_publ_author_name
 'M. A. Hepworth'
 'K. H. Jack'
 'R. D. Peacock'
 'G. J. Westland'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 10
_journal_year 1957
_journal_page_first 63
_journal_page_last 69
_publ_Section_title
;
 The crystal structures of the trifluorides of iron, cobalt, ruthenium,
      ↪ rhodium, palladium and iridium
;

_aflow_title 'FeF$_{3}$ ($D0_{12}$) Structure'
_aflow_proto 'A3B_hR8_167_e_b'
_aflow_params 'a,c/a,x_{2}'
_aflow_params_values '5.198,2.56464024625,0.664'
```

```
_aflow_Strukturbericht '$D0_{12}$'
_aflow_Pearson 'hR8'

_symmetry_space_group_name_H-M "R -3 2/c (hexagonal axes)"
_symmetry_Int_Tables_number 167

_cell_length_a 5.19800
_cell_length_b 5.19800
_cell_length_c 13.33100
_cell_angle_alpha 90.00000
_cell_angle_beta 90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 y,x,-z+1/2
5 -x,-x+y,-z+1/2
6 x-y,-y,-z+1/2
7 -x,-y,-z
8 y,-x+y,-z
9 x-y,x,-z
10 -y,-x,z+1/2
11 x,x-y,z+1/2
12 -x+y,y,z+1/2
13 x+1/3,y+2/3,z+2/3
14 -y+1/3,x-y+2/3,z+2/3
15 -x+y+1/3,-x+2/3,z+2/3
16 y+1/3,x+2/3,-z+1/6
17 -x+1/3,-x+y+2/3,-z+1/6
18 x-y+1/3,-y+2/3,-z+1/6
19 -x+1/3,-y+2/3,-z+2/3
20 y+1/3,-x+y+2/3,-z+2/3
21 x-y+1/3,x+2/3,-z+2/3
22 -y+1/3,-x+2/3,z+1/6
23 x+1/3,x-y+2/3,z+1/6
24 -x+y+1/3,y+2/3,z+1/6
25 x+2/3,y+1/3,z+1/3
26 -y+2/3,x-y+1/3,z+1/3
27 -x+y+2/3,-x+1/3,z+1/3
28 y+2/3,x+1/3,-z+5/6
29 -x+2/3,-x+y+1/3,-z+5/6
30 x-y+2/3,-y+1/3,-z+5/6
31 -x+2/3,-y+1/3,-z+1/3
32 y+2/3,-x+y+1/3,-z+1/3
33 x-y+2/3,x+1/3,-z+1/3
34 -y+2/3,-x+1/3,z+5/6
35 x+2/3,x-y+1/3,z+5/6
36 -x+y+2/3,y+1/3,z+5/6

loop_
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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Fe1 Fe 6 b 0.00000 0.00000 0.00000 1.00000
F1 F 18 e 0.41400 0.00000 0.25000 1.00000
```

FeF_3 ($D0_{12}$): A3B_hR8_167_e_b - POSCAR

```
A3B_hR8_167_e_b & a,c/a,x2 --params=5.198,2.564604024625,0.664 & R-3c D_{3d}^{6} #167 (be) & hR8 & $D0_{12}$ & F3Fe & F3Fe & M. A. Hepworth et al., Acta Cryst. 10, 63-69 (1957)
   1.00000000000000
   2.59900000000000  -1.50053334962384   4.44366666666667
   0.00000000000000   3.00106669924768   4.44366666666667
  -2.59900000000000  -1.50053334962384   4.44366666666667
     F     Fe
     6      2
Direct
   0.66400000000000  -0.16400000000000   0.25000000000000     F    (6e)
   0.25000000000000   0.66400000000000  -0.16400000000000     F    (6e)
  -0.16400000000000   0.25000000000000   0.66400000000000     F    (6e)
  -0.66400000000000   1.16400000000000   0.75000000000000     F    (6e)
   0.75000000000000  -0.66400000000000   1.16400000000000     F    (6e)
   1.16400000000000   0.75000000000000  -0.66400000000000     F    (6e)
   0.00000000000000   0.00000000000000   0.00000000000000    Fe    (2b)
   0.50000000000000   0.50000000000000   0.50000000000000    Fe    (2b)
```

Rinneite ($K_3NaFeCl_6$): A6BC3D_hR22_167_f_b_e_a - CIF

```
# CIF file
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_audit_creation_method FINDSYM

_chemical_name_mineral 'Rinneite'
_chemical_formula_sum 'Cl6 Fe K3 Na'

loop_
_publ_author_name
 'B. N. Figgis'
 'A. N. Sobolev'
 'E. S. Kucharski'
 'V. Broughton'
_journal_name_full_name
;
 Acta Crystallographica Section C: Structural Chemistry
;
_journal_volume 56
_journal_year 2000
_journal_page_first e228
_journal_page_last e229
_publ_Section_title
;
 Rinneite, K$_{3}$Na[FeCl$_{6}$], at 293, 84 and 9.5~K
;

_aflow_title 'Rinneite (K$_{3}$NaFeCl$_{6}$) Structure'
_aflow_proto 'A6BC3D_hR22_167_f_b_e_a'
_aflow_params 'a,c/a,x_{3},x_{4},y_{4},z_{4}'
_aflow_params_values '12.033,1.15208177512,0.62611,0.78748,-0.41951,0.45463'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hR22'

_symmetry_space_group_name_H-M "R -3 2/c (hexagonal axes)"
_symmetry_Int_Tables_number 167

_cell_length_a 12.03300
_cell_length_b 12.03300
_cell_length_c 13.86300
_cell_angle_alpha 90.00000
_cell_angle_beta 90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 y,x,-z+1/2
5 -x,-x+y,-z+1/2
6 x-y,-y,-z+1/2
7 -x,-y,-z
8 y,-x+y,-z
9 x-y,x,-z
10 -y,-x,z+1/2
11 x,x-y,z+1/2
12 -x+y,y,z+1/2
13 x+1/3,y+2/3,z+2/3
14 -y+1/3,x-y+2/3,z+2/3
15 -x+y+1/3,-x+2/3,z+2/3
16 y+1/3,x+2/3,-z+1/6
17 -x+1/3,-x+y+2/3,-z+1/6
18 x-y+1/3,-y+2/3,-z+1/6
19 -x+1/3,-y+2/3,-z+2/3
20 y+1/3,-x+y+2/3,-z+2/3
21 x-y+1/3,x+2/3,-z+2/3
22 -y+1/3,-x+2/3,z+1/6
23 x+1/3,x-y+2/3,z+1/6
24 -x+y+1/3,y+2/3,z+1/6
25 x+2/3,y+1/3,z+1/3
26 -y+2/3,x-y+1/3,z+1/3
27 -x+y+2/3,-x+1/3,z+1/3
28 y+2/3,x+1/3,-z+5/6
29 -x+2/3,-x+y+1/3,-z+5/6
30 x-y+2/3,-y+1/3,-z+5/6
31 -x+2/3,-y+1/3,-z+1/3
32 y+2/3,-x+y+1/3,-z+1/3
33 x-y+2/3,x+1/3,-z+1/3
34 -y+2/3,-x+1/3,z+5/6
35 x+2/3,x-y+1/3,z+5/6
36 -x+y+2/3,y+1/3,z+5/6

loop_
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_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Na1 Na 6 a 0.00000 0.00000 0.25000 1.00000
Fe1 Fe 6 b 0.00000 0.00000 0.00000 1.00000
K1 K 18 e 0.62389 0.00000 0.25000 1.00000
Cl1 Cl 36 f 0.51328 0.69371 0.27420 1.00000
```

Rinneite ($K_3NaFeCl_6$): A6BC3D_hR22_167_f_b_e_a - POSCAR

```
A6BC3D_hR22_167_f_b_e_a & a,c/a,x3,x4,y4,z4 --params=12.033,1.15208177512,0.62611,0.78748,-0.41951,0.45463 & R-3c D_{3d}^{6} #167 (abef) & hR22 & None & Cl6FeK3Na & Rinneite & B. N. Figgis et al., Acta Crystallogr. C 56, e228-e229 (2000)
   1.00000000000000
   6.01650000000000  -3.47362789457938   4.62100000000000
   0.00000000000000   6.94725578915877   4.62100000000000
  -6.01650000000000  -3.47362789457938   4.62100000000000
     Cl     Fe     K     Na
     12      2      6      2
Direct
   0.78748000000000  -0.41951000000000   0.45463000000000     Cl    (12f)
   0.45463000000000   0.78748000000000  -0.41951000000000     Cl    (12f)
  -0.41951000000000   0.45463000000000   0.78748000000000     Cl    (12f)
   0.04537000000000   0.91951000000000  -0.28748000000000     Cl    (12f)
   0.91951000000000  -0.28748000000000   0.04537000000000     Cl    (12f)
  -0.28748000000000   0.04537000000000   0.91951000000000     Cl    (12f)
  -0.78748000000000   0.41951000000000  -0.45463000000000     Cl    (12f)
  -0.45463000000000  -0.78748000000000   0.41951000000000     Cl    (12f)
   0.41951000000000  -0.45463000000000  -0.78748000000000     Cl    (12f)
   0.95463000000000   0.08049000000000   1.28748000000000     Cl    (12f)
   0.08049000000000   1.28748000000000   0.95463000000000     Cl    (12f)
```

```
   1.28748000000000   0.95463000000000   0.08049000000000   Cl   (12f)
   0.00000000000000   0.00000000000000   0.00000000000000   Fe    (2b)
   0.50000000000000   0.50000000000000   0.50000000000000   Fe    (2b)
   0.62611000000000  -0.12611000000000   0.25000000000000   K     (6e)
   0.25000000000000   0.62611000000000  -0.12611000000000   K     (6e)
  -0.12611000000000   0.25000000000000   0.62611000000000   K     (6e)
  -0.62611000000000   0.12611000000000   0.75000000000000   K     (6e)
   0.75000000000000  -0.62611000000000   0.12611000000000   K     (6e)
   0.12611000000000   0.75000000000000  -0.62611000000000   K     (6e)
   0.25000000000000   0.25000000000000   0.25000000000000   Na    (2a)
   0.75000000000000   0.75000000000000   0.75000000000000   Na    (2a)
```

$Cs_3Tl_2Cl_9$ ($K7_2$): A9B3C2_hR28_167_ef_e_c - CIF

```
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_chemical_name_mineral 'Cl9Cs3Tl2'
_chemical_formula_sum 'Cl9 Cs3 Tl2'

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_publ_author_name
 'J. L. Hoard'
 'L. Goldstein'
_journal_name_full_name
;
 Journal of Chemical Physics
;
_journal_volume 3
_journal_year 1935
_journal_page_first 199
_journal_page_last 202
_publ_Section_title
;
 The Crystal Structure of Cesium Enneachlordithalliate, Cs$_{3}$Tl$_{2}
    ↪ $Cl$_{9}$
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'Cs$_{3}$Tl$_{2}$Cl$_{9}$ ($K7_{2}$) Structure'
_aflow_proto 'A9B3C2_hR28_167_ef_e_c'
_aflow_params 'a,c/a,x_{1},x_{2},x_{3},x_{4},y_{4},z_{4}'
_aflow_params_values '12.79567,1.4299790476,0.348,0.097,0.58333,0.586,-
    ↪ 0.755,0.42601'
_aflow_Strukturbericht '$K7_{2}$'
_aflow_Pearson 'hR28'

_symmetry_space_group_name_H-M "R -3 2/c (hexagonal axes)"
_symmetry_Int_Tables_number 167

_cell_length_a    12.79567
_cell_length_b    12.79567
_cell_length_c    18.29754
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 y,x,-z+1/2
5 -x,-x+y,-z+1/2
6 x-y,-y,-z+1/2
7 -x,-y,-z
8 y,-x+y,-z
9 x-y,x,-z
10 -y,-x,z+1/2
11 x,x-y,z+1/2
12 -x+y,y,z+1/2
13 x+1/3,y+2/3,z+2/3
14 -y+1/3,x-y+2/3,z+2/3
15 -x+y+1/3,-x+2/3,z+2/3
16 y+1/3,x+2/3,-z+1/6
17 -x+1/3,-x+y+2/3,-z+1/6
18 x-y+1/3,-y+2/3,-z+1/6
19 -x+1/3,-y+2/3,-z+2/3
20 y+1/3,-x+y+2/3,-z+2/3
21 x-y+1/3,x+2/3,-z+2/3
22 -y+1/3,-x+2/3,z+1/6
23 x+1/3,x-y+2/3,z+1/6
24 -x+y+1/3,y+2/3,z+1/6
25 x+2/3,y+1/3,z+1/3
26 -y+2/3,x-y+1/3,z+1/3
27 -x+y+2/3,-x+1/3,z+1/3
28 y+2/3,x+1/3,-z+5/6
29 -x+2/3,-x+y+1/3,-z+5/6
30 x-y+2/3,-y+1/3,-z+5/6
31 -x+2/3,-y+1/3,-z+1/3
32 y+2/3,-x+y+1/3,-z+1/3
33 x-y+2/3,x+1/3,-z+1/3
34 -y+2/3,-x+1/3,z+5/6
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36 -x+y+2/3,y+1/3,z+5/6

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_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Tl1 Tl 12 c 0.00000 0.00000 0.34800 1.00000
Cl1 Cl 18 e 0.15300 0.00000 0.25000 1.00000
Cs1 Cs 18 e 0.66667 0.00000 0.25000 1.00000
Cl2 Cl 36 f 0.50033 0.84067 0.08567 1.00000
```

$Cs_3Tl_2Cl_9$ ($K7_2$): A9B3C2_hR28_167_ef_e_c - POSCAR

```
A9B3C2_hR28_167_ef_e_c & a,c/a,x1,x2,x3,x4,y4,z4 --params=12.79567,
    ↪ 1.4299790476,0.348,0.097,0.58333,0.586,-0.755,0.42601 & R-3c D_
    ↪ {3d}^{6} #167 (ce^2f) & hR28 & $K7_{2}$ & Cl9Cs3Tl2 & Cl9Cs3Tl2
    ↪ & J. L. Hoard and L. Goldstein, J. Chem. Phys. 3, 199-202 (
    ↪ 1935)
   1.00000000000000
   6.39783500000000  -3.69379175948081   6.09918000000000
   0.00000000000000   7.38758351896162   6.09918000000000
  -6.39783500000000  -3.69379175948081   6.09918000000000
     Cl    Cs    Tl
     18     6     4
Direct
   0.09700000000000   0.40300000000000   0.25000000000000   Cl    (6e)
   0.25000000000000   0.09700000000000   0.40300000000000   Cl    (6e)
   0.40300000000000   0.25000000000000   0.09700000000000   Cl    (6e)
  -0.09700000000000   0.59700000000000   0.75000000000000   Cl    (6e)
   0.75000000000000  -0.09700000000000   0.59700000000000   Cl    (6e)
   0.59700000000000   0.75000000000000  -0.09700000000000   Cl    (6e)
   0.58600000000000  -0.75500000000000   0.42601000000000   Cl   (12f)
   0.42601000000000   0.58600000000000  -0.75500000000000   Cl   (12f)
  -0.75500000000000   0.42601000000000   0.58600000000000   Cl   (12f)
   0.07399000000000   1.25500000000000  -0.08600000000000   Cl   (12f)
   1.25500000000000  -0.08600000000000   0.07399000000000   Cl   (12f)
  -0.08600000000000   0.07399000000000   1.25500000000000   Cl   (12f)
  -0.58600000000000   0.75500000000000  -0.42601000000000   Cl   (12f)
  -0.42601000000000  -0.58600000000000   0.75500000000000   Cl   (12f)
   0.75500000000000  -0.42601000000000  -0.58600000000000   Cl   (12f)
   0.92601000000000  -0.25500000000000   1.08600000000000   Cl   (12f)
  -0.25500000000000   1.08600000000000   0.92601000000000   Cl   (12f)
   1.08600000000000   0.92601000000000  -0.25500000000000   Cl   (12f)
   0.58333000000000  -0.08333000000000   0.25000000000000   Cs    (6e)
   0.25000000000000   0.58333000000000  -0.08333000000000   Cs    (6e)
  -0.08333000000000   0.25000000000000   0.58333000000000   Cs    (6e)
  -0.58333000000000   0.08333000000000   0.75000000000000   Cs    (6e)
   0.75000000000000  -0.58333000000000   0.08333000000000   Cs    (6e)
   0.08333000000000   0.75000000000000  -0.58333000000000   Cs    (6e)
   0.34800000000000   0.34800000000000   0.34800000000000   Tl    (4c)
   0.15200000000000   0.15200000000000   0.15200000000000   Tl    (4c)
  -0.34800000000000  -0.34800000000000  -0.34800000000000   Tl    (4c)
   0.84800000000000   0.84800000000000   0.84800000000000   Tl    (4c)
```

Crancrinite ($Na_6Ca_2Al_6Si_6O_{24}(CO_3)_2$, $S3_3$ (I)): A3BCD3E15F3_hP52_173_c_b_b_c_5c_c - CIF

```
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_audit_creation_method FINDSYM

_chemical_name_mineral 'Crancrinite'
_chemical_formula_sum 'Al3 C Ca Na3 O15 Si3'

loop_
_publ_author_name
 'S. K\^{o}zu'
 'K. Takan\\'{e}'
_journal_name_full_name
;
 Proceedings of the Imperial Academy (Japan)
;
_journal_volume 9
_journal_year 1933
_journal_page_first 56
_journal_page_last 59
_publ_Section_title
;
 Crystal Structure of Cancrinite from D\^{o}d\^{o
;

# Found in Strukturbericht Band III 1933-1935, 1937

_aflow_title 'Crancrinite (Na$_{6}$Ca$_{2}$Al$_{6}$Si$_{6}$O$_{24}$(CO$_
    ↪ {3}$)$_{2}$, $S3_{3}$ (I)) Structure'
_aflow_proto 'A3BCD3E15F3_hP52_173_c_b_b_c_5c_c'
_aflow_params 'a,c/a,z_{1},z_{2},x_{3},y_{3},z_{3},x_{4},y_{4},z_{4},x_{
    ↪ 5},y_{5},z_{5},x_{6},y_{6},z_{6},x_{7},y_{7},z_{7},x_{8},y_{8},
    ↪ z_{8},x_{9},y_{9},z_{9},x_{10},y_{10},z_{10}'
_aflow_params_values '12.72,0.407232704403,0.36,0.86,0.26,0.23,0.24,0.5,
    ↪ 0.5,0.22,0.05,0.36,0.01,0.36,0.32,-0.03,0.17,0.27,0.26,0.87,
    ↪ 0.16,0.24,0.2,0.64,0.36,0.033,0.26,0.26'
_aflow_Strukturbericht '$S3_{3}$ (I)'
_aflow_Pearson 'hP52'

_symmetry_space_group_name_H-M "P 63"
_symmetry_Int_Tables_number 173

_cell_length_a    12.72000
_cell_length_b    12.72000
_cell_length_c    5.18000
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
```

```
2 x-y,x,z+1/2
3 -y,x-y,z
4 -x,-y,z+1/2
5 -x+y,-x,z
6 y,-x+y,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
C1 C 2 b 0.33333 0.66667 0.36000 1.00000
Ca1 Ca 2 b 0.33333 0.66667 0.86000 1.00000
Al1 Al 6 c 0.26000 0.23000 0.24000 1.00000
Na1 Na 6 c 0.50000 0.50000 0.22000 1.00000
O1 O 6 c 0.05000 0.36000 0.01000 1.00000
O2 O 6 c 0.36000 0.32000 -0.03000 1.00000
O3 O 6 c 0.17000 0.27000 0.26000 1.00000
O4 O 6 c 0.87000 0.16000 0.24000 1.00000
O5 O 6 c 0.20000 0.64000 0.36000 1.00000
Si1 Si 6 c 0.03300 0.26000 0.26000 1.00000
```

Crancrinite ($Na_6Ca_2Al_6Si_6O_{24}(CO_3)_2$, $S3_3$ (I)): A3BCD3E15F3_hP52_173_c_b_b_c_5c_c - POSCAR

```
A3BCD3E15F3_hP52_173_c_b_b_c_5c_c & a,c/a,z1,z2,x3,y3,z3,x4,y4,z4,x5,y5,
    ↪ z5,x6,y6,z6,x7,y7,z7,x8,y8,z8,x9,y9,z9,x10,y10,z10 --params=
    ↪ 12.72,0.407232704403,0.36,0.86,0.26,0.23,0.24,0.5,0.5,0.22,0.05
    ↪ ,0.36,0.01,0.36,0.32,-0.03,0.17,0.27,0.26,0.87,0.16,0.24,0.2,
    ↪ 0.64,0.36,0.033,0.26,0.26 & P6_{3} C_{6}^{6} #173 (b^2c^8) &
    ↪ hP52 & $S3_{3}$ (I) & Al3CCaNa3O15Si3 & Crancrinite & S. K\^{o}
    ↪ zu and K. Takan\'{e}, Proceedings of the Imperial Academy (
    ↪ Japan) 9, 56-59 (1933)
   1.00000000000000
   6.36000000000000 -11.01584313613810    0.00000000000000
   6.36000000000000  11.01584313613810    0.00000000000000
   0.00000000000000   0.00000000000000    5.18000000000000
     Al    C    Ca    Na    O    Si
      6    2     2     6    30    6
Direct
   0.26000000000000   0.23000000000000   0.24000000000000   Al   (6c)
  -0.23000000000000   0.03000000000000   0.24000000000000   Al   (6c)
  -0.03000000000000  -0.26000000000000   0.24000000000000   Al   (6c)
  -0.26000000000000  -0.23000000000000   0.74000000000000   Al   (6c)
   0.23000000000000  -0.03000000000000   0.74000000000000   Al   (6c)
   0.03000000000000   0.26000000000000   0.74000000000000   Al   (6c)
   0.33333333333333   0.66666666666667   0.36000000000000   C    (2b)
   0.66666666666667   0.33333333333333   0.86000000000000   C    (2b)
   0.33333333333333   0.66666666666667   0.86000000000000   Ca   (2b)
   0.66666666666667   0.33333333333333   1.36000000000000   Ca   (2b)
   0.50000000000000   0.50000000000000   0.22000000000000   Na   (6c)
  -0.50000000000000   0.00000000000000   0.22000000000000   Na   (6c)
   0.00000000000000  -0.50000000000000   0.22000000000000   Na   (6c)
  -0.50000000000000  -0.50000000000000   0.72000000000000   Na   (6c)
   0.50000000000000   0.00000000000000   0.72000000000000   Na   (6c)
   0.00000000000000   0.50000000000000   0.72000000000000   Na   (6c)
   0.05000000000000   0.36000000000000   0.01000000000000   O    (6c)
  -0.36000000000000  -0.31000000000000   0.01000000000000   O    (6c)
   0.31000000000000  -0.05000000000000   0.01000000000000   O    (6c)
  -0.05000000000000  -0.36000000000000   0.51000000000000   O    (6c)
   0.36000000000000   0.31000000000000   0.51000000000000   O    (6c)
  -0.31000000000000   0.05000000000000   0.51000000000000   O    (6c)
   0.36000000000000   0.32000000000000  -0.03000000000000   O    (6c)
  -0.32000000000000   0.04000000000000  -0.03000000000000   O    (6c)
  -0.04000000000000  -0.36000000000000  -0.03000000000000   O    (6c)
  -0.36000000000000  -0.32000000000000   0.47000000000000   O    (6c)
   0.32000000000000  -0.04000000000000   0.47000000000000   O    (6c)
   0.04000000000000   0.36000000000000   0.47000000000000   O    (6c)
   0.17000000000000   0.27000000000000   0.26000000000000   O    (6c)
  -0.27000000000000  -0.10000000000000   0.26000000000000   O    (6c)
   0.10000000000000  -0.17000000000000   0.26000000000000   O    (6c)
  -0.17000000000000  -0.27000000000000   0.76000000000000   O    (6c)
   0.27000000000000   0.10000000000000   0.76000000000000   O    (6c)
  -0.10000000000000   0.17000000000000   0.76000000000000   O    (6c)
   0.87000000000000   0.16000000000000   0.24000000000000   O    (6c)
  -0.16000000000000   0.71000000000000   0.24000000000000   O    (6c)
  -0.71000000000000  -0.87000000000000   0.24000000000000   O    (6c)
  -0.87000000000000  -0.16000000000000   0.74000000000000   O    (6c)
   0.16000000000000  -0.71000000000000   0.74000000000000   O    (6c)
   0.71000000000000   0.87000000000000   0.74000000000000   O    (6c)
   0.20000000000000   0.64000000000000   0.36000000000000   O    (6c)
  -0.64000000000000  -0.44000000000000   0.36000000000000   O    (6c)
   0.44000000000000  -0.20000000000000   0.36000000000000   O    (6c)
  -0.20000000000000  -0.64000000000000   0.86000000000000   O    (6c)
   0.64000000000000   0.44000000000000   0.86000000000000   O    (6c)
  -0.44000000000000   0.20000000000000   0.86000000000000   O    (6c)
   0.03300000000000   0.26000000000000   0.26000000000000   Si   (6c)
  -0.26000000000000  -0.22700000000000   0.26000000000000   Si   (6c)
   0.22700000000000  -0.03300000000000   0.26000000000000   Si   (6c)
  -0.03300000000000  -0.26000000000000   0.76000000000000   Si   (6c)
   0.26000000000000   0.22700000000000   0.76000000000000   Si   (6c)
  -0.22700000000000   0.03300000000000   0.76000000000000   Si   (6c)
```

La_3CuSiS_7: AB3C7D_hP24_173_a_c_b2c_b - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'CuLa3S7Si'
_chemical_formula_sum 'Cu La3 S7 Si'

loop_
_publ_author_name
 'G. Collin'
 'P. Laruelle'
_journal_name_full_name
;
 Bulletin de la Societ{\'e} fran{\c{c}}aise de Mineralogie et de
    ↪ Crystallographie
;
_journal_volume 94
_journal_year 1971
_journal_page_first 175
_journal_page_last 176
_publ_Section_title
;
 Structure de La$_{6}$Cu$_{2}$Si$_{2}$S$_{14}$
;

# Found in Crystal structure and magnetic properties of Ce$_{3}$CuSnSe$_
    ↪ {7}$, 2005

_aflow_title 'La$_{3}$CuSiS$_{7}$ Structure'
_aflow_proto 'AB3C7D_hP24_173_a_c_b2c_b'
_aflow_params 'a,c/a,z_{1},z_{2},z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},z_{
    ↪ 5},x_{6},y_{6},z_{6}'
_aflow_params_values '10.31,0.561978661494,0.278,0.024,0.664,0.123,0.357
    ↪ ,0.25,0.085,0.25,0.761,0.41,0.526,0.523'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hP24'

_symmetry_space_group_name_H-M "P 63"
_symmetry_Int_Tables_number 173

_cell_length_a     10.31000
_cell_length_b     10.31000
_cell_length_c     5.79400
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z+1/2
3 -y,x-y,z
4 -x,-y,z+1/2
5 -x+y,-x,z
6 y,-x+y,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cu1 Cu 2 a 0.00000 0.00000 0.27800 1.00000
S1 S 2 b 0.33333 0.66667 0.02400 1.00000
Si1 Si 2 b 0.33333 0.66667 0.66400 1.00000
La1 La 6 c 0.12300 0.35700 0.25000 1.00000
S2 S 6 c 0.08500 0.25000 0.76100 1.00000
S3 S 6 c 0.41000 0.52600 0.52300 1.00000
```

La_3CuSiS_7: AB3C7D_hP24_173_a_c_b2c_b - POSCAR

```
AB3C7D_hP24_173_a_c_b2c_b & a,c/a,z1,z2,z3,x4,y4,z4,x5,y5,z5,x6,y6,z6 --
    ↪ params=10.31,0.561978661494,0.278,0.024,0.664,0.123,0.357,0.25,
    ↪ 0.085,0.25,0.761,0.41,0.526,0.523 & P6_{3} C_{6}^{6} #173 (ab^
    ↪ 2c^3) & hP24 & None & CuLa3S7Si & CuLa3S7Si & G. Collin and P.
    ↪ Laruelle, Bull. Soc. fr. Min'eral. Crystallogr. 94, 175-176 (
    ↪ 1971)
   1.00000000000000
   5.15500000000000  -8.92872191301756    0.00000000000000
   5.15500000000000   8.92872191301756    0.00000000000000
   0.00000000000000   0.00000000000000    5.79400000000000
     Cu    La    S    Si
      2     6    14    2
Direct
   0.00000000000000   0.00000000000000   0.27800000000000   Cu   (2a)
   0.00000000000000   0.00000000000000   0.77800000000000   Cu   (2a)
   0.12300000000000   0.35700000000000   0.25000000000000   La   (6c)
  -0.35700000000000  -0.23400000000000   0.25000000000000   La   (6c)
   0.23400000000000  -0.12300000000000   0.25000000000000   La   (6c)
  -0.12300000000000  -0.35700000000000   0.75000000000000   La   (6c)
   0.35700000000000   0.23400000000000   0.75000000000000   La   (6c)
  -0.23400000000000   0.12300000000000   0.75000000000000   La   (6c)
   0.33333333333333   0.66666666666667   0.02400000000000   S    (2b)
   0.66666666666667   0.33333333333333   0.52400000000000   S    (2b)
   0.08500000000000   0.25000000000000   0.76100000000000   S    (6c)
  -0.25000000000000  -0.16500000000000   0.76100000000000   S    (6c)
   0.16500000000000  -0.08500000000000   0.76100000000000   S    (6c)
  -0.08500000000000  -0.25000000000000   1.26100000000000   S    (6c)
   0.25000000000000   0.16500000000000   1.26100000000000   S    (6c)
  -0.16500000000000   0.08500000000000   1.26100000000000   S    (6c)
   0.41000000000000   0.52600000000000   0.52300000000000   S    (6c)
  -0.52600000000000  -0.11600000000000   0.52300000000000   S    (6c)
   0.11600000000000  -0.41000000000000   0.52300000000000   S    (6c)
  -0.41000000000000  -0.52600000000000   1.02300000000000   S    (6c)
   0.52600000000000   0.11600000000000   1.02300000000000   S    (6c)
  -0.11600000000000   0.41000000000000   1.02300000000000   S    (6c)
   0.33333333333333   0.66666666666667   0.66400000000000   Si   (2b)
   0.66666666666667   0.33333333333333   1.16400000000000   Si   (2b)
```

La_3BWO_9 ($P6_3$): AB3C9D_hP28_173_a_c_3c_b - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'BLa3O9W'
_chemical_formula_sum 'B La3 O9 W'

loop_
_publ_author_name
 'J. Han'
 'F. Pan'
 'M. S. Molokeev'
 'J. Dai'
 'M. Peng'
 'W. Zhou'
 'J. Wang'
_journal_name_full_name
;
 ACS Applied Materials and Interfaces
;
_journal_volume 10
_journal_year 2018
_journal_page_first 13660
_journal_page_last 13668
_publ_Section_title
;
 Redefinition of Crystal Structure and Bi$^{3+}$ Yellow Luminescence
       ↪ with Strong Near-Ultraviolet Excitation in La$_{3}$BWO$_{9}$:
       ↪ Bi$^{3+}$ Phosphor for White Light-Emitting Diodes
;

_aflow_title 'La$_{3}$BWO$_{9}$ ($P6_{3}$) Structure'
_aflow_proto 'AB3C9D_hP28_173_a_c_3c_b'
_aflow_params 'a,c/a,z_{1},z_{2},x_{3},y_{3},z_{3},x_{4},y_{4},z_{4},x_{
      ↪ 5},y_{5},z_{5},x_{6},y_{6},z_{6}'
_aflow_params_values '8.84324,0.630487242232,0.358,0.25,0.3636,0.0854,
      ↪ 0.2213,0.17,0.047,0.861,0.192,0.483,0.044,0.142,0.517,0.468'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hP28'

_symmetry_space_group_name_H-M "P 63"
_symmetry_Int_Tables_number 173

_cell_length_a    8.84324
_cell_length_b    8.84324
_cell_length_c    5.57555
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z+1/2
3 -y,x-y,z
4 -x,-y,z+1/2
5 -x+y,-x,z
6 y,-x+y,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
B1 B 2 a 0.00000 0.00000 0.35800 1.00000
W1 W 2 b 0.33333 0.66667 0.25000 1.00000
La1 La 6 c 0.36360 0.08540 0.22130 1.00000
O1 O 6 c 0.17000 0.04700 0.86100 1.00000
O2 O 6 c 0.19200 0.48300 0.04400 1.00000
O3 O 6 c 0.14200 0.51700 0.46800 1.00000
```

La_3BWO_9 ($P6_3$): AB3C9D_hP28_173_a_c_3c_b - POSCAR

```
AB3C9D_hP28_173_a_c_3c_b & a,c/a,z1,z2,x3,y3,z3,x4,y4,z4,x5,y5,z5,x6,y6,
      ↪ z6 --params=8.84324,0.630487242232,0.358,0.25,0.3636,0.0854,
      ↪ 0.2213,0.17,0.047,0.861,0.192,0.483,0.044,0.142,0.517,0.468 &
      ↪ P6_{3} C_{6}^{6} #173 (abc^4) & hP28 & None & BLa3O9W & BLa3O9W
      ↪ & J. Han et al., ACS Appl. Mater. Interfaces 10, 13660-13668 (
      ↪ 2018)
   1.00000000000000
   4.42162000000000   -7.65847049176270    0.00000000000000
   4.42162000000000    7.65847049176270    0.00000000000000
   0.00000000000000    0.00000000000000    5.57555000000000
     B     La     O     W
     2      6     18     2
Direct
   0.00000000000000    0.00000000000000    0.35800000000000    B    (2a)
   0.00000000000000    0.00000000000000    0.85800000000000    B    (2a)
   0.36360000000000    0.08540000000000    0.22130000000000    La   (6c)
  -0.08540000000000    0.27820000000000    0.22130000000000    La   (6c)
  -0.27820000000000   -0.36360000000000    0.22130000000000    La   (6c)
  -0.36360000000000   -0.08540000000000    0.72130000000000    La   (6c)
   0.08540000000000   -0.27820000000000    0.72130000000000    La   (6c)
   0.27820000000000    0.36360000000000    0.72130000000000    La   (6c)
   0.17000000000000    0.04700000000000    0.86100000000000    O    (6c)
  -0.04700000000000    0.12300000000000    0.86100000000000    O    (6c)
  -0.12300000000000   -0.17000000000000    0.86100000000000    O    (6c)
  -0.17000000000000   -0.04700000000000    1.36100000000000    O    (6c)
   0.04700000000000   -0.12300000000000    1.36100000000000    O    (6c)
   0.12300000000000    0.17000000000000    1.36100000000000    O    (6c)
   0.19200000000000    0.48300000000000    0.04400000000000    O    (6c)
  -0.48300000000000   -0.29100000000000    0.04400000000000    O    (6c)
   0.29100000000000   -0.19200000000000    0.04400000000000    O    (6c)
  -0.19200000000000   -0.48300000000000    0.54400000000000    O    (6c)
   0.48300000000000    0.29100000000000    0.54400000000000    O    (6c)
  -0.29100000000000    0.19200000000000    0.54400000000000    O    (6c)
   0.14200000000000    0.51700000000000    0.46800000000000    O    (6c)
  -0.51700000000000   -0.37500000000000    0.46800000000000    O    (6c)
   0.37500000000000   -0.14200000000000    0.46800000000000    O    (6c)
  -0.14200000000000   -0.51700000000000    0.96800000000000    O    (6c)
   0.51700000000000    0.37500000000000    0.96800000000000    O    (6c)
  -0.37500000000000    0.14200000000000    0.96800000000000    O    (6c)
   0.33333333333333    0.66666666666667    0.25000000000000    W    (2b)
   0.66666666666667    0.33333333333333    0.75000000000000    W    (2b)
```

α-$LiIO_3$: ABC3_hP10_173_b_a_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'ILiO3'
_chemical_formula_sum 'I Li O3'

loop_
_publ_author_name
 'A. Rosenzweig'
 'B. Morosin'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 20
_journal_year 1966
_journal_page_first 758
_journal_page_last 761
_publ_Section_title
;
 A reinvestigation of the crystal structure of LiIO$_{3}$
;

_aflow_title '$\alpha$-LiIO$_{3}$ Structure'
_aflow_proto 'ABC3_hP10_173_b_a_c'
_aflow_params 'a,c/a,z_{1},z_{2},x_{3},y_{3},z_{3}'
_aflow_params_values '5.1815,0.997954260349,0.8907,0.0,0.09358,0.34396,
      ↪ 0.1698'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hP10'

_symmetry_space_group_name_H-M "P 63"
_symmetry_Int_Tables_number 173

_cell_length_a    5.18150
_cell_length_b    5.18150
_cell_length_c    5.17090
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z+1/2
3 -y,x-y,z
4 -x,-y,z+1/2
5 -x+y,-x,z
6 y,-x+y,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Li1 Li 2 a 0.00000 0.00000 0.89070 1.00000
I1 I 2 b 0.33333 0.66667 0.00000 1.00000
O1 O 6 c 0.09358 0.34396 0.16980 1.00000
```

α-$LiIO_3$: ABC3_hP10_173_b_a_c - POSCAR

```
ABC3_hP10_173_b_a_c & a,c/a,z1,z2,x3,y3,z3 --params=5.1815,
      ↪ 0.997954260349,0.8907,0.0,0.09358,0.34396,0.1698 & P6_{3} C_{6
      ↪ }^{6} #173 (abc) & hP10 & None & ILiO3 & ILiO3 & A. Rosenzweig
      ↪ and B. Morosin, Acta Cryst. 20, 758-761 (1966)
   1.00000000000000
   2.59075000000000   -4.48731062970907    0.00000000000000
   2.59075000000000    4.48731062970907    0.00000000000000
   0.00000000000000    0.00000000000000    5.17090000000000
     I     Li     O
     2      2      6
Direct
   0.33333333333333    0.66666666666667    0.00000000000000    I    (2b)
   0.66666666666667    0.33333333333333    0.50000000000000    I    (2b)
   0.00000000000000    0.00000000000000    0.89070000000000    Li   (2a)
   0.00000000000000    0.00000000000000    1.39070000000000    Li   (2a)
   0.09358000000000    0.34396000000000    0.16980000000000    O    (6c)
  -0.34396000000000   -0.25038000000000    0.16980000000000    O    (6c)
   0.25038000000000   -0.09358000000000    0.16980000000000    O    (6c)
  -0.09358000000000   -0.34396000000000    0.66980000000000    O    (6c)
```

```
   0.34396000000000   0.25038000000000   0.66980000000000   O   (6c)
  -0.25038000000000   0.09358000000000   0.66980000000000   O   (6c)
```

$LiKSO_4$ ($H1_4$): ABC4D_hP14_173_a_b_bc_b - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'KLiO4S'
_chemical_formula_sum 'K Li O4 S'

loop_
_publ_author_name
 'S. {Bhakay-Tamhane}'
 'A. Sequiera'
 'R. Chidambaram'
_journal_name_full_name
;
 Acta Crystallographica Section C: Structural Chemistry
;
_journal_volume 40
_journal_year 1984
_journal_page_first 1648
_journal_page_last 1651
_publ_Section_title
;
 Structure of lithium potassium sulphate, LiKSO${4}$: a neutron
      ↪ diffraction study
;

_aflow_title 'LiKSO${4}$ ($H1_{4}$) Structure'
_aflow_proto 'ABC4D_hP14_173_a_b_bc_b'
_aflow_params 'a,c/a,z_{1},z_{2},z_{3},z_{4},x_{5},y_{5},z_{5}'
_aflow_params_values '5.146,1.6781966576,0.0,0.3146,0.5358,0.7045,0.3446
      ↪ ,0.4031,0.7583'
_aflow_Strukturbericht '$H1_{4}$'
_aflow_Pearson 'hP14'

_symmetry_space_group_name_H-M "P 63"
_symmetry_Int_Tables_number 173

_cell_length_a     5.14600
_cell_length_b     5.14600
_cell_length_c     8.63600
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z+1/2
3 -y,x-y,z
4 -x,-y,z+1/2
5 -x+y,-x,z
6 y,-x+y,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
K1 K 2 a 0.00000 0.00000 0.00000 1.00000
Li1 Li 2 b 0.33333 0.66667 0.31460 1.00000
O1 O 2 b 0.33333 0.66667 0.53580 1.00000
S1 S 2 b 0.33333 0.66667 0.70450 1.00000
O2 O 6 c 0.34460 0.40310 0.75830 1.00000
```

$LiKSO_4$ ($H1_4$): ABC4D_hP14_173_a_b_bc_b - POSCAR

```
ABC4D_hP14_173_a_b_bc_b & a,c/a,z1,z2,z3,z4,x5,y5,z5 --params=5.146,
      ↪ 1.6781966576,0.0,0.3146,0.5358,0.7045,0.3446,0.4031,0.7583 &
      ↪ P6_{3} C_{6}^{6} #173 (ab^3c) & hP14 & $H1_{4}$ & KLiO4S &
      ↪ KLiO4S & S. {Bhakay-Tamhane} and A. Sequiera and R. Chidambaram
      ↪ , Acta Crystallogr. C 40, 1648-1651 (1984)
   1.00000000000000
   2.57300000000000  -4.45656672787472   0.00000000000000
   2.57300000000000   4.45656672787472   0.00000000000000
   0.00000000000000   0.00000000000000   8.63600000000000
     K    Li    O    S
     2     2     8     2
Direct
   0.00000000000000   0.00000000000000   0.00000000000000   K   (2a)
   0.00000000000000   0.00000000000000   0.50000000000000   K   (2a)
   0.33333333333333   0.66666666666667   0.31460000000000   Li  (2b)
   0.66666666666667   0.33333333333333   0.81460000000000   Li  (2b)
   0.33333333333333   0.66666666666667   0.53580000000000   O   (2b)
   0.66666666666667   0.33333333333333   1.03580000000000   O   (2b)
   0.34460000000000   0.40310000000000   0.75830000000000   O   (6c)
  -0.40310000000000  -0.05850000000000   0.75830000000000   O   (6c)
   0.05850000000000  -0.34460000000000   0.75830000000000   O   (6c)
  -0.34460000000000  -0.40310000000000   1.25830000000000   O   (6c)
   0.40310000000000   0.05850000000000   1.25830000000000   O   (6c)
  -0.05850000000000   0.34460000000000   1.25830000000000   O   (6c)
   0.33333333333333   0.66666666666667   0.70450000000000   S   (2b)
   0.66666666666667   0.33333333333333   1.20450000000000   S   (2b)
```

$Rh_{20}Si_{13}$: A10B7_hP34_176_c3h_b2h - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Rh20Si13'
_chemical_formula_sum 'Rh10 Si7'

loop_
_publ_author_name
 'I. Engstr\"{o}m'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 19
_journal_year 1965
_journal_page_first 1924
_journal_page_last 1932
_publ_Section_title
;
 The Crystal Structure of Rh${20}$Si${13}$
;

_aflow_title 'Rh${20}$Si${13}$ Structure'
_aflow_proto 'A10B7_hP34_176_c3h_b2h'
_aflow_params 'a,c/a,x_{3},y_{3},x_{4},y_{4},x_{5},y_{5},x_{6},y_{6},x_{
      ↪ 7},y_{7}'
_aflow_params_values '11.85,0.305738396624,0.8124,0.8157,0.5992,0.8717,
      ↪ 0.39,-0.0713,0.7667,0.1058,0.5497,0.1678'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hP34'

_symmetry_space_group_name_H-M "P 63/m"
_symmetry_Int_Tables_number 176

_cell_length_a     11.85000
_cell_length_b     11.85000
_cell_length_c     3.62300
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z+1/2
3 -y,x-y,z
4 -x,-y,z+1/2
5 -x+y,-x,z
6 y,-x+y,z+1/2
7 -x,-y,-z
8 -x+y,-x,-z+1/2
9 y,-x+y,-z
10 x,y,-z+1/2
11 x-y,x,-z
12 -y,x-y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Si1 Si 2 b 0.00000 0.00000 0.00000 0.50000
Rh1 Rh 2 c 0.33333 0.66667 0.25000 1.00000
Rh2 Rh 6 h 0.81240 0.81570 0.25000 1.00000
Rh3 Rh 6 h 0.59920 0.87170 0.25000 1.00000
Rh4 Rh 6 h 0.39000 -0.07130 0.25000 1.00000
Si2 Si 6 h 0.76670 0.10580 0.25000 1.00000
Si3 Si 6 h 0.54970 0.16780 0.25000 1.00000
```

$Rh_{20}Si_{13}$: A10B7_hP34_176_c3h_b2h - POSCAR

```
A10B7_hP34_176_c3h_b2h & a,c/a,x3,y3,x4,y4,x5,y5,x6,y6,x7,y7 --params=
      ↪ 11.85,0.305738396624,0.8124,0.8157,0.5992,0.8717,0.39,-0.0713,
      ↪ 0.7667,0.1058,0.5497,0.1678 & P6_{3}/m C_{6h}^{2} #176 (bch^5)
      ↪ & hP34 & None & Rh20Si13 & Rh20Si13 & I. Engstr\"{o}m, Acta
      ↪ Chem. Scand. 19, 1924-1932 (1965)
   1.00000000000000
   5.92500000000000 -10.26240103484560   0.00000000000000
   5.92500000000000  10.26240103484560   0.00000000000000
   0.00000000000000   0.00000000000000   3.62300000000000
    Rh    Si
    20    14
Direct
   0.33333333333333   0.66666666666667   0.25000000000000   Rh   (2c)
   0.66666666666667   0.33333333333333   0.75000000000000   Rh   (2c)
   0.81240000000000   0.81570000000000   0.25000000000000   Rh   (6h)
  -0.81570000000000  -0.00330000000000   0.25000000000000   Rh   (6h)
   0.00330000000000  -0.81240000000000   0.25000000000000   Rh   (6h)
  -0.81240000000000  -0.81570000000000   0.75000000000000   Rh   (6h)
   0.81570000000000   0.00330000000000   0.75000000000000   Rh   (6h)
  -0.00330000000000   0.81240000000000   0.75000000000000   Rh   (6h)
   0.59920000000000   0.87170000000000   0.25000000000000   Rh   (6h)
  -0.87170000000000  -0.27250000000000   0.25000000000000   Rh   (6h)
   0.27250000000000  -0.59920000000000   0.25000000000000   Rh   (6h)
  -0.59920000000000  -0.87170000000000   0.75000000000000   Rh   (6h)
   0.87170000000000   0.27250000000000   0.75000000000000   Rh   (6h)
  -0.27250000000000   0.59920000000000   0.75000000000000   Rh   (6h)
   0.39000000000000  -0.07130000000000   0.25000000000000   Rh   (6h)
   0.07130000000000   0.46130000000000   0.25000000000000   Rh   (6h)
```

```
 −0.46130000000000  −0.39000000000000   0.25000000000000  Rh  (6h)
 −0.39000000000000   0.07130000000000   0.75000000000000  Rh  (6h)
 −0.07130000000000  −0.46130000000000   0.75000000000000  Rh  (6h)
  0.46130000000000   0.39000000000000   0.75000000000000  Rh  (6h)
  0.00000000000000   0.00000000000000   0.00000000000000  Si  (2b)
  0.00000000000000   0.00000000000000   0.50000000000000  Si  (2b)
  0.76670000000000   0.10580000000000   0.25000000000000  Si  (6h)
 −0.10580000000000   0.66090000000000   0.25000000000000  Si  (6h)
 −0.66090000000000  −0.76670000000000   0.25000000000000  Si  (6h)
 −0.76670000000000  −0.10580000000000   0.75000000000000  Si  (6h)
  0.10580000000000  −0.66090000000000   0.75000000000000  Si  (6h)
  0.66090000000000   0.76670000000000   0.75000000000000  Si  (6h)
  0.54970000000000   0.16780000000000   0.25000000000000  Si  (6h)
 −0.16780000000000   0.38190000000000   0.25000000000000  Si  (6h)
 −0.38190000000000  −0.54970000000000   0.25000000000000  Si  (6h)
 −0.54970000000000  −0.16780000000000   0.75000000000000  Si  (6h)
  0.16780000000000  −0.38190000000000   0.75000000000000  Si  (6h)
  0.38190000000000   0.54970000000000   0.75000000000000  Si  (6h)
```

Th_7S_{12} ($D8_k$): A3B2_hP20_176_2h_ah - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Th7S12'
_chemical_formula_sum 'S3 Th2'

loop_
_publ_author_name
 'W. H. Zachariasen'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 2
_journal_year 1949
_journal_page_first 288
_journal_page_last 291
_publ_Section_title
;
 Crystal chemical studies of the 5f-series of elements. IX. The crystal
      ↪ structure of Th$_{7}$S$_{12}$
;

_aflow_title 'Th$_{7}$S$_{12}$ ($D8_{k}$) Structure'
_aflow_proto 'A3B2_hP20_176_2h_ah'
_aflow_params 'a,c/a,x_{2},y_{2},x_{3},y_{3},x_{4},y_{4}'
_aflow_params_values '11.064,0.361713665944,0.375,0.861,0.0,0.765,0.717,
      ↪ 0.564'
_aflow_Strukturbericht '$D8_{k}$'
_aflow_Pearson 'hP20'

_symmetry_space_group_name_H-M "P 63/m"
_symmetry_Int_Tables_number 176

_cell_length_a     11.06400
_cell_length_b     11.06400
_cell_length_c     4.00200
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z+1/2
3 -y,x-y,z
4 -x,-y,z+1/2
5 -x+y,-x,z
6 y,-x+y,z+1/2
7 -x,-y,-z
8 -x+y,-x,-z+1/2
9 y,-x+y,-z
10 x,y,-z+1/2
11 x-y,x,-z
12 -y,x-y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Th1 Th 2 a 0.00000 0.00000 0.25000 0.50000
S1 S 6 h 0.37500 0.86100 0.25000 1.00000
S2 S 6 h 0.00000 0.76500 0.25000 1.00000
Th2 Th 6 h 0.71700 0.56400 0.25000 1.00000
```

Th_7S_{12} ($D8_k$): A3B2_hP20_176_2h_ah - POSCAR

```
A3B2_hP20_176_2h_ah & a,c/a,x2,y2,x3,y3,x4,y4 --params=11.064,
      ↪ 0.361713665944,0.375,0.861,0.0,0.765,0.717,0.564 & P6_{3}/m C_{
      ↪ 6h}^{2} #176 (ah^3) & hP20 & $D8_{k}$ & Th7S12 & Th7S12 & W. H.
      ↪ Zachariasen, Acta Cryst. 2, 288-291 (1949)
   1.00000000000000
   5.53200000000000  -9.58170506747103   0.00000000000000
   5.53200000000000   9.58170506747103   0.00000000000000
   0.00000000000000   0.00000000000000   4.00200000000000
     S     Th
    12      8
Direct
   0.37500000000000   0.86100000000000   0.25000000000000  S  (6h)
  -0.86100000000000  -0.48600000000000   0.25000000000000  S  (6h)
   0.48600000000000  -0.37500000000000   0.25000000000000  S  (6h)
  -0.37500000000000  -0.86100000000000   0.75000000000000  S  (6h)
   0.86100000000000   0.48600000000000   0.75000000000000  S  (6h)
  -0.48600000000000   0.37500000000000   0.75000000000000  S  (6h)
   0.00000000000000   0.76500000000000   0.25000000000000  S  (6h)
  -0.76500000000000  -0.76500000000000   0.25000000000000  S  (6h)
   0.76500000000000   0.00000000000000   0.25000000000000  S  (6h)
   0.00000000000000  -0.76500000000000   0.75000000000000  S  (6h)
   0.76500000000000   0.76500000000000   0.75000000000000  S  (6h)
  -0.76500000000000   0.00000000000000   0.75000000000000  S  (6h)
   0.00000000000000   0.00000000000000   0.25000000000000  Th (2a)
   0.00000000000000   0.00000000000000   0.75000000000000  Th (2a)
   0.71700000000000   0.56400000000000   0.25000000000000  Th (6h)
  -0.56400000000000   0.15300000000000   0.25000000000000  Th (6h)
  -0.15300000000000  -0.71700000000000   0.25000000000000  Th (6h)
  -0.71700000000000  -0.56400000000000   0.75000000000000  Th (6h)
   0.56400000000000  -0.15300000000000   0.75000000000000  Th (6h)
   0.15300000000000   0.71700000000000   0.75000000000000  Th (6h)
```

β-Si_3N_4: A4B3_hP14_176_ch_h - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'N4Si3'
_chemical_formula_sum 'N4 Si3'

loop_
_publ_author_name
 'P. Yang'
 'H.-K. Fun'
 'I. {Ab. Rahman}'
 'I. Saleh'
_journal_name_full_name
;
 Ceramics International
;
_journal_volume 21
_journal_year 1995
_journal_page_first 137
_journal_page_last 142
_publ_Section_title
;
 Two phase refinements of the structures of $\alpha$-Si$_{3}$N$_{4}$ and
      ↪ $\beta$-Si$_{3}$N$_{4}$ made from rice husk by Rietveld
      ↪ analysis
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title '$\beta$-Si$_{3}$N$_{4}$ Structure'
_aflow_proto 'A4B3_hP14_176_ch_h'
_aflow_params 'a,c/a,x_{2},y_{2},x_{3},y_{3}'
_aflow_params_values '7.6093,0.382150789166,0.329,0.039,0.1742,0.7678'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hP14'

_symmetry_space_group_name_H-M "P 63/m"
_symmetry_Int_Tables_number 176

_cell_length_a     7.60930
_cell_length_b     7.60930
_cell_length_c     2.90790
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z+1/2
3 -y,x-y,z
4 -x,-y,z+1/2
5 -x+y,-x,z
6 y,-x+y,z+1/2
7 -x,-y,-z
8 -x+y,-x,-z+1/2
9 y,-x+y,-z
10 x,y,-z+1/2
11 x-y,x,-z
12 -y,x-y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
N1 N 2 c 0.33333 0.66667 0.25000 1.00000
N2 N 6 h 0.32900 0.03900 0.25000 1.00000
Si1 Si 6 h 0.17420 0.76780 0.25000 1.00000
```

β-Si_3N_4: A4B3_hP14_176_ch_h - POSCAR

```
A4B3_hP14_176_ch_h & a,c/a,x2,y2,x3,y3 --params=7.6093,0.382150789166,
      ↪ 0.329,0.039,0.1742,0.7678 & P6_{3}/m C_{6h}^{2} #176 (ch^2) &
```

```
      ↪ hP14 & None & N4Si3 & N4Si3 & P. Yang et al., Ceram. Int. 21,
      ↪ 137-142 (1995)
   1.00000000000000
   3.80465000000000  -6.58984710501693   0.00000000000000
   3.80465000000000   6.58984710501693   0.00000000000000
   0.00000000000000   0.00000000000000   2.90790000000000
     N    Si
     8     6
Direct
   0.33333333333333   0.66666666666667   0.25000000000000   N   (2c)
   0.66666666666667   0.33333333333333   0.75000000000000   N   (2c)
   0.32900000000000   0.03900000000000   0.25000000000000   N   (6h)
  -0.03900000000000   0.29000000000000   0.25000000000000   N   (6h)
  -0.29000000000000  -0.32900000000000   0.25000000000000   N   (6h)
  -0.32900000000000  -0.03900000000000   0.75000000000000   N   (6h)
   0.03900000000000  -0.29000000000000   0.75000000000000   N   (6h)
   0.29000000000000   0.32900000000000   0.75000000000000   N   (6h)
   0.17420000000000   0.76780000000000   0.25000000000000   Si  (6h)
  -0.76780000000000  -0.59360000000000   0.25000000000000   Si  (6h)
   0.59360000000000  -0.17420000000000   0.25000000000000   Si  (6h)
  -0.17420000000000  -0.76780000000000   0.75000000000000   Si  (6h)
   0.76780000000000   0.59360000000000   0.75000000000000   Si  (6h)
  -0.59360000000000   0.17420000000000   0.75000000000000   Si  (6h)
```

Fluorapatite [$Ca_5F(PO_4)_3$, $H5_7$]: A5BC12D3_hP42_176_fh_a_2hi_h - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Fluorapatite'
_chemical_formula_sum 'Ca5 F O12 P3'

loop_
_publ_author_name
 'J. M. Hughes'
 'J. Rakovan'
_journal_name_full_name
;
 Reviews in Mineralogy and Geochemistry
;
_journal_volume 48
_journal_year 2002
_journal_page_first 1
_journal_page_last 12
_publ_Section_title
;
 The Crystal Structure of Apatite, Ca${5}$(PO${4}$)${3}$(F,OH,Cl)
;

_aflow_title 'Fluorapatite [Ca${5}$F(PO${4}$)${3}$, $H5_{7}$]
      ↪ Structure'
_aflow_proto 'A5BC12D3_hP42_176_fh_a_2hi_h'
_aflow_params 'a,c/a,z_{2},x_{3},y_{3},x_{4},y_{4},x_{5},y_{5},x_{6},y_{
      ↪ 6},x_{7},y_{7},z_{7}'
_aflow_params_values '9.397,0.731935724167,-0.001,-0.00712,0.24227,
      ↪ 0.4849,0.3273,0.4667,0.5875,0.36895,0.3985,0.2575,0.3421,0.0705
      ↪ '
_aflow_Strukturbericht '$H5_{7}$'
_aflow_Pearson 'hP42'

_symmetry_space_group_name_H-M "P 63/m"
_symmetry_Int_Tables_number 176

_cell_length_a     9.39700
_cell_length_b     9.39700
_cell_length_c     6.87800
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z+1/2
3 -y,x-y,z
4 -x,-y,z+1/2
5 -x+y,-x,z
6 y,-x+y,z+1/2
7 -x,-y,-z
8 -x+y,-x,-z+1/2
9 y,-x+y,-z
10 x,y,-z+1/2
11 x-y,x,-z
12 -y,x-y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
F1 F 2 a 0.00000 0.00000 0.25000 1.00000
Ca1 Ca 4 f 0.33333 0.66667 -0.00100 1.00000
Ca2 Ca 6 h -0.00712 0.24227 0.25000 1.00000
O1 O 6 h 0.48490 0.32730 0.25000 1.00000
O2 O 6 h 0.46670 0.58750 0.25000 1.00000
P1 P 6 h 0.36895 0.39850 0.25000 1.00000
O3 O 12 i 0.25750 0.34210 0.07050 1.00000
```

Fluorapatite [$Ca_5F(PO_4)_3$, $H5_7$]: A5BC12D3_hP42_176_fh_a_2hi_h - POSCAR

```
A5BC12D3_hP42_176_fh_a_2hi_h & a,c/a,z2,x3,y3,x4,y4,x5,y5,x6,y6,x7,y7,z7
      ↪ --params=9.397,0.731935724167,-0.001,-0.00712,0.24227,0.4849,
      ↪ 0.3273,0.4667,0.5875,0.36895,0.3985,0.2575,0.3421,0.0705 & P6_{
      ↪ 3}/m C_{6h}^{2} #176 (afh^4i) & hP42 & $H5_{7}$ & Ca5FO12P &
      ↪ Fluorapatite & J. M. Hughes and J. Rakovan, Rev. Mineral.
      ↪ Geochem. 48, 1-12 (2002)
   1.00000000000000
   4.69850000000000  -8.13804071936237   0.00000000000000
   4.69850000000000   8.13804071936237   0.00000000000000
   0.00000000000000   0.00000000000000   6.87800000000000
    Ca    F    O    P
    10    2   24    6
Direct
   0.33333333333333   0.66666666666667  -0.00100000000000   Ca   (4f)
   0.66666666666667   0.33333333333333   0.49900000000000   Ca   (4f)
   0.66666666666667   0.33333333333333   0.00100000000000   Ca   (4f)
   0.33333333333333   0.66666666666667   0.50100000000000   Ca   (4f)
  -0.00712000000000   0.24227000000000   0.25000000000000   Ca   (6h)
  -0.24227000000000  -0.24939000000000   0.25000000000000   Ca   (6h)
   0.24939000000000   0.00712000000000   0.25000000000000   Ca   (6h)
   0.00712000000000  -0.24227000000000   0.75000000000000   Ca   (6h)
   0.24227000000000   0.24939000000000   0.75000000000000   Ca   (6h)
  -0.24939000000000  -0.00712000000000   0.75000000000000   Ca   (6h)
   0.00000000000000   0.00000000000000   0.25000000000000   F    (2a)
   0.00000000000000   0.00000000000000   0.75000000000000   F    (2a)
   0.48490000000000   0.32730000000000   0.25000000000000   O    (6h)
  -0.32730000000000   0.15760000000000   0.25000000000000   O    (6h)
  -0.15760000000000  -0.48490000000000   0.25000000000000   O    (6h)
  -0.48490000000000  -0.32730000000000   0.75000000000000   O    (6h)
   0.32730000000000  -0.15760000000000   0.75000000000000   O    (6h)
   0.15760000000000   0.48490000000000   0.75000000000000   O    (6h)
   0.46670000000000   0.58750000000000   0.25000000000000   O    (6h)
  -0.58750000000000  -0.12080000000000   0.25000000000000   O    (6h)
   0.12080000000000  -0.46670000000000   0.25000000000000   O    (6h)
  -0.46670000000000  -0.58750000000000   0.75000000000000   O    (6h)
   0.58750000000000   0.12080000000000   0.75000000000000   O    (6h)
  -0.12080000000000   0.46670000000000   0.75000000000000   O    (6h)
   0.25750000000000   0.34210000000000   0.07050000000000   O   (12i)
  -0.34210000000000  -0.08460000000000   0.07050000000000   O   (12i)
   0.08460000000000  -0.25750000000000   0.07050000000000   O   (12i)
  -0.25750000000000  -0.34210000000000   0.57050000000000   O   (12i)
   0.34210000000000   0.08460000000000   0.57050000000000   O   (12i)
  -0.08460000000000   0.25750000000000   0.57050000000000   O   (12i)
  -0.25750000000000  -0.34210000000000  -0.07050000000000   O   (12i)
   0.34210000000000   0.08460000000000  -0.07050000000000   O   (12i)
  -0.08460000000000   0.25750000000000  -0.07050000000000   O   (12i)
   0.25750000000000   0.34210000000000   0.42950000000000   O   (12i)
  -0.34210000000000  -0.08460000000000   0.42950000000000   O   (12i)
   0.08460000000000  -0.25750000000000   0.42950000000000   O   (12i)
   0.36895000000000   0.39850000000000   0.25000000000000   P    (6h)
  -0.39850000000000  -0.02955000000000   0.25000000000000   P    (6h)
   0.02955000000000  -0.36895000000000   0.25000000000000   P    (6h)
  -0.36895000000000  -0.39850000000000   0.75000000000000   P    (6h)
   0.39850000000000   0.02955000000000   0.75000000000000   P    (6h)
  -0.02955000000000   0.36895000000000   0.75000000000000   P    (6h)
```

$Fe_2(CO)_9$ ($F4_1$): A9B2C9_hP40_176_hi_f_hi - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'C9Fe2O9'
_chemical_formula_sum 'C9 Fe2 O9'

loop_
_publ_author_name
 'F. A. Cotton'
 'J. M. Troup'
_journal_name_full_name
;
 Journal of the Chemical Society, Dalton Transactions
;
_journal_volume
_journal_year 1974
_journal_page_first 800
_journal_page_last 802
_publ_Section_title
;
 Accurate determination of a classic structure in the metal carbonyl
      ↪ field: nonacarbonyldi-iron
;

# Found in Examining the structural changes in Fe${2}$(CO)${9}$ under
      ↪ high external pressures by Raman spectroscopy, 2007

_aflow_title 'Fe${2}$(CO)${9}$ ($F4_{1}$) Structure'
_aflow_proto 'A9B2C9_hP40_176_hi_f_hi'
_aflow_params 'a,c/a,z_{1},x_{2},y_{2},x_{3},y_{3},x_{4},y_{4},z_{4},x_{
      ↪ 5},y_{5},z_{5}'
_aflow_params_values '6.436,2.50512740833,0.17175,0.332,0.9098,0.331,
      ↪ 0.0895,0.338,0.4243,0.1133,0.3441,0.2817,0.0747'
_aflow_Strukturbericht '$F4_{1}$'
_aflow_Pearson 'hP40'

_symmetry_space_group_name_H-M "P 63/m"
_symmetry_Int_Tables_number 176

_cell_length_a     6.43600
_cell_length_b     6.43600
_cell_length_c     16.12300
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000
```

```
loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z+1/2
3 -y,x-y,z
4 -x,-y,z+1/2
5 -x+y,-x,z
6 y,-x+y,z+1/2
7 -x,-y,-z
8 -x+y,-x,-z+1/2
9 y,-x+y,-z
10 x,y,-z+1/2
11 x-y,x,-z
12 -y,x-y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Fe1 Fe 4 f 0.33333 0.66667 0.17175 1.00000
C1 C 6 h 0.33200 0.90980 0.25000 1.00000
O1 O 6 h 0.33100 0.08950 0.25000 1.00000
C2 C 12 i 0.33800 0.42430 0.11330 1.00000
O2 O 12 i 0.34410 0.28170 0.07470 1.00000
```

$Fe_2(CO)_9$ ($F4_1$): A9B2C9_hP40_176_hi_f_hi - POSCAR

```
A9B2C9_hP40_176_hi_f_hi & a,c/a,z1,x2,y2,x3,y3,x4,y4,z4,x5,y5,z5 --
    ↪ params=6.436,2.50512740833,0.17175,0.332,0.9098,0.331,0.0895,
    ↪ 0.338,0.4243,0.1133,0.3441,0.2817,0.0747 & P6_{3}/m C_{6h}^{2}
    ↪ #176 (fh^2i^2) & hP40 & $F4_{1}$ & C9Fe2O9 & C9Fe2O9 & F. A.
    ↪ Cotton and J. M. Troup, {J. Chem. Soc. Dalton Trans., 800-802
    ↪ (1974)
   1.00000000000000
   3.21800000000000  -5.57373949875665   0.00000000000000
   3.21800000000000   5.57373949875665   0.00000000000000
   0.00000000000000   0.00000000000000  16.12300000000000
     C     Fe     O
    18     4    18
Direct
   0.33200000000000   0.90980000000000   0.25000000000000  C   (6h)
  -0.90980000000000  -0.57780000000000   0.25000000000000  C   (6h)
   0.57780000000000  -0.33200000000000   0.25000000000000  C   (6h)
  -0.33200000000000  -0.90980000000000   0.75000000000000  C   (6h)
   0.90980000000000   0.57780000000000   0.75000000000000  C   (6h)
  -0.57780000000000   0.33200000000000   0.75000000000000  C   (6h)
   0.33800000000000   0.42430000000000   0.11330000000000  C  (12i)
  -0.42430000000000  -0.08630000000000   0.11330000000000  C  (12i)
   0.08630000000000  -0.33800000000000   0.11330000000000  C  (12i)
  -0.33800000000000  -0.42430000000000   0.61330000000000  C  (12i)
   0.42430000000000   0.08630000000000   0.61330000000000  C  (12i)
  -0.08630000000000   0.33800000000000   0.61330000000000  C  (12i)
  -0.33800000000000  -0.42430000000000  -0.11330000000000  C  (12i)
   0.42430000000000   0.08630000000000  -0.11330000000000  C  (12i)
  -0.08630000000000   0.33800000000000  -0.11330000000000  C  (12i)
   0.33800000000000   0.42430000000000   0.38670000000000  C  (12i)
  -0.42430000000000  -0.08630000000000   0.38670000000000  C  (12i)
   0.08630000000000  -0.33800000000000   0.38670000000000  C  (12i)
   0.33333333333333   0.66666666666667   0.17175000000000  Fe  (4f)
   0.66666666666667   0.33333333333333   0.67175000000000  Fe  (4f)
   0.66666666666667   0.33333333333333  -0.17175000000000  Fe  (4f)
   0.33333333333333   0.66666666666667   0.32825000000000  Fe  (4f)
   0.33100000000000   0.08950000000000   0.25000000000000  O   (6h)
  -0.08950000000000   0.24150000000000   0.25000000000000  O   (6h)
  -0.24150000000000  -0.33100000000000   0.25000000000000  O   (6h)
  -0.33100000000000  -0.08950000000000   0.75000000000000  O   (6h)
   0.08950000000000  -0.24150000000000   0.75000000000000  O   (6h)
   0.24150000000000   0.33100000000000   0.75000000000000  O   (6h)
   0.34410000000000   0.28170000000000   0.07470000000000  O  (12i)
  -0.28170000000000   0.06240000000000   0.07470000000000  O  (12i)
  -0.06240000000000  -0.34410000000000   0.07470000000000  O  (12i)
  -0.34410000000000  -0.28170000000000   0.57470000000000  O  (12i)
   0.28170000000000  -0.06240000000000   0.57470000000000  O  (12i)
   0.06240000000000   0.34410000000000   0.57470000000000  O  (12i)
  -0.34410000000000  -0.28170000000000  -0.07470000000000  O  (12i)
   0.28170000000000  -0.06240000000000  -0.07470000000000  O  (12i)
   0.06240000000000   0.34410000000000  -0.07470000000000  O  (12i)
   0.34410000000000   0.28170000000000   0.42530000000000  O  (12i)
  -0.28170000000000   0.06240000000000   0.42530000000000  O  (12i)
  -0.06240000000000  -0.34410000000000   0.42530000000000  O  (12i)
```

$K_3W_2Cl_9$ ($K7_1$): A9B3C2_hP28_176_hi_af_f - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Cl9K3W2'
_chemical_formula_sum 'Cl9 K3 W2'

loop_
_publ_author_name
 'W. H. {Watson, Jr.}'
 'J. Waser'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 11
_journal_year 1958
_journal_page_first 689
_journal_page_last 692
_publ_Section_title
;
 Refinement of the structure of tripotassiumditungsten enneachloride,
    ↪ K$_{3}$W$_{2}$Cl$_{9}$
;

_aflow_title 'K$_{3}$W$_{2}$Cl$_{9}$ ($K7_{1}$) Structure'
_aflow_proto 'A9B3C2_hP28_176_hi_af_f'
_aflow_params 'a,c/a,z_{2},z_{3},x_{4},y_{4},x_{5},y_{5},z_{5}'
_aflow_params_values '7.17,2.26638772664,0.5718,0.3241,0.4588,0.4472,
    ↪ 0.1348,0.3506,0.4074'
_aflow_Strukturbericht '$K7_{1}$'
_aflow_Pearson 'hP28'

_symmetry_space_group_name_H-M "P 63/m"
_symmetry_Int_Tables_number 176

_cell_length_a    7.17000
_cell_length_b    7.17000
_cell_length_c    16.25000
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z+1/2
3 -y,x-y,z
4 -x,-y,z+1/2
5 -x+y,-x,z
6 y,-x+y,z+1/2
7 -x,-y,-z
8 -x+y,-x,-z+1/2
9 y,-x+y,-z
10 x,y,-z+1/2
11 x-y,x,-z
12 -y,x-y,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
K1 K 2 a 0.00000 0.00000 0.25000 1.00000
K2 K 4 f 0.33333 0.66667 0.57180 1.00000
W1 W 4 f 0.33333 0.66667 0.32410 1.00000
Cl1 Cl 6 h 0.45880 0.44720 0.25000 1.00000
Cl2 Cl 12 i 0.13480 0.35060 0.40740 1.00000
```

$K_3W_2Cl_9$ ($K7_1$): A9B3C2_hP28_176_hi_af_f - POSCAR

```
A9B3C2_hP28_176_hi_af_f & a,c/a,z2,z3,x4,y4,x5,y5,z5 --params=7.17,
    ↪ 2.26638772664,0.5718,0.3241,0.4588,0.4472,0.1348,0.3506,0.4074
    ↪ & P6_{3}/m C_{6h}^{2} #176 (af^2hi) & hP28 & $K7_{1}$ & Cl9K3W2
    ↪ & Cl9K3W2 & W. H. {Watson, Jr.} and J. Waser, Acta Cryst. 11,
    ↪ 689-692 (1958)
   1.00000000000000
   3.58500000000000  -6.20940214513442   0.00000000000000
   3.58500000000000   6.20940214513442   0.00000000000000
   0.00000000000000   0.00000000000000  16.25000000000000
     Cl     K     W
    18     6     4
Direct
   0.45880000000000   0.44720000000000   0.25000000000000  Cl   (6h)
  -0.44720000000000   0.01160000000000   0.25000000000000  Cl   (6h)
  -0.01160000000000  -0.45880000000000   0.25000000000000  Cl   (6h)
  -0.45880000000000  -0.44720000000000   0.75000000000000  Cl   (6h)
   0.44720000000000  -0.01160000000000   0.75000000000000  Cl   (6h)
   0.01160000000000   0.45880000000000   0.75000000000000  Cl   (6h)
   0.13480000000000   0.35060000000000   0.40740000000000  Cl  (12i)
  -0.35060000000000  -0.21580000000000   0.40740000000000  Cl  (12i)
   0.21580000000000  -0.13480000000000   0.40740000000000  Cl  (12i)
  -0.13480000000000  -0.35060000000000   0.90740000000000  Cl  (12i)
   0.35060000000000   0.21580000000000   0.90740000000000  Cl  (12i)
  -0.21580000000000   0.13480000000000   0.90740000000000  Cl  (12i)
  -0.13480000000000  -0.35060000000000  -0.40740000000000  Cl  (12i)
   0.35060000000000   0.21580000000000  -0.40740000000000  Cl  (12i)
  -0.21580000000000   0.13480000000000  -0.40740000000000  Cl  (12i)
   0.13480000000000   0.35060000000000   0.09260000000000  Cl  (12i)
  -0.35060000000000  -0.21580000000000   0.09260000000000  Cl  (12i)
   0.21580000000000  -0.13480000000000   0.09260000000000  Cl  (12i)
   0.00000000000000   0.00000000000000   0.25000000000000  K   (2a)
   0.00000000000000   0.00000000000000   0.75000000000000  K   (2a)
   0.33333333333333   0.66666666666667   0.57180000000000  K   (4f)
   0.66666666666667   0.33333333333333   1.07180000000000  K   (4f)
   0.66666666666667   0.33333333333333  -0.57180000000000  K   (4f)
   0.33333333333333   0.66666666666667  -0.07180000000000  K   (4f)
   0.33333333333333   0.66666666666667   0.32410000000000  W   (4f)
   0.66666666666667   0.33333333333333   0.82410000000000  W   (4f)
   0.66666666666667   0.33333333333333  -0.32410000000000  W   (4f)
   0.33333333333333   0.66666666666667   0.17590000000000  W   (4f)
```

Hg_2O_2NaI: A2BCD2_hP18_180_f_c_b_i - CIF

```
# CIF file
data_findsym-output
```

```
_audit_creation_method FINDSYM

_chemical_name_mineral 'Hg2INaO2'
_chemical_formula_sum 'Hg2 I Na O2'

loop_
_publ_author_name
 'K. Aurivillius'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 18
_journal_year 1964
_journal_page_first 1305
_journal_page_last 1306
_publ_Section_title
;
 Least-Squares Refinement of the Crystal Structures of Orthorhombic HgO
      ↪ and of Hg$_{2}$O$_{2}$NaI
;

_aflow_title 'Hg$_{2}$O$_{2}$NaI Structure'
_aflow_proto 'A2BCD2_hP18_180_f_c_b_i'
_aflow_params 'a,c/a,z_{3},x_{4}'
_aflow_params_values '6.667,1.50802459877,0.3333,0.1521'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hP18'

_symmetry_space_group_name_H-M "P 62 2 2"
_symmetry_Int_Tables_number 180

_cell_length_a     6.66700
_cell_length_b     6.66700
_cell_length_c     10.05400
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z+1/3
3 -y,x-y,z+2/3
4 -x,-y,z
5 -x+y,-x,z+1/3
6 y,-x+y,z+2/3
7 x-y,-y,-z
8 x,x-y,-z+1/3
9 y,x,-z+2/3
10 -x+y,y,-z
11 -x,-x+y,-z+1/3
12 -y,-x,-z+2/3

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Na1 Na 3 b 0.00000 0.00000 0.50000 1.00000
I1 I 3 c 0.50000 0.00000 0.00000 1.00000
Hg1 Hg 6 f 0.50000 0.00000 0.33330 1.00000
O1 O 6 i 0.15210 0.30420 0.00000 1.00000
```

Hg_2O_2NaI: A2BCD2_hP18_180_f_c_b_i - POSCAR

```
A2BCD2_hP18_180_f_c_b_i & a,c/a,z3,x4 --params=6.667,1.50802459877,
      ↪ 0.3333,0.1521 & P6_{2}22 D_{6}^{4} #180 (bcfi) & hP18 & None &
      ↪ Hg2INaO2 & Hg2INaO2 & K. Aurivillius, Acta Chem. Scand. 18,
      ↪ 1305-1306 (1964)
   1.00000000000000
   3.33350000000000  -5.77379136703085   0.00000000000000
   3.33350000000000   5.77379136703085   0.00000000000000
   0.00000000000000   0.00000000000000  10.05400000000000
    Hg    I    Na    O
     6    3     3    6
Direct
   0.50000000000000   0.00000000000000   0.33330000000000  Hg  (6f)
   0.00000000000000   0.50000000000000   0.99996666666667  Hg  (6f)
   0.50000000000000   0.50000000000000   0.66663333333333  Hg  (6f)
   0.00000000000000   0.50000000000000   0.33336666666667  Hg  (6f)
   0.50000000000000   0.00000000000000  -0.33330000000000  Hg  (6f)
   0.50000000000000   0.50000000000000   0.00003333333333  Hg  (6f)
   0.50000000000000   0.00000000000000   0.00000000000000   I  (3c)
   0.00000000000000   0.50000000000000   0.66666666666667   I  (3c)
   0.50000000000000   0.50000000000000   0.33333333333333   I  (3c)
   0.00000000000000   0.00000000000000   0.50000000000000  Na  (3b)
   0.00000000000000   0.00000000000000   0.16666666666667  Na  (3b)
   0.00000000000000   0.00000000000000   0.83333333333333  Na  (3b)
   0.15210000000000   0.30420000000000   0.00000000000000   O  (6i)
  -0.30420000000000  -0.15210000000000   0.66666666666667   O  (6i)
   0.15210000000000  -0.15210000000000   0.33333333333333   O  (6i)
  -0.15210000000000  -0.30420000000000   0.00000000000000   O  (6i)
   0.30420000000000   0.15210000000000   0.66666666666667   O  (6i)
  -0.15210000000000   0.15210000000000   0.33333333333333   O  (6i)
```

$BaAl_2O_4$ ($H2_8$): A2BC6_hP18_182_f_b_gh - CIF

```
# CIF file
data_findsym-output
```

```
_audit_creation_method FINDSYM

_chemical_name_mineral 'Al2BaO4'
_chemical_formula_sum 'Al2 Ba O6'

loop_
_publ_author_name
 'A. J. Perrotta'
 'J. V. Smith'
_journal_name_full_name
;
 Bulletin de la Societ{\'e} fran{\c{c}}aise de Mineralogie et de
      ↪ Crystallographie
;
_journal_volume 91
_journal_year 1968
_journal_page_first 85
_journal_page_last 87
_publ_Section_title
;
 The Crystal Structure of BaAl$_{2}$O$_{4}$
;

_aflow_title 'BaAl$_{2}$O$_{4}$ ($H2_{8}$) Structure'
_aflow_proto 'A2BC6_hP18_182_f_b_gh'
_aflow_params 'a,c/a,z_{2},x_{3},x_{4}'
_aflow_params_values '5.218,1.68282866999,0.054,0.36,0.371'
_aflow_Strukturbericht '$H2_{8}$'
_aflow_Pearson 'hP18'

_symmetry_space_group_name_H-M "P 63 2 2"
_symmetry_Int_Tables_number 182

_cell_length_a     5.21800
_cell_length_b     5.21800
_cell_length_c     8.78100
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z+1/2
3 -y,x-y,z
4 -x,-y,z+1/2
5 -x+y,-x,z
6 y,-x+y,z+1/2
7 x-y,-y,-z
8 x,x-y,-z+1/2
9 y,x,-z
10 -x+y,y,-z+1/2
11 -x,-x+y,-z
12 -y,-x,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ba1 Ba 2 b 0.00000 0.00000 0.25000 1.00000
Al1 Al 4 f 0.33333 0.66667 0.05400 1.00000
O1 O 6 g 0.36000 0.00000 0.00000 1.00000
O2 O 6 h 0.37100 0.74200 0.25000 0.33333
```

$BaAl_2O_4$ ($H2_8$): A2BC6_hP18_182_f_b_gh - POSCAR

```
A2BC6_hP18_182_f_b_gh & a,c/a,z2,x3,x4 --params=5.218,1.68282866999,
      ↪ 0.054,0.36,0.371 & P6_{3}22 D_{6}^{6} #182 (bfgh) & hP18 & $H2_
      ↪ {8}$ & Al2BaO4 & Al2BaO4 & A. J. Perrotta and J. V. Smith,
      ↪ Bull. Soc. fr. Min'eral. Crystallogr. 91, 85-87 (1968)
   1.00000000000000
   2.60900000000000  -4.51892055694720   0.00000000000000
   2.60900000000000   4.51892055694720   0.00000000000000
   0.00000000000000   0.00000000000000   8.78100000000000
    Al    Ba    O
     4     2    12
Direct
   0.33333333333333   0.66666666666667   0.05400000000000  Al  (4f)
   0.66666666666667   0.33333333333333   0.55400000000000  Al  (4f)
   0.66666666666667   0.33333333333333  -0.05400000000000  Al  (4f)
   0.33333333333333   0.66666666666667   0.44600000000000  Al  (4f)
   0.00000000000000   0.00000000000000   0.25000000000000  Ba  (2b)
   0.00000000000000   0.00000000000000   0.75000000000000  Ba  (2b)
   0.36000000000000   0.00000000000000   0.00000000000000   O  (6g)
   0.00000000000000   0.36000000000000   0.00000000000000   O  (6g)
  -0.36000000000000  -0.36000000000000   0.00000000000000   O  (6g)
  -0.36000000000000   0.00000000000000   0.50000000000000   O  (6g)
   0.00000000000000  -0.36000000000000   0.50000000000000   O  (6g)
   0.36000000000000   0.36000000000000   0.50000000000000   O  (6g)
   0.37100000000000   0.74200000000000   0.25000000000000   O  (6h)
  -0.74200000000000  -0.37100000000000   0.25000000000000   O  (6h)
   0.37100000000000  -0.37100000000000   0.25000000000000   O  (6h)
  -0.37100000000000  -0.74200000000000   0.75000000000000   O  (6h)
   0.74200000000000   0.37100000000000   0.75000000000000   O  (6h)
  -0.37100000000000   0.37100000000000   0.75000000000000   O  (6h)
```

$E2_3$ ($LiIO_3$) (*obsolete*): ABC3_hP10_182_c_b_g - CIF

```
# CIF file
```

```
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'ILiO3'
_chemical_formula_sum 'I Li O3'

loop_
_publ_author_name
 'W. H. Zachariasen'
 'F. A. Barta'
_journal_name_full_name
;
 Physical Review
;
_journal_volume 37
_journal_year 1931
_journal_page_first 1626
_journal_page_last 1630
_publ_Section_title
;
 Crystal Structure of Lithium Iodate
;

# Found in A reinvestigation of the crystal structure of LiIO$_{3}$,
      ↪ 1966

_aflow_title '$E2_{3}$ (LiIO$_{3}$) ({\em{obsolete}}) Structure'
_aflow_proto 'ABC3_hP10_182_c_b_g'
_aflow_params 'a,c/a,x_{3}'
_aflow_params_values '5.469,0.942585481807,0.33333'
_aflow_Strukturbericht '$E2_{3}$'
_aflow_Pearson 'hP10'

_symmetry_space_group_name_H-M "P 63 2 2"
_symmetry_Int_Tables_number 182

_cell_length_a     5.46900
_cell_length_b     5.46900
_cell_length_c     5.15500
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z+1/2
3 -y,x-y,z
4 -x,-y,z+1/2
5 -x+y,-x,z
6 y,-x+y,z+1/2
7 x-y,-y,-z
8 x,x-y,-z+1/2
9 y,x,-z
10 -x+y,y,-z+1/2
11 -x,-x+y,-z
12 -y,-x,-z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Li1 Li 2 b 0.00000 0.00000 0.25000 1.00000
I1 I 2 c 0.33333 0.66667 0.25000 1.00000
O1 O 6 g 0.33333 0.00000 0.00000 1.00000
```

$E2_3$ ($LiIO_3$) (*obsolete*): ABC3_hP10_182_c_b_g - POSCAR

```
ABC3_hP10_182_c_b_g & a,c/a,x3 --params=5.469,0.942585481807,0.33333 &
      ↪ P6_{3}22 D_{6}^{6} #182 (bcg) & hP10 & $E2_{3}$ & ILiO3 & ILiO3
      ↪ & W. H. Zachariasen and F. A. Barta, Phys. Rev. 37, 1626-1630
      ↪ (1931)
   1.00000000000000
   2.73450000000000  -4.73629293329710   0.00000000000000
   2.73450000000000   4.73629293329710   0.00000000000000
   0.00000000000000   0.00000000000000   5.15500000000000
     I     Li     O
     2     2     6
Direct
   0.33333333333333   0.66666666666667   0.25000000000000   I   (2c)
   0.66666666666667   0.33333333333333   0.75000000000000   I   (2c)
   0.00000000000000   0.00000000000000   0.25000000000000   Li  (2b)
   0.00000000000000   0.00000000000000   0.75000000000000   Li  (2b)
   0.33333000000000   0.00000000000000   0.00000000000000   O   (6g)
   0.00000000000000   0.33333000000000   0.00000000000000   O   (6g)
  -0.33333000000000  -0.33333000000000   0.00000000000000   O   (6g)
  -0.33333000000000   0.00000000000000   0.50000000000000   O   (6g)
   0.00000000000000  -0.33333000000000   0.50000000000000   O   (6g)
   0.33333000000000   0.33333000000000   0.50000000000000   O   (6g)
```

$Zn_2Mo_3O_8$: A3B8C2_hP26_186_c_ab2c_2b - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Mo3O8Zn2'
_chemical_formula_sum 'Mo3 O8 Zn2'

loop_
_publ_author_name
 'G. B. Ansell'
 'L. Katz'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 21
_journal_year 1966
_journal_page_first 482
_journal_page_last 485
_publ_Section_title
;
 A Refinement of the Crystal Structure of Zinc Molybdenum(IV) Oxide,
      ↪ Zn$_{2}$Mo$_{3}$O$_{8}$
;

# Found in Structure of Kamiokite, 1986

_aflow_title 'Zn$_{2}$Mo$_{3}$O$_{8}$ Structure'
_aflow_proto 'A3B8C2_hP26_186_c_ab2c_2b'
_aflow_params 'a,c/a,z_{1},z_{2},z_{3},z_{4},x_{5},z_{5},x_{6},z_{6},x_{
      ↪ 7},z_{7}'
_aflow_params_values '5.775,1.71688311688,0.3886,0.147,-0.0535,0.5132,
      ↪ 0.1461,0.25,0.4861,0.3639,0.1647,0.6354'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hP26'

_symmetry_space_group_name_H-M "P 63 m c"
_symmetry_Int_Tables_number 186

_cell_length_a     5.77500
_cell_length_b     5.77500
_cell_length_c     9.91500
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z+1/2
3 -y,x-y,z
4 -x,-y,z+1/2
5 -x+y,-x,z
6 y,-x+y,z+1/2
7 -x+y,y,z
8 -x,-x+y,z+1/2
9 -y,-x,z
10 x-y,-y,z+1/2
11 x,x-y,z
12 y,x,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 2 a 0.00000 0.00000 0.38860 1.00000
O2 O 2 b 0.33333 0.66667 0.14700 1.00000
Zn1 Zn 2 b 0.33333 0.66667 -0.05350 1.00000
Zn2 Zn 2 b 0.33333 0.66667 0.51320 1.00000
Mo1 Mo 6 c 0.14610 0.85390 0.25000 1.00000
O3 O 6 c 0.48610 0.51390 0.36390 1.00000
O4 O 6 c 0.16470 0.83530 0.63540 1.00000
```

$Zn_2Mo_3O_8$: A3B8C2_hP26_186_c_ab2c_2b - POSCAR

```
A3B8C2_hP26_186_c_ab2c_2b & a,c/a,z1,z2,z3,z4,x5,z5,x6,z6,x7,z7 --params
      ↪ =5.775,1.71688311688,0.3886,0.147,-0.0535,0.5132,0.1461,0.25,
      ↪ 0.4861,0.3639,0.1647,0.6354 & P6_{3}mc C_{6v}^{4} #186 (ab^3c^3
      ↪ ) & hP26 & None & Mo3O8Zn2 & Mo3O8Zn2 & G. B. Ansell and L.
      ↪ Katz, Acta Cryst. 21, 482-485 (1966)
   1.00000000000000
   2.88750000000000  -5.00129670685513   0.00000000000000
   2.88750000000000   5.00129670685513   0.00000000000000
   0.00000000000000   0.00000000000000   9.91500000000000
     Mo     O     Zn
      6    16     4
Direct
   0.14610000000000  -0.14610000000000   0.25000000000000   Mo  (6c)
   0.14610000000000   0.29220000000000   0.25000000000000   Mo  (6c)
  -0.29220000000000  -0.14610000000000   0.25000000000000   Mo  (6c)
  -0.14610000000000   0.14610000000000   0.75000000000000   Mo  (6c)
  -0.14610000000000  -0.29220000000000   0.75000000000000   Mo  (6c)
   0.29220000000000   0.14610000000000   0.75000000000000   Mo  (6c)
   0.00000000000000   0.00000000000000   0.38860000000000   O   (2a)
   0.00000000000000   0.00000000000000   0.88860000000000   O   (2a)
   0.33333333333333   0.66666666666667   0.14700000000000   O   (2b)
   0.66666666666667   0.33333333333333   0.64700000000000   O   (2b)
   0.48610000000000  -0.48610000000000   0.36390000000000   O   (6c)
   0.48610000000000   0.97220000000000   0.36390000000000   O   (6c)
  -0.97220000000000  -0.48610000000000   0.36390000000000   O   (6c)
  -0.48610000000000   0.48610000000000   0.86390000000000   O   (6c)
  -0.48610000000000  -0.97220000000000   0.86390000000000   O   (6c)
   0.97220000000000   0.48610000000000   0.86390000000000   O   (6c)
   0.16470000000000  -0.16470000000000   0.63540000000000   O   (6c)
   0.16470000000000   0.32940000000000   0.63540000000000   O   (6c)
  -0.32940000000000  -0.16470000000000   0.63540000000000   O   (6c)
```

```
 −0.16470000000000   0.16470000000000   1.13540000000000   O   (6c)
 −0.16470000000000  −0.32940000000000   1.13540000000000   O   (6c)
  0.32940000000000   0.16470000000000   1.13540000000000   O   (6c)
  0.33333333333333   0.66666666666667  −0.05350000000000   Zn  (2b)
  0.66666666666667   0.33333333333333   0.44650000000000   Zn  (2b)
  0.33333333333333   0.66666666666667   0.51320000000000   Zn  (2b)
  0.66666666666667   0.33333333333333   1.01320000000000   Zn  (2b)
```

$Nd(BrO_3)_3{\cdot}9H_2O$ ($G2_2$): A3B9CD9_hP44_186_c_3c_b_cd - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Neodymium bromate enneahydrate'
_chemical_formula_sum 'Br3 (H2O)9 Nd O9'

loop_
_publ_author_name
 'L. Helmholz'
_journal_name_full_name
;
 Journal of the American Chemical Society
;
_journal_volume 61
_journal_year 1939
_journal_page_first 1544
_journal_page_last 1550
_publ_Section_title
;
 The Crystal Structure of Neodymium Bromate Enneahydrate, Nd(BrO$_{3}$)
      ↪ $_{3}$$\cdot$9H$_{2}$O
;

# Found in Strukturbericht Band VII 1939, 1943

_aflow_title 'Nd(BrO$_{3}$)$_{3}$$\cdot$9H$_{2}$O ($G2_{2}$) Structure'
_aflow_proto 'A3B9CD9_hP44_186_c_3c_b_cd'
_aflow_params 'a,c/a,z_{1},x_{2},z_{2},x_{3},z_{3},x_{4},z_{4},x_{5},z_{
      ↪ 5},x_{6},z_{6},x_{7},y_{7},z_{7}'
_aflow_params_values '11.73,0.576300085251,0.25,0.13,0.73,0.425,0.49,
      ↪ 0.425,0.01,0.215,0.25,0.105,0.53,0.065,0.365,0.75'
_aflow_Strukturbericht '$G2_{2}$'
_aflow_Pearson 'hP44'

_symmetry_space_group_name_H-M "P 63 m c"
_symmetry_Int_Tables_number 186

_cell_length_a     11.73000
_cell_length_b     11.73000
_cell_length_c     6.76000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z+1/2
3 -y,x-y,z
4 -x,-y,z+1/2
5 -x+y,-x,z
6 y,-x+y,z+1/2
7 -x+y,y,z
8 -x,-x+y,z+1/2
9 -y,-x,z
10 x-y,-y,z+1/2
11 x,x-y,z
12 y,x,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Nd1 Nd 2 b 0.33333 0.66667 0.25000 1.00000
Br1 Br 6 c 0.13000 0.87000 0.73000 1.00000
H2O1 H2O 6 c 0.42500 0.57500 0.49000 1.00000
H2O2 H2O 6 c 0.42500 0.57500 0.01000 1.00000
H2O3 H2O 6 c 0.21500 0.78500 0.25000 1.00000
O1 O 6 c 0.10500 0.89500 0.53000 1.00000
O2 O 12 d 0.06500 0.36500 0.75000 1.00000
```

$Nd(BrO_3)_3{\cdot}9H_2O$ ($G2_2$): A3B9CD9_hP44_186_c_3c_b_cd - POSCAR

```
A3B9CD9_hP44_186_c_3c_b_cd & a,c/a,z1,x2,z2,x3,z3,x4,z4,x5,z5,x6,z6,x7,
      ↪ y7,z7 --params=11.73,0.576300085251,0.25,0.13,0.73,0.425,0.49,
      ↪ 0.425,0.01,0.215,0.25,0.105,0.53,0.065,0.365,0.75 & P6_{3}mc C_
      ↪ {6v}^{4} #186 (bc^5d) & hP44 & $G2_{2}$ & Br(H2O)9NdO9 &
      ↪ Neodymium bromate enneahydrate & L. Helmholz, J. Am. Chem. Soc.
      ↪ 61, 1544-1550 (1939)
   1.00000000000000
   5.86500000000000 -10.15847798639150    0.00000000000000
   5.86500000000000  10.15847798639150    0.00000000000000
   0.00000000000000   0.00000000000000    6.76000000000000
     Br    H2O     Nd      O
      6     18      2     18
Direct
   0.13000000000000  -0.13000000000000   0.73000000000000   Br   (6c)
   0.13000000000000   0.26000000000000   0.73000000000000   Br   (6c)
  -0.26000000000000  -0.13000000000000   0.73000000000000   Br   (6c)
  -0.13000000000000   0.13000000000000   1.23000000000000   Br   (6c)
  -0.13000000000000  -0.26000000000000   1.23000000000000   Br   (6c)
   0.26000000000000   0.13000000000000   1.23000000000000   Br   (6c)
   0.42500000000000  -0.42500000000000   0.49000000000000   H2O  (6c)
   0.42500000000000   0.85000000000000   0.49000000000000   H2O  (6c)
  -0.85000000000000  -0.42500000000000   0.49000000000000   H2O  (6c)
  -0.42500000000000   0.42500000000000   0.99000000000000   H2O  (6c)
  -0.42500000000000  -0.85000000000000   0.99000000000000   H2O  (6c)
   0.85000000000000   0.42500000000000   0.99000000000000   H2O  (6c)
   0.42500000000000  -0.42500000000000   0.01000000000000   H2O  (6c)
   0.42500000000000   0.85000000000000   0.01000000000000   H2O  (6c)
  -0.85000000000000  -0.42500000000000   0.01000000000000   H2O  (6c)
  -0.42500000000000   0.42500000000000   0.51000000000000   H2O  (6c)
  -0.42500000000000  -0.85000000000000   0.51000000000000   H2O  (6c)
   0.85000000000000   0.42500000000000   0.51000000000000   H2O  (6c)
   0.21500000000000  -0.21500000000000   0.25000000000000   H2O  (6c)
   0.21500000000000   0.43000000000000   0.25000000000000   H2O  (6c)
  -0.43000000000000  -0.21500000000000   0.25000000000000   H2O  (6c)
  -0.21500000000000   0.21500000000000   0.75000000000000   H2O  (6c)
  -0.21500000000000  -0.43000000000000   0.75000000000000   H2O  (6c)
   0.43000000000000   0.21500000000000   0.75000000000000   H2O  (6c)
   0.33333333333333   0.66666666666667   0.25000000000000   Nd   (2b)
   0.66666666666667   0.33333333333333   0.75000000000000   Nd   (2b)
   0.10500000000000  -0.10500000000000   0.53000000000000   O    (6c)
   0.10500000000000   0.21000000000000   0.53000000000000   O    (6c)
  -0.21000000000000  -0.10500000000000   0.53000000000000   O    (6c)
  -0.10500000000000   0.10500000000000   1.03000000000000   O    (6c)
  -0.10500000000000  -0.21000000000000   1.03000000000000   O    (6c)
   0.21000000000000   0.10500000000000   1.03000000000000   O    (6c)
   0.06500000000000   0.36500000000000   0.75000000000000   O    (12d)
  -0.36500000000000  -0.30000000000000   0.75000000000000   O    (12d)
   0.30000000000000  -0.06500000000000   0.75000000000000   O    (12d)
  -0.06500000000000  -0.36500000000000   1.25000000000000   O    (12d)
   0.36500000000000   0.30000000000000   1.25000000000000   O    (12d)
  -0.30000000000000   0.06500000000000   1.25000000000000   O    (12d)
  -0.36500000000000  -0.06500000000000   0.75000000000000   O    (12d)
   0.30000000000000   0.36500000000000   0.75000000000000   O    (12d)
   0.06500000000000  -0.30000000000000   0.75000000000000   O    (12d)
   0.36500000000000   0.06500000000000   1.25000000000000   O    (12d)
  -0.30000000000000  -0.36500000000000   1.25000000000000   O    (12d)
  -0.06500000000000   0.30000000000000   1.25000000000000   O    (12d)
```

Swedenborgite ($NaBe_4SbO_7$, $E9_2$): A4BC7D_hP26_186_ac_b_a2c_b - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Swedenborgite'
_chemical_formula_sum 'Be4 Na O7 Sb'

loop_
_publ_author_name
 'D. M. C. Huminicki'
 'F. C. Hawthorne'
_journal_name_full_name
;
 Canadian Mineralogist
;
_journal_volume 39
_journal_year 2001
_journal_page_first 153
_journal_page_last 158
_publ_Section_title
;
 Refinement of the Crystal Structure of Swedenborgite
;

_aflow_title 'Swedenborgite (NaBe$_{4}$SbO$_{7}$, $E9_{2}$) Structure'
_aflow_proto 'A4BC7D_hP26_186_ac_b_a2c_b'
_aflow_params 'a,c/a,z_{1},z_{2},z_{3},z_{4},x_{5},z_{5},x_{6},z_{6},x_{
      ↪ 7},z_{7}'
_aflow_params_values '5.4317,1.63063129407,0.0629,0.3728,0.6245,0.0,
      ↪ 0.1664,0.3126,0.4961,0.3706,0.1616,0.1269'
_aflow_Strukturbericht '$E9_{2}$'
_aflow_Pearson 'hP26'

_symmetry_space_group_name_H-M "P 63 m c"
_symmetry_Int_Tables_number 186

_cell_length_a     5.43170
_cell_length_b     5.43170
_cell_length_c     8.85710
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z+1/2
3 -y,x-y,z
4 -x,-y,z+1/2
5 -x+y,-x,z
6 y,-x+y,z+1/2
7 -x+y,y,z
8 -x,-x+y,z+1/2
9 -y,-x,z
10 x-y,-y,z+1/2
11 x,x-y,z
12 y,x,z+1/2

loop_
```

```
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Be1 Be 2 a 0.00000 0.00000 0.06290 1.00000
O1 O 2 a 0.00000 0.00000 0.37280 1.00000
Na1 Na 2 b 0.33333 0.66667 0.62450 1.00000
Sb1 Sb 2 b 0.33333 0.66667 0.00000 1.00000
Be2 Be 6 c 0.16640 0.83360 0.31260 1.00000
O2 O 6 c 0.49610 0.50390 0.37060 1.00000
O3 O 6 c 0.16160 0.83840 0.12690 1.00000
```

Swedenborgite ($NaBe_4SbO_7$, $E9_2$): A4BC7D_hP26_186_ac_b_a2c_b - POSCAR

```
A4BC7D_hP26_186_ac_b_a2c_b & a,c/a,z1,z2,z3,z4,x5,z5,x6,z6,x7,z7 --
    ↪ params=5.4317,1.63063129407,0.0629,0.3728,0.6245,0.0,0.1664,
    ↪ 0.3126,0.4961,0.3706,0.1616,0.1269 & P6_{3}mc C_{6v}^{4} #186 (
    ↪ a^2b^2c^3) & hP26 & $E9_{2}$ & Be4NaO7Sb & Swedenborgite & D.
    ↪ M. C. Huminicki and F. C. Hawthorne, Can. Mineral. 39, 153-158
    ↪ (2001)
   1.00000000000000
   2.71585000000000  -4.70399018573594   0.00000000000000
   2.71585000000000   4.70399018573594   0.00000000000000
   0.00000000000000   0.00000000000000   8.85710000000000
    Be    Na    O    Sb
     8     2    14     2
Direct
   0.00000000000000   0.00000000000000   0.06290000000000  Be   (2a)
   0.00000000000000   0.00000000000000   0.56290000000000  Be   (2a)
   0.16640000000000  -0.16640000000000   0.31260000000000  Be   (6c)
   0.16640000000000   0.33280000000000   0.31260000000000  Be   (6c)
  -0.33280000000000  -0.16640000000000   0.31260000000000  Be   (6c)
  -0.16640000000000   0.16640000000000   0.81260000000000  Be   (6c)
  -0.16640000000000  -0.33280000000000   0.81260000000000  Be   (6c)
   0.33280000000000   0.16640000000000   0.81260000000000  Be   (6c)
   0.33333333333333   0.66666666666667   0.62450000000000  Na   (2b)
   0.66666666666667   0.33333333333333   1.12450000000000  Na   (2b)
   0.00000000000000   0.00000000000000   0.37280000000000  O    (2a)
   0.00000000000000   0.00000000000000   0.87280000000000  O    (2a)
   0.49610000000000  -0.49610000000000   0.37060000000000  O    (6c)
   0.49610000000000   0.99220000000000   0.37060000000000  O    (6c)
  -0.99220000000000  -0.49610000000000   0.37060000000000  O    (6c)
  -0.49610000000000   0.49610000000000   0.87060000000000  O    (6c)
  -0.49610000000000  -0.99220000000000   0.87060000000000  O    (6c)
   0.99220000000000   0.49610000000000   0.87060000000000  O    (6c)
   0.16160000000000  -0.16160000000000   0.12690000000000  O    (6c)
   0.16160000000000   0.32320000000000   0.12690000000000  O    (6c)
  -0.32320000000000  -0.16160000000000   0.12690000000000  O    (6c)
  -0.16160000000000   0.16160000000000   0.62690000000000  O    (6c)
  -0.16160000000000  -0.32320000000000   0.62690000000000  O    (6c)
   0.32320000000000   0.16160000000000   0.62690000000000  O    (6c)
   0.33333333333333   0.66666666666667   0.00000000000000  Sb   (2b)
   0.66666666666667   0.33333333333333   0.50000000000000  Sb   (2b)
```

$C27$ (CdI_2) (*questionable*): AB2_hP6_186_b_ab - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'CdI2'
_chemical_formula_sum 'Cd I2'

loop_
_publ_author_name
 'O. Hassel'
_journal_name_full_name
;
 Zeitschrift f\"{u}r Physikalische Chemie
;
_journal_volume 22B
_journal_year 1933
_journal_page_first 333
_journal_page_last 334
_publ_Section_title
;
 Zur Kristallstruktur des Cadmiumjodids CdJ$_{2}$
;

# Found in Strukturbericht Band III 1933-1935, 1937

_aflow_title '$C27$ (CdI$_{2}$) ({\em{questionable}}) Structure'
_aflow_proto 'AB2_hP6_186_b_ab'
_aflow_params 'a,c/a,z_{1},z_{2},z_{3}'
_aflow_params_values '4.24,3.22405660377,0.375,0.0,0.625'
_aflow_Strukturbericht '$C27$'
_aflow_Pearson 'hP6'

_symmetry_space_group_name_H-M "P 63 m c"
_symmetry_Int_Tables_number 186

_cell_length_a    4.24000
_cell_length_b    4.24000
_cell_length_c    13.67000
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z+1/2
3 -y,x-y,z
4 -x,-y,z+1/2
5 -x+y,-x,z
6 y,-x+y,z+1/2
7 -x+y,y,z
8 -x,-x+y,z+1/2
9 -y,-x,z
10 x-y,-y,z+1/2
11 x,x-y,z
12 y,x,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
I1 I 2 a 0.00000 0.00000 0.37500 1.00000
Cd1 Cd 2 b 0.33333 0.66667 0.00000 1.00000
I2 I 2 b 0.33333 0.66667 0.62500 1.00000
```

$C27$ (CdI_2) (*questionable*): AB2_hP6_186_b_ab - POSCAR

```
AB2_hP6_186_b_ab & a,c/a,z1,z2,z3 --params=4.24,3.22405660377,0.375,0.0,
    ↪ 0.625 & P6_{3}mc C_{6v}^{4} #186 (ab^2) & hP6 & $C27$ & CdI2 &
    ↪ CdI2 & O. Hassel, Z. Phys. Chem. 22B, 333-334 (1933)
   1.00000000000000
   2.12000000000000  -3.67194771204602   0.00000000000000
   2.12000000000000   3.67194771204602   0.00000000000000
   0.00000000000000   0.00000000000000  13.67000000000000
    Cd    I
     2    4
Direct
   0.33333333333333   0.66666666666667   0.00000000000000  Cd   (2b)
   0.66666666666667   0.33333333333333   0.50000000000000  Cd   (2b)
   0.00000000000000   0.00000000000000   0.37500000000000  I    (2a)
   0.00000000000000   0.00000000000000   0.87500000000000  I    (2a)
   0.33333333333333   0.66666666666667   0.62500000000000  I    (2b)
   0.66666666666667   0.33333333333333   1.12500000000000  I    (2b)
```

$LiClO_4{\cdot}3H_2O$ ($H4_{18}$): AB6CD7_hP30_186_b_d_a_b2c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'ClH6LiO7'
_chemical_formula_sum 'Cl H6 Li O7'

loop_
_publ_author_name
 'J.-O. Lundgren'
 'R. Liminga'
 'R. Tellgren'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 38
_journal_year 1982
_journal_page_first 15
_journal_page_last 20
_publ_Section_title
;
 Neutron diffraction refinement of pyroelectric lithium perchlorate
    ↪ trihydrate
;

# Found in The OH stretching frequency in LiClO$_{4}$$\cdot$3H$_{2}$O(s)
    ↪ from ab initio and model potential calculations, 1992

_aflow_title 'LiClO$_{4}$$\cdot$3H$_{2}$O ($H4_{18}$) Structure'
_aflow_proto 'AB6CD7_hP30_186_b_d_a_b2c'
_aflow_params 'a,c/a,z_{1},z_{2},z_{3},x_{4},z_{4},x_{5},z_{5},x_{6},y_{
    ↪ 6},z_{6}'
_aflow_params_values '7.7192,0.706433309151,0.27671,0.5,0.2381,0.56534,
    ↪ 0.08902,0.12232,0.02787,-0.06662,0.26326,0.53382'
_aflow_Strukturbericht '$H4_{18}$'
_aflow_Pearson 'hP30'

_symmetry_space_group_name_H-M "P 63 m c"
_symmetry_Int_Tables_number 186

_cell_length_a    7.71920
_cell_length_b    7.71920
_cell_length_c    5.45310
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z+1/2
3 -y,x-y,z
4 -x,-y,z+1/2
5 -x+y,-x,z
6 y,-x+y,z+1/2
7 -x+y,y,z
```

```
8 -x,-x+y,z+1/2
9 -y,-x,z
10 x-y,-y,z+1/2
11 x,x-y,z
12 y,x,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Li1 Li 2 a 0.00000 0.00000 0.27671 1.00000
Cl1 Cl 2 b 0.33333 0.66667 0.50000 1.00000
O1 O 2 b 0.33333 0.66667 0.23810 1.00000
O2 O 6 c 0.56534 0.43466 0.08902 1.00000
O3 O 6 c 0.12232 0.87768 0.02787 1.00000
H1 H 12 d -0.06662 0.26326 0.53382 1.00000
```

$LiClO_4{\cdot}3H_2O$ ($H4_{18}$): AB6CD7_hP30_186_b_d_a_b2c - POSCAR

```
AB6CD7_hP30_186_b_d_a_b2c & a,c/a,z1,z2,z3,x4,z4,x5,z5,x6,y6,z6 --params
    ↪ =7.7192,0.706433309151,0.27671,0.5,0.2381,0.56534,0.08902,
    ↪ 0.12232,0.02787,-0.06662,0.26326,0.53382 & P6_{3}mc C_{6v}^{4}
    ↪ #186 (ab^2c^2d) & hP30 & $H4_{18}$ & ClH6LiO7 & ClH6LiO7 &
    ↪ J.-O. Lundgren and R. Liminga and R. Tellgren, Acta
    ↪ Crystallogr. Sect. B Struct. Sci. 38, 15-20 (1982)
   1.00000000000000
   3.85960000000000  -6.68502329689284   0.00000000000000
   3.85960000000000   6.68502329689284   0.00000000000000
   0.00000000000000   0.00000000000000   5.45310000000000
    Cl    H    Li    O
     2   12     2   14
Direct
   0.33333333333333   0.66666666666667   0.50000000000000   Cl   (2b)
   0.66666666666667   0.33333333333333   1.00000000000000   Cl   (2b)
  -0.06662000000000   0.26326000000000   0.53382000000000   H   (12d)
  -0.26326000000000  -0.32988000000000   0.53382000000000   H   (12d)
   0.32988000000000   0.06662000000000   0.53382000000000   H   (12d)
   0.06662000000000  -0.26326000000000   1.03382000000000   H   (12d)
   0.26326000000000   0.32988000000000   1.03382000000000   H   (12d)
  -0.32988000000000  -0.06662000000000   1.03382000000000   H   (12d)
  -0.26326000000000   0.06662000000000   0.53382000000000   H   (12d)
   0.32988000000000   0.26326000000000   0.53382000000000   H   (12d)
  -0.06662000000000  -0.32988000000000   0.53382000000000   H   (12d)
   0.26326000000000  -0.06662000000000   1.03382000000000   H   (12d)
  -0.32988000000000  -0.26326000000000   1.03382000000000   H   (12d)
   0.06662000000000   0.32988000000000   1.03382000000000   H   (12d)
   0.00000000000000   0.00000000000000   0.27671000000000   Li   (2a)
   0.00000000000000   0.00000000000000   0.77671000000000   Li   (2a)
   0.33333333333333   0.66666666666667   0.23810000000000   O   (2b)
   0.66666666666667   0.33333333333333   0.73810000000000   O   (2b)
   0.56534000000000  -0.56534000000000   0.08902000000000   O   (6c)
   0.56534000000000   1.13068000000000   0.08902000000000   O   (6c)
  -1.13068000000000  -0.56534000000000   0.08902000000000   O   (6c)
  -0.56534000000000   0.56534000000000   0.58902000000000   O   (6c)
  -0.56534000000000  -1.13068000000000   0.58902000000000   O   (6c)
   1.13068000000000   0.56534000000000   0.58902000000000   O   (6c)
   0.12232000000000  -0.12232000000000   0.02787000000000   O   (6c)
   0.12232000000000   0.24464000000000   0.02787000000000   O   (6c)
  -0.24464000000000  -0.12232000000000   0.02787000000000   O   (6c)
  -0.12232000000000   0.12232000000000   0.52787000000000   O   (6c)
  -0.12232000000000  -0.24464000000000   0.52787000000000   O   (6c)
   0.24464000000000   0.12232000000000   0.52787000000000   O   (6c)
```

Cd(OH)Cl ($E0_3$): ABCD_hP8_186_b_b_a_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'CdClHO'
_chemical_formula_sum 'Cd Cl H O'

loop_
_publ_author_name
 'Y. Cudennec'
 'A. Riou'
 'Y. G\\'{e}rault'
 'A. Lecerf'
_journal_name_full_name
;
 Journal of Solid State Chemistry
;
_journal_volume 151
_journal_year 2000
_journal_page_first 308
_journal_page_last 312
_publ_Section_title
;
 Synthesis and Crystal Structures of Cd(OH)Cl and Cu(OH)Cl and
    ↪ Relationship to Brucite Type
;

_aflow_title 'Cd(OH)Cl ($E0_{3}$) Structure'
_aflow_proto 'ABCD_hP8_186_b_b_a_a'
_aflow_params 'a,c/a,z_{1},z_{2},z_{3},z_{4}'
_aflow_params_values '3.6648,2.79155752019,0.18,0.0892,0.0079,0.3433'
_aflow_Strukturbericht '$E0_{3}$'
_aflow_Pearson 'hP8'

_symmetry_space_group_name_H-M "P 63 m c"
_symmetry_Int_Tables_number 186

_cell_length_a    3.66480
_cell_length_b    3.66480
_cell_length_c    10.23050
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z+1/2
3 -y,x-y,z
4 -x,-y,z+1/2
5 -x+y,-x,z
6 y,-x+y,z+1/2
7 -x+y,y,z
8 -x,-x+y,z+1/2
9 -y,-x,z
10 x-y,-y,z+1/2
11 x,x-y,z
12 y,x,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
H1 H 2 a 0.00000 0.00000 0.18000 1.00000
O1 O 2 a 0.00000 0.00000 0.08920 1.00000
Cd1 Cd 2 b 0.33333 0.66667 0.00790 1.00000
Cl1 Cl 2 b 0.33333 0.66667 0.34330 1.00000
```

Cd(OH)Cl ($E0_3$): ABCD_hP8_186_b_b_a_a - POSCAR

```
ABCD_hP8_186_b_b_a_a & a,c/a,z1,z2,z3,z4 --params=3.6648,2.79155752019,
    ↪ 0.18,0.0892,0.0079,0.3433 & P6_{3}mc C_{6v}^{4} #186 (a^2b^2) &
    ↪ hP8 & $E0_{3}$ & CdClHO & CdClHO & Y. Cudennec et al., J.
    ↪ Solid State Chem. 151, 308-312 (2000)
   1.00000000000000
   1.83240000000000  -3.17380989978921   0.00000000000000
   1.83240000000000   3.17380989978921   0.00000000000000
   0.00000000000000   0.00000000000000  10.23050000000000
    Cd    Cl    H    O
     2     2    2    2
Direct
   0.33333333333333   0.66666666666667   0.00790000000000   Cd   (2b)
   0.66666666666667   0.33333333333333   0.50790000000000   Cd   (2b)
   0.33333333333333   0.66666666666667   0.34330000000000   Cl   (2b)
   0.66666666666667   0.33333333333333   0.84330000000000   Cl   (2b)
   0.00000000000000   0.00000000000000   0.18000000000000   H   (2a)
   0.00000000000000   0.00000000000000   0.68000000000000   H   (2a)
   0.00000000000000   0.00000000000000   0.08920000000000   O   (2a)
   0.00000000000000   0.00000000000000   0.58920000000000   O   (2a)
```

Cr-233 Quasi-One-Dimensional Superconductor ($K_2Cr_3As_3$): A3B3C2_hP16_187_jk_jk_ck - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'As3Cr3K2'
_chemical_formula_sum 'As3 Cr3 K2'

loop_
_publ_author_name
 'J.-K. Bao'
 'J.-Y. Liu'
 'C.-W. Ma'
 'Z.-H. Meng'
 'Z.-T. Tang'
 'Y.-L. Sun'
 'H.-F. Zhai'
 'H. Jiang'
 'H. Bai'
 'C.-M. Feng'
 'Z.-A. Xu'
 'G.-H. Cao'
_journal_name_full_name
;
 Physical Review X
;
_journal_volume 5
_journal_year 2015
_journal_page_first 011013
_journal_page_last 011013
_publ_Section_title
;
 Superconductivity in Quasi-One-Dimensional K$_{2}$Cr$_{3}$As$_{3}$ with
    ↪ Significant Electron Correlations
;

# Found in Q.-G. et al., Superconductivity at 10.4 K in a novel
    ↪ quasi-one-dimensional ternary molybdenum pnictide K$_{2}$Mo$_{3
    ↪ }$As$_{3}$, Sci. Bull. 63, 952-956 (2018)

_aflow_title 'Cr-233 Quasi-One-Dimensional Superconductor (K$_{2}$Cr$_{3
    ↪ }$As$_{3}$) Structure'
_aflow_proto 'A3B3C2_hP16_187_jk_jk_ck'
```

```
_aflow_params 'a,c/a,x_{2},x_{3},x_{4},x_{5},x_{6}'
_aflow_params_values '9.9832,0.423751903197,0.8339,0.0898,0.1676,-0.0873
    ↪ ,0.5387'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hP16'

_symmetry_space_group_name_H-M "P -6 m 2"
_symmetry_Int_Tables_number 187

_cell_length_a 9.98320
_cell_length_b 9.98320
_cell_length_c 4.23040
_cell_angle_alpha 90.00000
_cell_angle_beta 90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 x,x-y,-z
5 -x+y,y,-z
6 -y,-x,-z
7 -x+y,-x,-z
8 x,y,-z
9 -y,x-y,-z
10 -x+y,y,z
11 -y,-x,z
12 x,x-y,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
K1 K 1 c 0.33333 0.66667 0.00000 1.00000
As1 As 3 j 0.83390 0.16610 0.00000 1.00000
Cr1 Cr 3 j 0.08980 -0.08980 0.00000 1.00000
As2 As 3 k 0.16760 0.83240 0.50000 1.00000
Cr2 Cr 3 k -0.08730 0.08730 0.50000 1.00000
K2 K 3 k 0.53870 0.46130 0.50000 1.00000
```

Cr-233 Quasi-One-Dimensional Superconductor ($K_2Cr_3As_3$): A3B3C2_hP16_187_jk_jk_ck - POSCAR

```
A3B3C2_hP16_187_jk_jk_ck & a,c/a,x2,x3,x4,x5,x6 --params=9.9832,
    ↪ 0.423751903197,0.8339,0.0898,0.1676,-0.0873,0.5387 & P-6m2 D_{
    ↪ 3h}^{1} #187 (cj^2k^3) & hP16 & None & As3Cr3K2 & As3Cr3K2 &
    ↪ J.-K. Bao et al., Phys. Rev. X 5, 011013(2015)
   1.00000000000000
   4.99160000000000  -8.64570481106081   0.00000000000000
   4.99160000000000   8.64570481106081   0.00000000000000
   0.00000000000000   0.00000000000000   4.23040000000000
    As  Cr  K
     6   6   4
Direct
   0.83390000000000  -0.83390000000000   0.00000000000000  As  (3j)
   0.83390000000000   1.66780000000000   0.00000000000000  As  (3j)
  -1.66780000000000  -0.83390000000000   0.00000000000000  As  (3j)
   0.16760000000000  -0.16760000000000   0.50000000000000  As  (3k)
   0.16760000000000   0.33520000000000   0.50000000000000  As  (3k)
  -0.33520000000000  -0.16760000000000   0.50000000000000  As  (3k)
   0.08980000000000  -0.08980000000000   0.00000000000000  Cr  (3j)
   0.08980000000000   0.17960000000000   0.00000000000000  Cr  (3j)
  -0.17960000000000  -0.08980000000000   0.00000000000000  Cr  (3j)
  -0.08730000000000   0.08730000000000   0.50000000000000  Cr  (3k)
  -0.08730000000000  -0.17460000000000   0.50000000000000  Cr  (3k)
   0.17460000000000   0.08730000000000   0.50000000000000  Cr  (3k)
   0.33333333333333   0.66666666666667   0.00000000000000  K  (1c)
   0.53870000000000  -0.53870000000000   0.50000000000000  K  (3k)
   0.53870000000000   1.07740000000000   0.50000000000000  K  (3k)
  -1.07740000000000  -0.53870000000000   0.50000000000000  K  (3k)
```

Cs_7O: A7B_hP24_187_ai2j2kn_j - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Cs7O'
_chemical_formula_sum 'Cs7 O'

loop_
_publ_author_name
 'A. Simon'
_journal_name_full_name
;
 Zeitschrift fur Anorganische und Allgemeine Chemie
;
_journal_volume 422
_journal_year 1976
_journal_page_first 208
_journal_page_last 218
_publ_Section_title
;
 \"{U}ber Alkalimetall-Suboxide. VII. Das metallreichste C\"{a}
    ↪ siumoxid-Cs$_{7}$O
;
```

```
# Found in Binary Alloy Phase Diagrams, 1990 Found in Binary Alloy Phase
    ↪ Diagrams, {Cd-Ce to Hf-Rb}}

_aflow_title 'Cs$_{7}$O Structure'
_aflow_proto 'A7B_hP24_187_ai2j2kn_j'
_aflow_params 'a,c/a,z_{2},x_{3},x_{4},x_{5},x_{6},x_{7},x_{8},z_{8}'
_aflow_params_values '16.244,0.562977099237,0.20321,0.55085,0.21582,
    ↪ 0.74805,0.44928,0.11229,0.81357,0.21992'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hP24'

_symmetry_space_group_name_H-M "P -6 m 2"
_symmetry_Int_Tables_number 187

_cell_length_a 16.24400
_cell_length_b 16.24400
_cell_length_c 9.14500
_cell_angle_alpha 90.00000
_cell_angle_beta 90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 x,x-y,-z
5 -x+y,y,-z
6 -y,-x,-z
7 -x+y,-x,-z
8 x,y,-z
9 -y,x-y,-z
10 -x+y,y,z
11 -y,-x,z
12 x,x-y,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cs1 Cs 1 a 0.00000 0.00000 0.00000 1.00000
Cs2 Cs 2 i 0.66667 0.33333 0.20321 1.00000
Cs3 Cs 3 j 0.55085 0.44915 0.00000 1.00000
Cs4 Cs 3 j 0.21582 0.78418 0.00000 1.00000
O1 O 3 j 0.74805 0.25195 0.00000 1.00000
Cs5 Cs 3 k 0.44928 0.55072 0.50000 1.00000
Cs6 Cs 3 k 0.11229 0.88771 0.50000 1.00000
Cs7 Cs 6 n 0.81357 0.18643 0.21992 1.00000
```

Cs_7O: A7B_hP24_187_ai2j2kn_j - POSCAR

```
A7B_hP24_187_ai2j2kn_j & a,c/a,z2,x3,x4,x5,x6,x7,x8,z8 --params=16.244,
    ↪ 0.562977099237,0.20321,0.55085,0.21582,0.74805,0.44928,0.11229,
    ↪ 0.81357,0.21992 & P-6m2 D_{3h}^{1} #187 (aij^3k^2n) & hP24 &
    ↪ None & Cs7O & Cs7O & A. Simon, Z. Anorg. Allg. Chem. 422,
    ↪ 208-218 (1976)
   1.00000000000000
   8.12200000000000  -14.06771665907440   0.00000000000000
   8.12200000000000   14.06771665907440   0.00000000000000
   0.00000000000000   0.00000000000000   9.14500000000000
    Cs  O
    21  3
Direct
   0.00000000000000   0.00000000000000   0.00000000000000  Cs  (1a)
   0.66666666666667   0.33333333333333   0.20321000000000  Cs  (2i)
   0.66666666666667   0.33333333333333  -0.20321000000000  Cs  (2i)
   0.55085000000000  -0.55085000000000   0.00000000000000  Cs  (3j)
   0.55085000000000   1.10170000000000   0.00000000000000  Cs  (3j)
  -1.10170000000000  -0.55085000000000   0.00000000000000  Cs  (3j)
   0.21582000000000  -0.21582000000000   0.00000000000000  Cs  (3j)
   0.21582000000000   0.43164000000000   0.00000000000000  Cs  (3j)
  -0.43164000000000  -0.21582000000000   0.00000000000000  Cs  (3j)
   0.44928000000000  -0.44928000000000   0.50000000000000  Cs  (3k)
   0.44928000000000   0.89856000000000   0.50000000000000  Cs  (3k)
  -0.89856000000000  -0.44928000000000   0.50000000000000  Cs  (3k)
   0.11229000000000  -0.11229000000000   0.50000000000000  Cs  (3k)
   0.11229000000000   0.22458000000000   0.50000000000000  Cs  (3k)
  -0.22458000000000  -0.11229000000000   0.50000000000000  Cs  (3k)
   0.81357000000000  -0.81357000000000   0.21992000000000  Cs  (6n)
   0.81357000000000   1.62714000000000   0.21992000000000  Cs  (6n)
  -1.62714000000000  -0.81357000000000   0.21992000000000  Cs  (6n)
   0.81357000000000  -0.81357000000000  -0.21992000000000  Cs  (6n)
   0.81357000000000   1.62714000000000  -0.21992000000000  Cs  (6n)
  -1.62714000000000  -0.81357000000000  -0.21992000000000  Cs  (6n)
   0.74805000000000  -0.74805000000000   0.00000000000000  O  (3j)
   0.74805000000000   1.49610000000000   0.00000000000000  O  (3j)
  -1.49610000000000  -0.74805000000000   0.00000000000000  O  (3j)
```

ZrNiAl: ABC_hP9_189_g_ad_f - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'AlNiZr'
_chemical_formula_sum 'Al Ni Zr'

loop_
_publ_author_name
```

```
 'O. Shved'
 'L. P. Salamakha'
 'S. Mudry'
 'O. Sologub'
 'P. F. Rogl'
 'E. Bauer'
_journal_name_full_name
;
 Journal of Alloys and Compounds
;
_journal_volume 821
_journal_year 2020
_journal_page_first 153326
_journal_page_last 153326
_publ_Section_title
;
 Zr-based nickel aluminides: crystal structure and electronic properties
;

_aflow_title 'ZrNiAl Structure'
_aflow_proto 'ABC_hP9_189_g_ad_f'
_aflow_params 'a,c/a,x_{3},x_{4}'
_aflow_params_values '6.91558,0.50180028284,0.59354,0.2491'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hP9'

_symmetry_space_group_name_H-M "P -6 2 m"
_symmetry_Int_Tables_number 189

_cell_length_a    6.91558
_cell_length_b    6.91558
_cell_length_c    3.47024
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 x-y,-y,-z
5 y,x,-z
6 -x,-x+y,-z
7 -x+y,-x,-z
8 x,y,-z
9 -y,x-y,-z
10 -x,-x+y,z
11 x-y,-y,z
12 y,x,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ni1 Ni 1 a 0.00000 0.00000 0.00000 1.00000
Ni2 Ni 2 d 0.33333 0.66667 0.50000 1.00000
Zr1 Zr 3 f 0.59354 0.00000 0.00000 1.00000
Al1 Al 3 g 0.24910 0.00000 0.50000 1.00000
```

ZrNiAl: ABC_hP9_189_g_ad_f - POSCAR

```
ABC_hP9_189_g_ad_f & a,c/a,x3,x4 --params=6.91558,0.50180028284,0.59354,
      ↪ 0.2491 & P-62m D_{3h}^{3} #189 (adfg) & hP9 & None & AlNiZr &
      ↪ AlNiZr & O. Shved et al., J. Alloys Compd. 821, 153326(2020)
   1.00000000000000
   3.45779000000000  -5.98906796190359   0.00000000000000
   3.45779000000000   5.98906796190359   0.00000000000000
   0.00000000000000   0.00000000000000   3.47024000000000
     Al    Ni    Zr
      3     3     3
Direct
   0.24910000000000   0.00000000000000   0.50000000000000   Al   (3g)
   0.00000000000000   0.24910000000000   0.50000000000000   Al   (3g)
  -0.24910000000000  -0.24910000000000   0.50000000000000   Al   (3g)
   0.00000000000000   0.00000000000000   0.00000000000000   Ni   (1a)
   0.33333333333333   0.66666666666667   0.50000000000000   Ni   (2d)
   0.66666666666667   0.33333333333333   0.50000000000000   Ni   (2d)
   0.59354000000000   0.00000000000000   0.00000000000000   Zr   (3f)
   0.00000000000000   0.59354000000000   0.00000000000000   Zr   (3f)
  -0.59354000000000  -0.59354000000000   0.00000000000000   Zr   (3f)
```

$CsSO_3$ ($K1_2$): AB3C_hP20_190_ac_i_f - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'CsO3S'
_chemical_formula_sum 'Cs O3 S'

loop_
_publ_author_name
 'G. H\"{a}gg'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Physikalische Chemie B
;
_journal_volume 18
_journal_year 1932
_journal_page_first 327
_journal_page_last 342
_publ_Section_title
;
 Die Kristallstruktur von Caesiumdithionat, Cs$_{2}$S$_{2}$O$_{6}$
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'CsSO$_{3}$ ($K1_{2}$) Structure'
_aflow_proto 'AB3C_hP20_190_ac_i_f'
_aflow_params 'a,c/a,z_{3},x_{4},y_{4},z_{4}'
_aflow_params_values '6.326,1.82342712615,0.66,0.44,0.33333,0.125'
_aflow_Strukturbericht '$K1_{2}$'
_aflow_Pearson 'hP20'

_symmetry_space_group_name_H-M "P -6 2 c"
_symmetry_Int_Tables_number 190

_cell_length_a    6.32600
_cell_length_b    6.32600
_cell_length_c    11.53500
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 x-y,-y,-z
5 y,x,-z
6 -x,-x+y,-z
7 -x+y,-x,-z+1/2
8 x,y,-z+1/2
9 -y,x-y,-z+1/2
10 -x,-x+y,z+1/2
11 x-y,-y,z+1/2
12 y,x,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cs1 Cs 2 a 0.00000 0.00000 0.00000 1.00000
Cs2 Cs 2 c 0.33333 0.66667 0.25000 1.00000
S1 S 4 f 0.33333 0.66667 0.66000 1.00000
O1 O 12 i 0.44000 0.33333 0.12500 1.00000
```

$CsSO_3$ ($K1_2$): AB3C_hP20_190_ac_i_f - POSCAR

```
AB3C_hP20_190_ac_i_f & a,c/a,z3,x4,y4,z4 --params=6.326,1.82342712615,
      ↪ 0.66,0.44,0.33333,0.125 & P-62c D_{3h}^{4} #190 (acfi) & hP20 &
      ↪ $K1_{2}$ & CsO3S & CsO3S & G. H\"{a}gg, Z. Physik. Chem. B 18,
      ↪ 327-342 (1932)
   1.00000000000000
   3.16300000000000  -5.47847670434036   0.00000000000000
   3.16300000000000   5.47847670434036   0.00000000000000
   0.00000000000000   0.00000000000000  11.53500000000000
     Cs    O    S
      4   12    4
Direct
   0.00000000000000   0.00000000000000   0.00000000000000   Cs   (2a)
   0.00000000000000   0.00000000000000   0.50000000000000   Cs   (2a)
   0.33333333333333   0.66666666666667   0.25000000000000   Cs   (2c)
   0.66666666666667   0.33333333333333   0.75000000000000   Cs   (2c)
   0.44000000000000   0.33333000000000   0.12500000000000   O   (12i)
  -0.33333000000000   0.10667000000000   0.12500000000000   O   (12i)
  -0.10667000000000  -0.44000000000000   0.12500000000000   O   (12i)
   0.44000000000000   0.33333000000000   0.37500000000000   O   (12i)
  -0.33333000000000   0.10667000000000   0.37500000000000   O   (12i)
  -0.10667000000000  -0.44000000000000   0.37500000000000   O   (12i)
   0.33333000000000   0.44000000000000  -0.12500000000000   O   (12i)
   0.10667000000000  -0.33333000000000  -0.12500000000000   O   (12i)
  -0.44000000000000  -0.10667000000000  -0.12500000000000   O   (12i)
   0.33333000000000   0.44000000000000   0.62500000000000   O   (12i)
   0.10667000000000  -0.33333000000000   0.62500000000000   O   (12i)
  -0.44000000000000  -0.10667000000000   0.62500000000000   O   (12i)
   0.33333333333333   0.66666666666667   0.66000000000000   S    (4f)
   0.33333333333333   0.66666666666667  -0.16000000000000   S    (4f)
   0.66666666666667   0.33333333333333  -0.66000000000000   S    (4f)
   0.66666666666667   0.33333333333333   1.16000000000000   S    (4f)
```

Bastnäsite [CeF(CO_3)]: ABCD3_hP36_190_h_g_af_hi - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Bastn\"{a}site'
_chemical_formula_sum 'C Ce F O3'

loop_
_publ_author_name
 'Y. Ni'
 'J. M. Hughes'
 'A. N. Mariano'
```

```
_journal_name_full_name
;
 American Mineralogist
;
_journal_volume 78
_journal_year 1993
_journal_page_first 415
_journal_page_last 418
_publ_Section_title
;
 The atomic arrangement of bastn\"{a}site-(Ce), Ce(CO$_{3}$)F, and
      ↪ structural elements of synchysite-(Ce), r{\"o}ntgenite-(Ce),
      ↪ and parisite-(Ce)
;

_aflow_title 'Bastn\"{a}site [CeF(CO$_{3}$)] Structure'
_aflow_proto 'ABCD3_hP36_190_h_g_af_hi'
_aflow_params 'a,c/a,z_{2},x_{3},x_{4},y_{4},x_{5},y_{5},x_{6},y_{6},z_{
      ↪ 6}'
_aflow_params_values '7.1175,1.37153494907,0.449,0.33941,0.032,0.71,
      ↪ 0.207,0.891,0.3245,0.3828,0.6354'
_aflow_Strukturbericht '$G7_{1}$'
_aflow_Pearson 'hP36'

_symmetry_space_group_name_H-M "P -6 2 c"
_symmetry_Int_Tables_number 190

_cell_length_a    7.11750
_cell_length_b    7.11750
_cell_length_c    9.76190
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 -y,x-y,z
3 -x+y,-x,z
4 x-y,-y,-z
5 y,x,-z
6 -x,-x+y,-z
7 -x+y,-x,-z+1/2
8 x,y,-z+1/2
9 -y,x-y,-z+1/2
10 -x,-x+y,z+1/2
11 x-y,-y,z+1/2
12 y,x,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
F1 F 2 a 0.00000 0.00000 0.00000 1.00000
F2 F 4 f 0.33333 0.66667 0.44900 1.00000
Ce1 Ce 6 g 0.33941 0.00000 0.00000 1.00000
C1 C 6 h 0.03200 0.71000 0.25000 1.00000
O1 O 6 h 0.20700 0.89100 0.25000 1.00000
O2 O 12 i 0.32450 0.38280 0.63540 1.00000
```

Bastnäsite [CeF(CO_3)]: ABCD3_hP36_190_h_g_af_hi - POSCAR

```
ABCD3_hP36_190_h_g_af_hi & a,c/a,z2,x3,x4,y4,x5,y5,x6,y6,z6 --params=
      ↪ 7.1175,1.37153494907,0.449,0.33941,0.032,0.71,0.207,0.891,
      ↪ 0.3245,0.3828,0.6354 & P-62c D_{3h}^{4} #190 (afgh^2i) & hP36 &
      ↪  $G7_{1}$ & CCeFO3 & Bastn\"{a}site & Y. Ni and J. M. Hughes
      ↪ and A. N. Mariano, Am. Mineral. 78, 415-418 (1993)
   1.00000000000000
   3.55875000000000   -6.16393581143574    0.00000000000000
   3.55875000000000    6.16393581143574    0.00000000000000
   0.00000000000000    0.00000000000000    9.76190000000000
     C     Ce     F     O
     6     6     6     18
Direct
   0.03200000000000    0.71000000000000    0.25000000000000     C    (6h)
  -0.71000000000000   -0.67800000000000    0.25000000000000     C    (6h)
   0.67800000000000   -0.03200000000000    0.25000000000000     C    (6h)
   0.71000000000000    0.03200000000000    0.75000000000000     C    (6h)
  -0.67800000000000   -0.71000000000000    0.75000000000000     C    (6h)
  -0.03200000000000    0.67800000000000    0.75000000000000     C    (6h)
   0.33941000000000    0.00000000000000    0.00000000000000    Ce    (6g)
   0.00000000000000    0.33941000000000    0.00000000000000    Ce    (6g)
  -0.33941000000000   -0.33941000000000    0.00000000000000    Ce    (6g)
   0.33941000000000    0.00000000000000    0.50000000000000    Ce    (6g)
   0.00000000000000    0.33941000000000    0.50000000000000    Ce    (6g)
  -0.33941000000000   -0.33941000000000    0.50000000000000    Ce    (6g)
   0.00000000000000    0.00000000000000    0.00000000000000     F    (2a)
   0.00000000000000    0.00000000000000    0.50000000000000     F    (2a)
   0.33333333333333    0.66666666666667    0.44900000000000     F    (4f)
   0.33333333333333    0.66666666666667    0.05100000000000     F    (4f)
   0.66666666666667    0.33333333333333   -0.44900000000000     F    (4f)
   0.66666666666667    0.33333333333333    0.94900000000000     F    (4f)
   0.20700000000000    0.89100000000000    0.25000000000000     O    (6h)
  -0.89100000000000   -0.68400000000000    0.25000000000000     O    (6h)
   0.68400000000000   -0.20700000000000    0.25000000000000     O    (6h)
   0.89100000000000    0.20700000000000    0.75000000000000     O    (6h)
  -0.68400000000000   -0.89100000000000    0.75000000000000     O    (6h)
  -0.20700000000000    0.68400000000000    0.75000000000000     O    (6h)
   0.32450000000000    0.38280000000000    0.63540000000000     O   (12i)
  -0.38280000000000   -0.05830000000000    0.63540000000000     O   (12i)
   0.05830000000000   -0.32450000000000    0.63540000000000     O   (12i)
   0.32450000000000    0.38280000000000   -0.13540000000000     O   (12i)
  -0.38280000000000   -0.05830000000000   -0.13540000000000     O   (12i)
   0.05830000000000   -0.32450000000000   -0.13540000000000     O   (12i)
   0.38280000000000    0.32450000000000   -0.63540000000000     O   (12i)
  -0.05830000000000   -0.38280000000000   -0.63540000000000     O   (12i)
  -0.32450000000000    0.05830000000000   -0.63540000000000     O   (12i)
   0.38280000000000    0.32450000000000    1.13540000000000     O   (12i)
  -0.05830000000000   -0.38280000000000    1.13540000000000     O   (12i)
  -0.32450000000000    0.05830000000000    1.13540000000000     O   (12i)
```

$TiBe_{12}$ (approximate, $D2_a$): A12B_hP13_191_cdei_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Be12Ti'
_chemical_formula_sum 'Be12 Ti'

loop_
_publ_author_name
 'R. F. Raeuchle'
 'R. E. Rundle'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 5
_journal_year 1952
_journal_page_first 85
_journal_page_last 93
_publ_Section_title
;
 The Structure of TiBe$_{12}$
;

# Found in Resolving the structure of TiBe$_{12}$, 2016

_aflow_title 'TiBe$_{12}$ (approximate, $D2_{a}$) Structure'
_aflow_proto 'A12B_hP13_191_cdei_a'
_aflow_params 'a,c/a,z_{4},z_{5}'
_aflow_params_values '4.23,1.73286052009,0.29,0.25'
_aflow_Strukturbericht '$D2_{a}$'
_aflow_Pearson 'hP13'

_symmetry_space_group_name_H-M "P 6/m 2/m 2/m"
_symmetry_Int_Tables_number 191

_cell_length_a    4.23000
_cell_length_b    4.23000
_cell_length_c    7.33000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z
3 -y,x-y,z
4 -x,-y,z
5 -x+y,-x,z
6 y,-x+y,z
7 x-y,-y,-z
8 x,x-y,-z
9 y,x,-z
10 -x+y,y,-z
11 -x,-x+y,-z
12 -y,-x,-z
13 -x,-y,-z
14 -x+y,-x,-z
15 y,-x+y,-z
16 x,y,-z
17 x-y,x,-z
18 -y,x-y,-z
19 -x+y,y,z
20 -x,-x+y,z
21 -y,-x,z
22 x-y,-y,z
23 x,x-y,z
24 y,x,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ti1 Ti 1 a 0.00000 0.00000 0.00000 1.00000
Be1 Be 2 c 0.33333 0.66667 0.00000 1.00000
Be2 Be 2 d 0.33333 0.66667 0.50000 1.00000
Be3 Be 2 e 0.00000 0.00000 0.29000 1.00000
Be4 Be 6 i 0.50000 0.00000 0.25000 1.00000
```

$TiBe_{12}$ (approximate, $D2_a$): A12B_hP13_191_cdei_a - POSCAR

```
A12B_hP13_191_cdei_a & a,c/a,z4,z5 --params=4.23,1.73286052009,0.29,0.25
      ↪ & P6/mmm D_{6h}^{1} #191 (acdei) & hP13 & $D2_{a}$ & Be12Ti &
```

```
    ↪ Be12Ti & R. F. Raeuchle and R. E. Rundle, Acta Cryst. 5, 85-93
    ↪ (1952)
   1.00000000000000
   2.11500000000000  -3.66328745800818   0.00000000000000
   2.11500000000000   3.66328745800818   0.00000000000000
   0.00000000000000   0.00000000000000   7.33000000000000
    Be    Ti
    12     1
Direct
   0.33333333333333   0.66666666666667   0.00000000000000   Be   (2c)
   0.66666666666667   0.33333333333333   0.00000000000000   Be   (2c)
   0.33333333333333   0.66666666666667   0.50000000000000   Be   (2d)
   0.66666666666667   0.33333333333333   0.50000000000000   Be   (2d)
   0.00000000000000   0.00000000000000   0.29000000000000   Be   (2e)
   0.00000000000000   0.00000000000000  -0.29000000000000   Be   (2e)
   0.50000000000000   0.00000000000000   0.25000000000000   Be   (6i)
   0.00000000000000   0.50000000000000   0.25000000000000   Be   (6i)
   0.50000000000000   0.50000000000000   0.25000000000000   Be   (6i)
   0.00000000000000   0.50000000000000  -0.25000000000000   Be   (6i)
   0.50000000000000   0.00000000000000  -0.25000000000000   Be   (6i)
   0.50000000000000   0.50000000000000  -0.25000000000000   Be   (6i)
   0.00000000000000   0.00000000000000   0.00000000000000   Ti   (1a)
```

Hexagonal WO_3: A3B_hP12_191_gl_f - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'O3W'
_chemical_formula_sum 'O3 W'

loop_
_publ_author_name
 'P. M. Woodward'
 'A. W. Sleight'
 'T. Vogt'
_journal_name_full_name
;
 Journal of Solid State Chemistry
;
_journal_volume 131
_journal_year 1997
_journal_page_first 9
_journal_page_last 17
_publ_Section_title
;
 Ferroelectric Tungsten Trioxide
;

_aflow_title 'Hexagonal WO$_{3}$ Structure'
_aflow_proto 'A3B_hP12_191_gl_f'
_aflow_params 'a,c/a,x_{3}'
_aflow_params_values '7.298,0.534255960537,0.212'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hP12'

_symmetry_space_group_name_H-M "P 6/m 2/m 2/m"
_symmetry_Int_Tables_number 191

_cell_length_a    7.29800
_cell_length_b    7.29800
_cell_length_c    3.89900
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z
3 -y,x-y,z
4 -x,-y,z
5 -x+y,-x,z
6 y,-x+y,z
7 x-y,-y,-z
8 x,x-y,-z
9 y,x,-z
10 -x+y,y,-z
11 -x,-x+y,-z
12 -y,-x,-z
13 -x,-y,-z
14 -x+y,-x,-z
15 y,-x+y,-z
16 x,y,-z
17 x-y,x,-z
18 -y,x-y,-z
19 -x+y,y,z
20 -x,-x+y,z
21 -y,-x,z
22 x-y,-y,z
23 x,x-y,z
24 y,x,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
W1 W 3 f 0.50000 0.00000 0.00000 1.00000
O1 O 3 g 0.50000 0.00000 0.50000 1.00000
O2 O 6 l 0.21200 0.42400 0.00000 1.00000
```

Hexagonal WO_3: A3B_hP12_191_gl_f - POSCAR

```
A3B_hP12_191_gl_f & a,c/a,x3 --params=7.298,0.534255960537,0.212 & P6/
    ↪ mmm D_{6h}^{1} #191 (fgl) & hP12 & None & O3W & O3W & P. M.
    ↪ Woodward and A. W. Sleight and T. Vogt, J. Solid State Chem.
    ↪ 131, 9-17 (1997)
   1.00000000000000
   3.64900000000000  -6.32025339681883   0.00000000000000
   3.64900000000000   6.32025339681883   0.00000000000000
   0.00000000000000   0.00000000000000   3.89900000000000
    O     W
    9     3
Direct
   0.50000000000000   0.00000000000000   0.50000000000000   O   (3g)
   0.00000000000000   0.50000000000000   0.50000000000000   O   (3g)
   0.50000000000000   0.50000000000000   0.50000000000000   O   (3g)
   0.21200000000000   0.42400000000000   0.00000000000000   O   (6l)
  -0.42400000000000  -0.21200000000000   0.00000000000000   O   (6l)
   0.21200000000000  -0.21200000000000   0.00000000000000   O   (6l)
  -0.21200000000000  -0.42400000000000   0.00000000000000   O   (6l)
   0.42400000000000   0.21200000000000   0.00000000000000   O   (6l)
  -0.21200000000000   0.21200000000000   0.00000000000000   O   (6l)
   0.50000000000000   0.00000000000000   0.00000000000000   W   (3f)
   0.00000000000000   0.50000000000000   0.00000000000000   W   (3f)
   0.50000000000000   0.50000000000000   0.00000000000000   W   (3f)
```

$D0_6$ (Tysonite, LaF_3) (*obsolete*): A3B_hP24_193_ack_g - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Tysonite'
_chemical_formula_sum 'F3 La'

loop_
_publ_author_name
 'I. Oftedal'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Physik B Condensed Matter
;
_journal_volume 13
_journal_year 1931
_journal_page_first 190
_journal_page_last 200
_publ_Section_title
;
 Zur Kristallstruktur von Tysonit (Ce, La, ...)F$_{3}$
;

# Found in Strukturbericht Band II 1928-1932, 1937

_aflow_title '$D0_{6}$ (Tysonite, LaF$_{3}$) ({\em{obsolete}}) Structure
    ↪ '
_aflow_proto 'A3B_hP24_193_ack_g'
_aflow_params 'a,c/a,x_{3},x_{4},z_{4}'
_aflow_params_values '7.12,1.02247191011,0.34,0.33333,0.075'
_aflow_Strukturbericht '$D0_{6}$'
_aflow_Pearson 'hP24'

_symmetry_space_group_name_H-M "P 63/m 2/c 2/m"
_symmetry_Int_Tables_number 193

_cell_length_a    7.12000
_cell_length_b    7.12000
_cell_length_c    7.28000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z+1/2
3 -y,x-y,z
4 -x,-y,z+1/2
5 -x+y,-x,z
6 y,-x+y,z+1/2
7 x-y,-y,-z+1/2
8 x,x-y,-z
9 y,x,-z+1/2
10 -x+y,y,-z
11 -x,-x+y,-z+1/2
12 -y,-x,-z
13 -x,-y,-z
14 -x+y,-x,-z+1/2
15 y,-x+y,-z
16 x,y,-z+1/2
17 x-y,x,-z
18 -y,x-y,-z+1/2
19 -x+y,y,z+1/2
20 -x,-x+y,z
21 -y,-x,z+1/2
22 x-y,-y,z
23 x,x-y,z+1/2
24 y,x,z

loop_
_atom_site_label
_atom_site_type_symbol
```

```
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
F1 F 2 a 0.00000 0.00000 0.25000 1.00000
F2 F 4 c 0.33333 0.66667 0.25000 1.00000
La1 La 6 g 0.34000 0.00000 0.25000 1.00000
F3 F 12 k 0.33333 0.00000 0.07500 1.00000
```

$D0_6$ (Tysonite, LaF_3) (*obsolete*): A3B_hP24_193_ack_g - POSCAR

```
A3B_hP24_193_ack_g & a,c/a,x3,x4,z4 --params=7.12,1.02247191011,0.34,
    ↪ 0.33333,0.075 & P6_{3}/mcm D_{6h}^{3} #193 (acgk) & hP24 & $D0_
    ↪ {6}$ & F3La & Tysonite & I. Oftedal, Z. Phys. B: Condens.
    ↪ Matter 13, 190–200 (1931)
   1.00000000000000
   3.56000000000000  -6.16610087494520   0.00000000000000
   3.56000000000000   6.16610087494520   0.00000000000000
   0.00000000000000   0.00000000000000   7.28000000000000
     F     La
    18      6
Direct
   0.00000000000000   0.00000000000000   0.25000000000000   F   (2a)
   0.00000000000000   0.00000000000000   0.75000000000000   F   (2a)
   0.33333333333333   0.66666666666667   0.25000000000000   F   (4c)
   0.66666666666667   0.33333333333333   0.75000000000000   F   (4c)
   0.66666666666667   0.33333333333333   0.25000000000000   F   (4c)
   0.33333333333333   0.66666666666667   0.75000000000000   F   (4c)
   0.33333000000000   0.00000000000000   0.07500000000000   F  (12k)
   0.00000000000000   0.33333000000000   0.07500000000000   F  (12k)
  -0.33333000000000  -0.33333000000000   0.07500000000000   F  (12k)
  -0.33333000000000   0.00000000000000   0.57500000000000   F  (12k)
   0.00000000000000  -0.33333000000000   0.57500000000000   F  (12k)
   0.33333000000000   0.33333000000000   0.57500000000000   F  (12k)
   0.00000000000000   0.33333000000000   0.42500000000000   F  (12k)
   0.33333000000000   0.00000000000000   0.42500000000000   F  (12k)
  -0.33333000000000  -0.33333000000000   0.42500000000000   F  (12k)
   0.00000000000000  -0.33333000000000  -0.07500000000000   F  (12k)
  -0.33333000000000   0.00000000000000  -0.07500000000000   F  (12k)
   0.33333000000000   0.33333000000000  -0.07500000000000   F  (12k)
   0.34000000000000   0.00000000000000   0.25000000000000  La   (6g)
   0.00000000000000   0.34000000000000   0.25000000000000  La   (6g)
  -0.34000000000000  -0.34000000000000   0.25000000000000  La   (6g)
  -0.34000000000000   0.00000000000000   0.75000000000000  La   (6g)
   0.00000000000000  -0.34000000000000   0.75000000000000  La   (6g)
   0.34000000000000   0.34000000000000   0.75000000000000  La   (6g)
```

Ti_5Ga_4: A4B5_hP18_193_bg_dg - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Ga4Ti5'
_chemical_formula_sum 'Ga4 Ti5'

loop_
_publ_author_name
 'K. Schubert'
 'H. G. Meissner'
 'M. P\"{o}tzschke'
 'W. Rossteutscher'
 'E. Stolz'
_journal_name_full_name
;
 Naturwissenschaften
;
_journal_volume 49
_journal_year 1962
_journal_page_first 57
_journal_page_last 57
_publ_Section_title
;
 Einige Strukturdaten metallischer Phasen (7)
;

# Found in Ti$_{5}$Ga$_{4}$ Crystal Structure, {PAULING FILE in:
    ↪ Inorganic Solid Phases, SpringerMaterials (online database),
    ↪ Springer, Heidelberg (ed.) SpringerMaterials }, 2016

_aflow_title 'Ti$_{5}$Ga$_{4}$ Structure'
_aflow_proto 'A4B5_hP18_193_bg_dg'
_aflow_params 'a,c/a,x_{3},x_{4}'
_aflow_params_values '7.861,0.693550438875,0.62,0.29'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hP18'

_symmetry_space_group_name_H-M "P 63/m 2/c 2/m"
_symmetry_Int_Tables_number 193

_cell_length_a    7.86100
_cell_length_b    7.86100
_cell_length_c    5.45200
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z+1/2
3 -y,x-y,z
4 -x,-y,z+1/2
5 -x+y,-x,z
6 y,-x+y,z+1/2
7 x-y,-y,-z+1/2
8 x,x-y,-z
9 y,x,-z+1/2
10 -x+y,y,-z
11 -x,-x+y,-z+1/2
12 -y,-x,-z
13 -x,-y,-z
14 -x+y,-x,-z+1/2
15 y,-x+y,-z
16 x,y,-z+1/2
17 x-y,x,-z
18 -y,x-y,-z+1/2
19 -x+y,y,z+1/2
20 -x,-x+y,z
21 -y,-x,z+1/2
22 x-y,-y,z
23 x,x-y,z+1/2
24 y,x,z

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ga1 Ga 2 b 0.00000 0.00000 0.00000 1.00000
Ti1 Ti 4 d 0.33333 0.66667 0.00000 1.00000
Ga2 Ga 6 g 0.62000 0.00000 0.25000 1.00000
Ti2 Ti 6 g 0.29000 0.00000 0.25000 1.00000
```

Ti_5Ga_4: A4B5_hP18_193_bg_dg - POSCAR

```
A4B5_hP18_193_bg_dg & a,c/a,x3,x4 --params=7.861,0.693550438875,0.62,
    ↪ 0.29 & P6_{3}/mcm D_{6h}^{3} #193 (bdg^2) & hP18 & None &
    ↪ Ga4Ti5 & Ga4Ti5 & K. Schubert et al., Naturwissenschaften 49,
    ↪ 57(1962)
   1.00000000000000
   3.93050000000000  -6.80782569914947   0.00000000000000
   3.93050000000000   6.80782569914947   0.00000000000000
   0.00000000000000   0.00000000000000   5.45200000000000
    Ga    Ti
     8    10
Direct
   0.00000000000000   0.00000000000000   0.00000000000000   Ga   (2b)
   0.00000000000000   0.00000000000000   0.50000000000000   Ga   (2b)
   0.62000000000000   0.00000000000000   0.25000000000000   Ga   (6g)
   0.00000000000000   0.62000000000000   0.25000000000000   Ga   (6g)
  -0.62000000000000  -0.62000000000000   0.25000000000000   Ga   (6g)
  -0.62000000000000   0.00000000000000   0.75000000000000   Ga   (6g)
   0.00000000000000  -0.62000000000000   0.75000000000000   Ga   (6g)
   0.62000000000000   0.62000000000000   0.75000000000000   Ga   (6g)
   0.33333333333333   0.66666666666667   0.00000000000000   Ti   (4d)
   0.66666666666667   0.33333333333333   0.50000000000000   Ti   (4d)
   0.66666666666667   0.33333333333333   0.00000000000000   Ti   (4d)
   0.33333333333333   0.66666666666667   0.50000000000000   Ti   (4d)
   0.29000000000000   0.00000000000000   0.25000000000000   Ti   (6g)
   0.00000000000000   0.29000000000000   0.25000000000000   Ti   (6g)
  -0.29000000000000  -0.29000000000000   0.25000000000000   Ti   (6g)
  -0.29000000000000   0.00000000000000   0.75000000000000   Ti   (6g)
   0.00000000000000  -0.29000000000000   0.75000000000000   Ti   (6g)
   0.29000000000000   0.29000000000000   0.75000000000000   Ti   (6g)
```

Proposed 300 GPa HfH_{10}: A10B_hP22_194_bhj_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'H10Hf'
_chemical_formula_sum 'H10 Hf'

loop_
_publ_author_name
 'H. Xie'
 'Y. Yao'
 'X. Feng'
 'D. Duan'
 'H. Song'
 'Z. Zhang'
 'S. Jiang'
 'S. A. T. Redfern'
 'V. Z. Kresin'
 'C. J. Pickard'
 'T. Cui'
_journal_name_full_name
;
 Physical Review Letters
;
_journal_volume 125
_journal_year 2020
_journal_page_first 217001
_journal_page_last 217001
_publ_Section_title
;
 Hydrogen pentagraphenelike structure stabilized by hafnium: a
    ↪ high-temperature conventional superconductor
;

_aflow_title 'Proposed 300~GPa HfH$_{10}$ Structure'
```

```
_aflow_proto 'A10B_hP22_194_bhj_c'
_aflow_params 'a,c/a,x_{3},x_{4},y_{4}'
_aflow_params_values '4.633,0.562702352687,0.115,0.375,0.077'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hP22'

_symmetry_space_group_name_H-M "P 63/m 2/m 2/c"
_symmetry_Int_Tables_number 194

_cell_length_a 4.63300
_cell_length_b 4.63300
_cell_length_c 2.60700
_cell_angle_alpha 90.00000
_cell_angle_beta 90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z+1/2
3 -y,x-y,z
4 -x,-y,z+1/2
5 -x+y,-x,z
6 y,-x+y,z+1/2
7 x-y,-y,-z
8 x,x-y,-z+1/2
9 y,x,-z
10 -x+y,y,-z+1/2
11 -x,-x+y,-z
12 -y,-x,-z+1/2
13 -x,-y,-z
14 -x+y,-x,-z+1/2
15 y,-x+y,-z
16 x,y,-z+1/2
17 x-y,x,-z
18 -y,x-y,-z+1/2
19 -x+y,y,z
20 -x,-x+y,z+1/2
21 -y,-x,z
22 x-y,-y,z+1/2
23 x,x-y,z
24 y,x,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
H1 H 2 b 0.00000 0.00000 0.25000 1.00000
Hf1 Hf 2 c 0.33333 0.66667 0.25000 1.00000
H2 H 6 h 0.11500 0.23000 0.25000 1.00000
H3 H 12 j 0.37500 0.07700 0.25000 1.00000
```

Proposed 300 GPa HfH_{10}: A10B_hP22_194_bhj_c - POSCAR

```
A10B_hP22_194_bhj_c & a,c/a,x3,x4,y4 --params=4.633,0.562702352687,0.115
    ↪ ,0.375,0.077 & P6_{3}/mmc D_{6h}^{4} #194 (bchj) & hP22 & None
    ↪ & H10Hf & H10Hf & H. Xie et al., Phys. Rev. Lett. 125, 217001 (
    ↪ 2020)
   1.00000000000000
   2.31650000000000  -4.01229569573330   0.00000000000000
   2.31650000000000   4.01229569573330   0.00000000000000
   0.00000000000000   0.00000000000000   2.60700000000000
     H     Hf
    20      2
Direct
   0.00000000000000   0.00000000000000   0.25000000000000   H   (2b)
   0.00000000000000   0.00000000000000   0.75000000000000   H   (2b)
   0.11500000000000   0.23000000000000   0.25000000000000   H   (6h)
  -0.23000000000000  -0.11500000000000   0.25000000000000   H   (6h)
   0.11500000000000  -0.11500000000000   0.25000000000000   H   (6h)
  -0.11500000000000  -0.23000000000000   0.75000000000000   H   (6h)
   0.23000000000000   0.11500000000000   0.75000000000000   H   (6h)
  -0.11500000000000   0.11500000000000   0.75000000000000   H   (6h)
   0.37500000000000   0.07700000000000   0.25000000000000   H   (12j)
  -0.07700000000000   0.29800000000000   0.25000000000000   H   (12j)
  -0.29800000000000  -0.37500000000000   0.25000000000000   H   (12j)
  -0.37500000000000  -0.07700000000000   0.75000000000000   H   (12j)
   0.07700000000000  -0.29800000000000   0.75000000000000   H   (12j)
   0.29800000000000   0.37500000000000   0.75000000000000   H   (12j)
   0.07700000000000   0.37500000000000   0.75000000000000   H   (12j)
   0.29800000000000  -0.07700000000000   0.75000000000000   H   (12j)
  -0.37500000000000  -0.29800000000000   0.75000000000000   H   (12j)
  -0.07700000000000  -0.37500000000000   0.25000000000000   H   (12j)
  -0.29800000000000   0.07700000000000   0.25000000000000   H   (12j)
   0.37500000000000   0.29800000000000   0.25000000000000   H   (12j)
   0.33333333333333   0.66666666666667   0.25000000000000   Hf   (2c)
   0.66666666666667   0.33333333333333   0.75000000000000   Hf   (2c)
```

Magnetoplumbite ($PbFe_{12}O_{19}$): A12B19C_hP64_194_ab2fk_efh2k_d - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Magnetoplumbite'
_chemical_formula_sum 'Fe12 O19 Pb'

loop_
_publ_author_name
 'R. Gerber'
 'Z. \v{S}im\v{s}a'
 'L. Jen\v{s}ovsk\\'{y}'
_journal_name_full_name
;
 Czechoslovak Journal of Physics
;
_journal_volume 44
_journal_year 1994
_journal_page_first 937
_journal_page_last 940
_publ_Section_title
;
 A note on the magnetoplumbite crystal structure
;

_aflow_title 'Magnetoplumbite (PbFe$_{12}$O$_{19}$) Structure'
_aflow_proto 'A12B19C_hP64_194_ab2fk_efh2k_d'
_aflow_params 'a,c/a,z_{4},z_{5},z_{6},z_{7},x_{8},x_{9},z_{9},x_{10},z_
    ↪ {10},x_{11},z_{11}'
_aflow_params_values '5.88,3.91496598639,0.15,0.028,0.19,-0.05,0.182,
    ↪ 0.167,0.892,0.167,0.05,0.5,0.15'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hP64'

_symmetry_space_group_name_H-M "P 63/m 2/m 2/c"
_symmetry_Int_Tables_number 194

_cell_length_a 5.88000
_cell_length_b 5.88000
_cell_length_c 23.02000
_cell_angle_alpha 90.00000
_cell_angle_beta 90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z+1/2
3 -y,x-y,z
4 -x,-y,z+1/2
5 -x+y,-x,z
6 y,-x+y,z+1/2
7 x-y,-y,-z
8 x,x-y,-z+1/2
9 y,x,-z
10 -x+y,y,-z+1/2
11 -x,-x+y,-z
12 -y,-x,-z+1/2
13 -x,-y,-z
14 -x+y,-x,-z+1/2
15 y,-x+y,-z
16 x,y,-z+1/2
17 x-y,x,-z
18 -y,x-y,-z+1/2
19 -x+y,y,z
20 -x,-x+y,z+1/2
21 -y,-x,z
22 x-y,-y,z+1/2
23 x,x-y,z
24 y,x,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Fe1 Fe 2 a 0.00000 0.00000 0.00000 1.00000
Fe2 Fe 2 b 0.00000 0.00000 0.25000 1.00000
Pb1 Pb 2 d 0.33333 0.66667 0.75000 1.00000
O1 O 4 e 0.00000 0.00000 0.15000 1.00000
Fe3 Fe 4 f 0.33333 0.66667 0.02800 1.00000
Fe4 Fe 4 f 0.33333 0.66667 0.19000 1.00000
O2 O 4 f 0.33333 0.66667 -0.05000 1.00000
O3 O 6 h 0.18200 0.36400 0.25000 1.00000
Fe5 Fe 12 k 0.16700 0.33400 0.89200 1.00000
O4 O 12 k 0.16700 0.33400 0.05000 1.00000
O5 O 12 k 0.50000 0.00000 0.15000 1.00000
```

Magnetoplumbite ($PbFe_{12}O_{19}$): A12B19C_hP64_194_ab2fk_efh2k_d - POSCAR

```
A12B19C_hP64_194_ab2fk_efh2k_d & a,c/a,z4,z5,z6,z7,x8,x9,z9,x10,z10,x11,
    ↪ z11 --params=5.88,3.91496598639,0.15,0.028,0.19,-0.05,0.182,
    ↪ 0.167,0.892,0.167,0.05,0.5,0.15 & P6_{3}/mmc D_{6h}^{4} #194 (
    ↪ abdef^3hk^3) & hP64 & None & Fe12O19Pb & Magnetoplumbite & R.
    ↪ Gerber and Z. \v{S}im\v{s}a and L. Jen\v{s}ovsk\'{y}, Czech. J.
    ↪ Phys. 44, 937-940 (1994)
   1.00000000000000
   2.94000000000000  -5.09222937425250   0.00000000000000
   2.94000000000000   5.09222937425250   0.00000000000000
   0.00000000000000   0.00000000000000  23.02000000000000
    Fe     O     Pb
    24    38     2
Direct
   0.00000000000000   0.00000000000000   0.00000000000000   Fe   (2a)
   0.00000000000000   0.00000000000000   0.50000000000000   Fe   (2a)
   0.00000000000000   0.00000000000000   0.25000000000000   Fe   (2b)
   0.00000000000000   0.00000000000000   0.75000000000000   Fe   (2b)
   0.33333333333333   0.66666666666667   0.02800000000000   Fe   (4f)
   0.66666666666667   0.33333333333333   0.52800000000000   Fe   (4f)
```

```
  0.66666666666667   0.33333333333333  -0.02800000000000   Fe   (4f)
  0.33333333333333   0.66666666666667   0.47200000000000   Fe   (4f)
  0.33333333333333   0.66666666666667   0.19000000000000   Fe   (4f)
  0.66666666666667   0.33333333333333   0.69000000000000   Fe   (4f)
  0.66666666666667   0.33333333333333  -0.19000000000000   Fe   (4f)
  0.33333333333333   0.66666666666667   0.31000000000000   Fe   (4f)
  0.16700000000000   0.33400000000000   0.89200000000000   Fe  (12k)
 -0.33400000000000  -0.16700000000000   0.89200000000000   Fe  (12k)
  0.16700000000000  -0.16700000000000   0.89200000000000   Fe  (12k)
 -0.16700000000000  -0.33400000000000   1.39200000000000   Fe  (12k)
  0.33400000000000   0.16700000000000   1.39200000000000   Fe  (12k)
 -0.16700000000000   0.16700000000000   1.39200000000000   Fe  (12k)
  0.33400000000000   0.16700000000000  -0.89200000000000   Fe  (12k)
 -0.16700000000000  -0.33400000000000  -0.89200000000000   Fe  (12k)
 -0.16700000000000   0.16700000000000  -0.89200000000000   Fe  (12k)
 -0.33400000000000  -0.16700000000000  -0.39200000000000   Fe  (12k)
  0.16700000000000   0.33400000000000  -0.39200000000000   Fe  (12k)
  0.16700000000000  -0.16700000000000  -0.39200000000000   Fe  (12k)
  0.00000000000000   0.00000000000000   0.15000000000000    O   (4e)
  0.00000000000000   0.00000000000000   0.65000000000000    O   (4e)
  0.00000000000000   0.00000000000000  -0.15000000000000    O   (4e)
  0.00000000000000   0.00000000000000   0.35000000000000    O   (4e)
  0.33333333333333   0.66666666666667  -0.05000000000000    O   (4f)
  0.66666666666667   0.33333333333333   0.45000000000000    O   (4f)
  0.66666666666667   0.33333333333333   0.05000000000000    O   (4f)
  0.33333333333333   0.66666666666667   0.55000000000000    O   (4f)
  0.18200000000000   0.36400000000000   0.25000000000000    O   (6h)
 -0.36400000000000  -0.18200000000000   0.25000000000000    O   (6h)
  0.18200000000000  -0.18200000000000   0.25000000000000    O   (6h)
 -0.18200000000000  -0.36400000000000   0.75000000000000    O   (6h)
  0.36400000000000   0.18200000000000   0.75000000000000    O   (6h)
 -0.18200000000000   0.18200000000000   0.75000000000000    O   (6h)
  0.16700000000000   0.33400000000000   0.05000000000000    O  (12k)
 -0.33400000000000  -0.16700000000000   0.05000000000000    O  (12k)
  0.16700000000000  -0.16700000000000   0.05000000000000    O  (12k)
 -0.16700000000000  -0.33400000000000   0.55000000000000    O  (12k)
  0.33400000000000   0.16700000000000   0.55000000000000    O  (12k)
 -0.16700000000000   0.16700000000000   0.55000000000000    O  (12k)
  0.33400000000000   0.16700000000000  -0.05000000000000    O  (12k)
 -0.16700000000000  -0.33400000000000  -0.05000000000000    O  (12k)
 -0.16700000000000   0.16700000000000  -0.05000000000000    O  (12k)
 -0.33400000000000  -0.16700000000000   0.45000000000000    O  (12k)
  0.16700000000000   0.33400000000000   0.45000000000000    O  (12k)
  0.16700000000000  -0.16700000000000   0.45000000000000    O  (12k)
  0.50000000000000   1.00000000000000   0.15000000000000    O  (12k)
 -1.00000000000000  -0.50000000000000   0.15000000000000    O  (12k)
  0.50000000000000  -0.50000000000000   0.15000000000000    O  (12k)
 -0.50000000000000  -1.00000000000000   0.65000000000000    O  (12k)
  1.00000000000000   0.50000000000000   0.65000000000000    O  (12k)
 -0.50000000000000   0.50000000000000   0.65000000000000    O  (12k)
  1.00000000000000   0.50000000000000  -0.15000000000000    O  (12k)
 -0.50000000000000  -1.00000000000000  -0.15000000000000    O  (12k)
 -0.50000000000000   0.50000000000000  -0.15000000000000    O  (12k)
 -1.00000000000000  -0.50000000000000   0.35000000000000    O  (12k)
  0.50000000000000   1.00000000000000   0.35000000000000    O  (12k)
  0.50000000000000  -0.50000000000000   0.35000000000000    O  (12k)
  0.33333333333333   0.66666666666667   0.75000000000000   Pb   (2d)
  0.66666666666667   0.33333333333333   0.25000000000000   Pb   (2d)
```

Pt_2Sn_3 ($D5_b$): A2B3_hP10_194_f_bf - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Pt2Sn3'
_chemical_formula_sum 'Pt2 Sn3'

loop_
_publ_author_name
 'K. Schubert'
 'H. Pfisterer'
_journal_name_full_name
;
 Zeitschrift fur Metallkunde
;
_journal_volume 40
_journal_year 1949
_journal_page_first 405
_journal_page_last 405
_publ_Section_title
;
 Kristallstruktur von Pt${2}$Sn${3}$
;

# Found in A Handbook of Lattice Spacings and Structures of Metals and
      ↪ Alloys, {N.~R.~C. No. 4303}, 1958

_aflow_title 'Pt${2}$Sn${3}$ ($D5_{b}$) Structure'
_aflow_proto 'A2B3_hP10_194_f_bf'
_aflow_params 'a,c/a,z_{2},z_{3}'
_aflow_params_values '4.36079,2.97196838188,0.143,-0.07'
_aflow_Strukturbericht '$D5_{b}$'
_aflow_Pearson 'hP10'

_symmetry_space_group_name_H-M "P 63/m 2/m 2/c"
_symmetry_Int_Tables_number 194

_cell_length_a     4.36079
_cell_length_b     4.36079
_cell_length_c     12.96013
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000
loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z+1/2
3 -y,x-y,z
4 -x,-y,z+1/2
5 -x+y,-x,z
6 y,-x+y,z+1/2
7 x-y,-y,-z
8 x,x-y,-z+1/2
9 y,x,-z
10 -x+y,y,-z+1/2
11 -x,-x+y,-z
12 -y,-x,-z+1/2
13 -x,-y,-z
14 -x+y,-x,-z+1/2
15 y,-x+y,-z
16 x,y,-z+1/2
17 x-y,x,-z
18 -y,x-y,-z+1/2
19 -x+y,y,z
20 -x,-x+y,z+1/2
21 -y,-x,z
22 x-y,-y,z+1/2
23 x,x-y,z
24 y,x,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Sn1 Sn 2 b 0.00000 0.00000 0.25000 1.00000
Pt1 Pt 4 f 0.33333 0.66667 0.14300 1.00000
Sn2 Sn 4 f 0.33333 0.66667 -0.07000 1.00000
```

Pt_2Sn_3 ($D5_b$): A2B3_hP10_194_f_bf - POSCAR

```
A2B3_hP10_194_f_bf & a,c/a,z2,z3 --params=4.36079,2.97196838188,0.143,-
      ↪ 0.07 & P6_{3}/mmc D_{6h}^{4} #194 (bf^2) & hP10 & $D5_{b}$ &
      ↪ Pt2Sn3 & Pt2Sn3 & K. Schubert and H. Pfisterer, Z. Metallkd. 40
      ↪ , 405(1949)
   1.00000000000000
   2.18039500000000  -3.77655492056914   0.00000000000000
   2.18039500000000   3.77655492056914   0.00000000000000
   0.00000000000000   0.00000000000000  12.96013000000000
    Pt    Sn
     4     6
Direct
   0.33333333333333   0.66666666666667   0.14300000000000   Pt   (4f)
   0.66666666666667   0.33333333333333   0.64300000000000   Pt   (4f)
   0.66666666666667   0.33333333333333  -0.14300000000000   Pt   (4f)
   0.33333333333333   0.66666666666667   0.35700000000000   Pt   (4f)
   0.00000000000000   0.00000000000000   0.25000000000000   Sn   (2b)
   0.00000000000000   0.00000000000000   0.75000000000000   Sn   (2b)
   0.33333333333333   0.66666666666667  -0.07000000000000   Sn   (4f)
   0.66666666666667   0.33333333333333   0.43000000000000   Sn   (4f)
   0.66666666666667   0.33333333333333   0.07000000000000   Sn   (4f)
   0.33333333333333   0.66666666666667   0.57000000000000   Sn   (4f)
```

β-Alumina (Al_2O_3, $D5_6$): A2B3_hP60_194_3fk_cdef2k - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral '$\beta$-alumina'
_chemical_formula_sum 'Al2 O3'

loop_
_publ_author_name
 'W. L. Bragg'
 'C. Gottfried'
 'J. West'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 77
_journal_year 1931
_journal_page_first 255
_journal_page_last 274
_publ_Section_title
;
 The Structure of $\beta$ Alumina
;

# Found in Strukturbericht Band II 1928-1932, 1937

_aflow_title '$\beta$-Alumina (Al$_{2}$O$_{3}$, $D5_{6}$) Structure'
_aflow_proto 'A2B3_hP60_194_3fk_cdef2k'
_aflow_params 'a,c/a,z_{3},z_{4},z_{5},z_{6},z_{7},x_{8},z_{8},x_{9},z_{
      ↪ 9},x_{10},z_{10}'
_aflow_params_values '5.56,4.05575539568,0.14,0.02,0.17,-0.17,-0.05,
      ↪ 0.83333,0.1,0.16667,0.05,0.5,0.14'
_aflow_Strukturbericht '$D5_{6}$'
_aflow_Pearson 'hP60'

_symmetry_space_group_name_H-M "P 63/m 2/m 2/c"
_symmetry_Int_Tables_number 194
```

```
_cell_length_a      5.56000
_cell_length_b      5.56000
_cell_length_c      22.55000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z+1/2
3 -y,x-y,z
4 -x,-y,z+1/2
5 -x+y,-x,z
6 y,-x+y,z+1/2
7 x-y,-y,-z
8 x,x-y,-z+1/2
9 y,x,-z
10 -x+y,y,-z+1/2
11 -x,-x+y,-z
12 -y,-x,-z+1/2
13 -x,-y,-z
14 -x+y,-x,-z+1/2
15 y,-x+y,-z
16 x,y,-z+1/2
17 x-y,x,-z
18 -y,x-y,-z+1/2
19 -x+y,y,z
20 -x,-x+y,z+1/2
21 -y,-x,z
22 x-y,-y,z+1/2
23 x,x-y,z
24 y,x,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 2 c 0.33333 0.66667 0.25000 1.00000
O2 O 2 d 0.33333 0.66667 0.75000 1.00000
O3 O 4 e 0.00000 0.00000 0.14000 1.00000
Al1 Al 4 f 0.33333 0.66667 0.02000 1.00000
Al2 Al 4 f 0.33333 0.66667 0.17000 1.00000
Al3 Al 4 f 0.33333 0.66667 -0.1700 1.00000
O4 O 4 f 0.33333 0.66667 -0.0500 1.00000
Al4 Al 12 k 0.83333 0.66667 0.10000 1.00000
O5 O 12 k 0.16667 0.33333 0.05000 1.00000
O6 O 12 k 0.50000 0.00000 0.14000 1.00000
```

β-Alumina (Al_2O_3, $D5_6$): A2B3_hP60_194_3fk_cdef2k - POSCAR

```
A2B3_hP60_194_3fk_cdef2k & a,c/a,z3,z4,z5,z6,z7,x8,z8,x9,z9,x10,z10 --
    ↪ params=5.56,4.05575539568,0.14,0.02,0.17,-0.17,-0.05,0.83333,
    ↪ 0.1,0.16667,0.05,0.5,0.14 & P6_{3}/mmc D_{6h}^{4} #194 (cdef^4k
    ↪ ^3) & hP60 & $D5_{6}$ & Al2O3 & $\beta$-alumina & W. L. Bragg
    ↪ and C. Gottfried and J. West, Zeitschrift f{"u}r
    ↪ Kristallographie - Crystalline Materials 77, 255-274 (1931)
   1.00000000000000
   2.78000000000000  -4.81510124504148   0.00000000000000
   2.78000000000000   4.81510124504148   0.00000000000000
   0.00000000000000   0.00000000000000  22.55000000000000
    Al   O
    24  36
Direct
   0.33333333333333   0.66666666666667   0.02000000000000  Al  (4f)
   0.66666666666667   0.33333333333333   0.52000000000000  Al  (4f)
   0.66666666666667   0.33333333333333  -0.02000000000000  Al  (4f)
   0.33333333333333   0.66666666666667   0.48000000000000  Al  (4f)
   0.33333333333333   0.66666666666667   0.17000000000000  Al  (4f)
   0.66666666666667   0.33333333333333   0.67000000000000  Al  (4f)
   0.66666666666667   0.33333333333333  -0.17000000000000  Al  (4f)
   0.33333333333333   0.66666666666667   0.33000000000000  Al  (4f)
   0.33333333333333   0.66666666666667  -0.17000000000000  Al  (4f)
   0.66666666666667   0.33333333333333   0.33000000000000  Al  (4f)
   0.66666666666667   0.33333333333333   0.17000000000000  Al  (4f)
   0.33333333333333   0.66666666666667   0.67000000000000  Al  (4f)
   0.83333000000000   1.66666000000000   0.10000000000000  Al (12k)
  -1.66666000000000  -0.83333000000000   0.10000000000000  Al (12k)
   0.83333000000000  -0.83333000000000   0.10000000000000  Al (12k)
  -0.83333000000000  -1.66666000000000   0.60000000000000  Al (12k)
   1.66666000000000   0.83333000000000   0.60000000000000  Al (12k)
  -0.83333000000000   0.83333000000000   0.60000000000000  Al (12k)
   1.66666000000000   0.83333000000000  -0.10000000000000  Al (12k)
  -0.83333000000000  -1.66666000000000  -0.10000000000000  Al (12k)
  -0.83333000000000   0.83333000000000  -0.10000000000000  Al (12k)
  -1.66666000000000  -0.83333000000000   0.40000000000000  Al (12k)
   0.83333000000000   1.66666000000000   0.40000000000000  Al (12k)
   0.83333000000000  -0.83333000000000   0.40000000000000  Al (12k)
   0.33333333333333   0.66666666666667   0.25000000000000   O  (2c)
   0.66666666666667   0.33333333333333   0.75000000000000   O  (2c)
   0.33333333333333   0.66666666666667   0.75000000000000   O  (2d)
   0.66666666666667   0.33333333333333   0.25000000000000   O  (2d)
   0.00000000000000   0.00000000000000   0.14000000000000   O  (4e)
   0.00000000000000   0.00000000000000   0.64000000000000   O  (4e)
   0.00000000000000   0.00000000000000  -0.14000000000000   O  (4e)
   0.00000000000000   0.00000000000000   0.36000000000000   O  (4e)
   0.33333333333333   0.66666666666667  -0.05000000000000   O  (4f)
   0.66666666666667   0.33333333333333   0.45000000000000   O  (4f)
   0.66666666666667   0.33333333333333   0.05000000000000   O  (4f)
   0.33333333333333   0.66666666666667   0.55000000000000   O  (4f)
   0.16667000000000   0.33334000000000   0.05000000000000   O (12k)
  -0.33334000000000  -0.16667000000000   0.05000000000000   O (12k)
   0.16667000000000  -0.16667000000000   0.05000000000000   O (12k)
  -0.16667000000000  -0.33334000000000   0.55000000000000   O (12k)
   0.33334000000000   0.16667000000000   0.55000000000000   O (12k)
  -0.16667000000000   0.16667000000000   0.55000000000000   O (12k)
   0.33334000000000   0.16667000000000  -0.05000000000000   O (12k)
  -0.16667000000000  -0.33334000000000  -0.05000000000000   O (12k)
  -0.16667000000000   0.16667000000000  -0.05000000000000   O (12k)
  -0.33334000000000  -0.16667000000000   0.45000000000000   O (12k)
   0.16667000000000   0.33334000000000   0.45000000000000   O (12k)
   0.16667000000000  -0.16667000000000   0.45000000000000   O (12k)
   0.50000000000000   1.00000000000000   0.14000000000000   O (12k)
  -1.00000000000000  -0.50000000000000   0.14000000000000   O (12k)
   0.50000000000000  -0.50000000000000   0.14000000000000   O (12k)
  -0.50000000000000  -1.00000000000000   0.64000000000000   O (12k)
   1.00000000000000   0.50000000000000   0.64000000000000   O (12k)
  -0.50000000000000   0.50000000000000   0.64000000000000   O (12k)
   1.00000000000000   0.50000000000000  -0.14000000000000   O (12k)
  -0.50000000000000  -1.00000000000000  -0.14000000000000   O (12k)
  -0.50000000000000   0.50000000000000  -0.14000000000000   O (12k)
  -1.00000000000000  -0.50000000000000   0.36000000000000   O (12k)
   0.50000000000000   1.00000000000000   0.36000000000000   O (12k)
   0.50000000000000  -0.50000000000000   0.36000000000000   O (12k)
```

$S3_4$ (II) (Catapleiite, $Na_2Zr(SiO_3)_3 \cdot H_2O$) (*obsolete*): A3B2C9D3E_hP36_194_g_f_hk_h_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Catapleiite'
_chemical_formula_sum '(H2O)3 Na2 O9 Si3 Zr'

loop_
_publ_author_name
 'B. Brunowski'
_journal_name_full_name
;
 Acta Physicochimica U.R.S.S.
;
_journal_volume 5
_journal_year 1936
_journal_page_first 863
_journal_page_last 892
_publ_Section_title
;
 Die Struktur des Katapleits
;

# Found in Strukturbericht Band V 1937, 1940

_aflow_title '$S3_{4}$ (II) (Catapleiite, Na$_{2}$Zr(SiO$_{2}$)$_{3}$$\
    ↪ cdot$ H$_{2}$O) ({\em{obsolete}}) Structure'
_aflow_proto 'A3B2C9D3E_hP36_194_g_f_hk_h_a'
_aflow_params 'a,c/a,z_{2},x_{4},x_{5},x_{6},z_{6}'
_aflow_params_values '7.39,1.3599458728,0.08,0.47,0.2,0.136,0.125'
_aflow_Strukturbericht '$S3_{4}$ (II)'
_aflow_Pearson 'hP36'

_symmetry_space_group_name_H-M "P 63/m 2/m 2/c"
_symmetry_Int_Tables_number 194

_cell_length_a      7.39000
_cell_length_b      7.39000
_cell_length_c      10.05000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z+1/2
3 -y,x-y,z
4 -x,-y,z+1/2
5 -x+y,-x,z
6 y,-x+y,z+1/2
7 x-y,-y,-z
8 x,x-y,-z+1/2
9 y,x,-z
10 -x+y,y,-z+1/2
11 -x,-x+y,-z
12 -y,-x,-z+1/2
13 -x,-y,-z
14 -x+y,-x,-z+1/2
15 y,-x+y,-z
16 x,y,-z+1/2
17 x-y,x,-z
18 -y,x-y,-z+1/2
19 -x+y,y,z
20 -x,-x+y,z+1/2
21 -y,-x,z
22 x-y,-y,z+1/2
23 x,x-y,z
24 y,x,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
```

```
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Zr1 Zr 2 a 0.00000 0.00000 0.00000 1.00000
Na1 Na 4 f 0.33333 0.66667 0.08000 1.00000
H2O1 H2O 6 g 0.50000 0.00000 0.00000 0.66667
O1 O 6 h 0.47000 0.94000 0.25000 1.00000
Si1 Si 6 h 0.20000 0.40000 0.25000 1.00000
O2 O 12 k 0.13600 0.27200 0.12500 1.00000
```

$S3_4$ (II) (Catapleiite, $Na_2Zr(SiO_3)_3 \cdot H_2O$) (*obsolete*): A3B2C9D3E_hP36_194_g_f_hk_h_a - POSCAR

```
A3B2C9D3E_hP36_194_g_f_hk_h_a & a,c/a,z2,x4,x5,x6,z6 --params=7.39,
      ↪ 1.3599458728,0.08,0.47,0.2,0.136,0.125 & P6_{3}/mmc D_{6h}^{4}
      ↪ #194 (afgh^2k) & hP36 & $S3_{4}$ (II) & H2Na2O6Si2Zr &
      ↪ Catapleiite & B. Brunowski, Acta Physicochim. USSR 5, 863-892 (
      ↪ 1936)
   1.00000000000000
   3.69500000000000  -6.39992773396700   0.00000000000000
   3.69500000000000   6.39992773396700   0.00000000000000
   0.00000000000000   0.00000000000000  10.05000000000000
   H2O     Na      O      Si      Zr
     6      4     18       6       2
Direct
   0.50000000000000   0.00000000000000   0.00000000000000  H2O    (6g)
   0.00000000000000   0.50000000000000   0.00000000000000  H2O    (6g)
   0.50000000000000   0.50000000000000   0.00000000000000  H2O    (6g)
   0.50000000000000   0.00000000000000   0.50000000000000  H2O    (6g)
   0.00000000000000   0.50000000000000   0.50000000000000  H2O    (6g)
   0.50000000000000   0.50000000000000   0.50000000000000  H2O    (6g)
   0.33333333333333   0.66666666666667   0.08000000000000  Na     (4f)
   0.66666666666667   0.33333333333333   0.58000000000000  Na     (4f)
   0.66666666666667   0.33333333333333  -0.08000000000000  Na     (4f)
   0.33333333333333   0.66666666666667   0.42000000000000  Na     (4f)
   0.47000000000000   0.94000000000000   0.25000000000000  O      (6h)
  -0.94000000000000  -0.47000000000000   0.25000000000000  O      (6h)
   0.47000000000000  -0.47000000000000   0.25000000000000  O      (6h)
  -0.47000000000000  -0.94000000000000   0.75000000000000  O      (6h)
   0.94000000000000   0.47000000000000   0.75000000000000  O      (6h)
  -0.47000000000000   0.47000000000000   0.75000000000000  O      (6h)
   0.13600000000000   0.27200000000000   0.12500000000000  O     (12k)
  -0.27200000000000  -0.13600000000000   0.12500000000000  O     (12k)
   0.13600000000000  -0.13600000000000   0.12500000000000  O     (12k)
  -0.13600000000000  -0.27200000000000   0.62500000000000  O     (12k)
   0.27200000000000   0.13600000000000   0.62500000000000  O     (12k)
  -0.13600000000000   0.13600000000000   0.62500000000000  O     (12k)
   0.27200000000000   0.13600000000000  -0.12500000000000  O     (12k)
  -0.13600000000000  -0.27200000000000  -0.12500000000000  O     (12k)
  -0.13600000000000   0.13600000000000  -0.12500000000000  O     (12k)
  -0.27200000000000  -0.13600000000000   0.37500000000000  O     (12k)
   0.13600000000000   0.27200000000000   0.37500000000000  O     (12k)
   0.13600000000000  -0.13600000000000   0.37500000000000  O     (12k)
   0.20000000000000   0.40000000000000   0.25000000000000  Si     (6h)
  -0.40000000000000  -0.20000000000000   0.25000000000000  Si     (6h)
   0.20000000000000  -0.20000000000000   0.25000000000000  Si     (6h)
  -0.20000000000000  -0.40000000000000   0.75000000000000  Si     (6h)
   0.40000000000000   0.20000000000000   0.75000000000000  Si     (6h)
  -0.20000000000000   0.20000000000000   0.75000000000000  Si     (6h)
   0.00000000000000   0.00000000000000   0.00000000000000  Zr     (2a)
   0.00000000000000   0.00000000000000   0.50000000000000  Zr     (2a)
```

ReB_3: A3B_hP8_194_af_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'B3Re'
_chemical_formula_sum 'B3 Re'

loop_
_publ_author_name
 'B. Aronsson'
 'E. Stenberg'
 'J. {\AA}selius'
_journal_name_full_name
;
 Acta Chemica Scandinavica
;
_journal_volume 14
_journal_year 1960
_journal_page_first 733
_journal_page_last 741
_publ_Section_title
;
 Borides of Rhenium and the Platinum Metals. The Crystal Structure of
      ↪ Re$_{7}$B$_{3}$, ReB$_{3}$, Rh$_{7}$B$_{3}$, RhB$_{\sim 1.1}$,
      ↪  IrB$_{\sim 1.1}$ and PtB
;

# Found in Pearson's Handbook of Crystallographic Data for Intermetallic
      ↪  Phases, 1991

_aflow_title 'ReB$_{3}$ Structure'
_aflow_proto 'A3B_hP8_194_af_c'
_aflow_params 'a,c/a,z_{3}'
_aflow_params_values '2.9,2.5775862069,0.55'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hP8'

_symmetry_space_group_name_H-M "P 63/m 2/m 2/c"
_symmetry_Int_Tables_number 194

_cell_length_a    2.90000
_cell_length_b    2.90000
_cell_length_c    7.47500
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z+1/2
3 -y,x-y,z
4 -x,-y,z+1/2
5 -x+y,-x,z
6 y,-x+y,z+1/2
7 x-y,-y,-z
8 x,x-y,-z+1/2
9 y,x,-z
10 -x+y,y,-z+1/2
11 -x,-x+y,-z
12 -y,-x,-z+1/2
13 -x,-y,-z
14 -x+y,-x,-z+1/2
15 y,-x+y,-z
16 x,y,-z+1/2
17 x-y,x,-z
18 -y,x-y,-z+1/2
19 -x+y,y,z
20 -x,-x+y,z+1/2
21 -y,-x,z
22 x-y,-y,z+1/2
23 x,x-y,z
24 y,x,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
B1 B 2 a 0.00000 0.00000 0.00000 1.00000
Re1 Re 2 c 0.33333 0.66667 0.25000 1.00000
B2 B 4 f 0.33333 0.66667 0.55000 1.00000
```

ReB_3: A3B_hP8_194_af_c - POSCAR

```
A3B_hP8_194_af_c & a,c/a,z3 --params=2.9,2.5775862069,0.55 & P6_{3}/mmc
      ↪ D_{6h}^{4} #194 (acf) & hP8 & None & B3Re & B3Re & B. Aronsson
      ↪ and E. Stenberg and J. {\AA}selius, Acta Chem. Scand. 14,
      ↪ 733-741 (1960)
   1.00000000000000
   1.45000000000000  -2.51147367097487   0.00000000000000
   1.45000000000000   2.51147367097487   0.00000000000000
   0.00000000000000   0.00000000000000   7.47500000000000
     B      Re
     6       2
Direct
   0.00000000000000   0.00000000000000   0.00000000000000     B     (2a)
   0.00000000000000   0.00000000000000   0.50000000000000     B     (2a)
   0.33333333333333   0.66666666666667   0.55000000000000     B     (4f)
   0.66666666666667   0.33333333333333   1.05000000000000     B     (4f)
   0.66666666666667   0.33333333333333  -0.55000000000000     B     (4f)
   0.33333333333333   0.66666666666667  -0.05000000000000     B     (4f)
   0.33333333333333   0.66666666666667   0.25000000000000    Re     (2c)
   0.66666666666667   0.33333333333333   0.75000000000000    Re     (2c)
```

$Cs_3Cr_2Cl_9$: A9B2C3_hP28_194_hk_f_bf - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Cl9Cr2Cs3'
_chemical_formula_sum 'Cl9 Cr2 Cs3'

loop_
_publ_author_name
 'G. G. Wessel'
 'D. J. W. IJdo'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 10
_journal_year 1957
_journal_page_first 466
_journal_page_last 468
_publ_Section_title
;
 The Crystal Structure of Cs$_{3}$Cr$_{2}$Cl$_{9}$
;

# Found in Structure of the dimer compounds Cs$_{3}$$R$$_{2}$Br$_{9}$ (
      ↪ $R$ = Tb, Dy, Ho, Er, Yb) at 8 and 295-K studied by neutron
      ↪ diffraction, 1989

_aflow_title 'Cs$_{3}$Cr$_{2}$Cl$_{9}$ Structure'
_aflow_proto 'A9B2C3_hP28_194_hk_f_bf'
_aflow_params 'a,c/a,z_{2},z_{3},x_{4},x_{5},z_{5}'
_aflow_params_values '7.22,2.48337950139,0.837,0.077,0.508,0.824,0.092'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hP28'
```

```
_symmetry_space_group_name_H-M "P 63/m 2/m 2/c"
_symmetry_Int_Tables_number 194

_cell_length_a 7.22000
_cell_length_b 7.22000
_cell_length_c 17.93000
_cell_angle_alpha 90.00000
_cell_angle_beta 90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z+1/2
3 -y,x-y,z
4 -x,-y,z+1/2
5 -x+y,-x,z
6 y,-x+y,z+1/2
7 x-y,-y,-z
8 x,x-y,-z+1/2
9 y,x,-z
10 -x+y,y,-z+1/2
11 -x,-x+y,-z
12 -y,-x,-z+1/2
13 -x,-y,-z
14 -x+y,-x,-z+1/2
15 y,-x+y,-z
16 x,y,-z+1/2
17 x-y,x,-z
18 -y,x-y,-z+1/2
19 -x+y,y,z
20 -x,-x+y,z+1/2
21 -y,-x,z
22 x-y,-y,z+1/2
23 x,x-y,z
24 y,x,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cs1 Cs 2 b 0.00000 0.00000 0.25000 1.00000
Cr1 Cr 4 f 0.33333 0.66667 0.83700 1.00000
Cs2 Cs 4 f 0.33333 0.66667 0.07700 1.00000
Cl1 Cl 6 h 0.50800 0.01600 0.25000 1.00000
Cl2 Cl 12 k 0.82400 0.64800 0.09200 1.00000
```

$Cs_3Cr_2Cl_9$: A9B2C3_hP28_194_hk_f_bf - POSCAR

```
A9B2C3_hP28_194_hk_f_bf & a,c/a,z2,z3,x4,x5,z5 --params=7.22,
    ↪ 2.48337950139,0.837,0.077,0.508,0.824,0.092 & P6_{3}/mmc D_{6h
    ↪ }^{4} #194 (bf^2hk) & hP28 & None & Cl9Cr2Cs3 & Cl9Cr2Cs3 & G.
    ↪ G. Wessel and D. J. W. IJdo, Acta Cryst. 10, 466-468 (1957)
   1.00000000000000
   3.61000000000000  -6.25270341532365   0.00000000000000
   3.61000000000000   6.25270341532365   0.00000000000000
   0.00000000000000   0.00000000000000  17.93000000000000
     Cl    Cr    Cs
     18     4     6
Direct
   0.50800000000000   1.01600000000000   0.25000000000000  Cl   (6h)
  -1.01600000000000  -0.50800000000000   0.25000000000000  Cl   (6h)
   0.50800000000000  -0.50800000000000   0.25000000000000  Cl   (6h)
  -0.50800000000000  -1.01600000000000   0.75000000000000  Cl   (6h)
   1.01600000000000   0.50800000000000   0.75000000000000  Cl   (6h)
  -0.50800000000000   0.50800000000000   0.75000000000000  Cl   (6h)
   0.82400000000000   1.64800000000000   0.09200000000000  Cl  (12k)
  -1.64800000000000  -0.82400000000000   0.09200000000000  Cl  (12k)
   0.82400000000000  -0.82400000000000   0.09200000000000  Cl  (12k)
  -0.82400000000000  -1.64800000000000   0.59200000000000  Cl  (12k)
   1.64800000000000   0.82400000000000   0.59200000000000  Cl  (12k)
  -0.82400000000000   0.82400000000000   0.59200000000000  Cl  (12k)
   1.64800000000000   0.82400000000000  -0.09200000000000  Cl  (12k)
  -0.82400000000000  -1.64800000000000  -0.09200000000000  Cl  (12k)
  -0.82400000000000   0.82400000000000  -0.09200000000000  Cl  (12k)
  -1.64800000000000  -0.82400000000000   0.40800000000000  Cl  (12k)
   0.82400000000000   1.64800000000000   0.40800000000000  Cl  (12k)
   0.82400000000000  -0.82400000000000   0.40800000000000  Cl  (12k)
   0.33333333333333   0.66666666666667   0.83700000000000  Cr   (4f)
   0.66666666666667   0.33333333333333   1.33700000000000  Cr   (4f)
   0.66666666666667   0.33333333333333  -0.83700000000000  Cr   (4f)
   0.33333333333333   0.66666666666667  -0.33700000000000  Cr   (4f)
   0.00000000000000   0.00000000000000   0.25000000000000  Cs   (2b)
   0.00000000000000   0.00000000000000   0.75000000000000  Cs   (2b)
   0.33333333333333   0.66666666666667   0.07700000000000  Cs   (4f)
   0.66666666666667   0.33333333333333   0.57700000000000  Cs   (4f)
   0.66666666666667   0.33333333333333  -0.07700000000000  Cs   (4f)
   0.33333333333333   0.66666666666667   0.42300000000000  Cs   (4f)
```

$Na_{0.74}CoO_2$: AB2C2_hP10_194_a_bc_f - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'CoNa0.74O2'
_chemical_formula_sum 'Co Na2 O2'

loop_
_publ_author_name
 'R. J. Balsys'
 'R. L. Davis'
_journal_name_full_name
;
 Solid State Ionics
;
_journal_volume 93
_journal_year 1997
_journal_page_first 279
_journal_page_last 282
_publ_Section_title
;
 Refinement of the structure of Na$_{0.74}$CoO$_{2}$ using neutron
    ↪ powder diffraction
;

# Found in S. Sk, J. Pati, R. S. Dhaka, and. Exploring the possibility
    ↪ of enhancing the figure-of-merit ($> 2$) of Na$_{0.74}$CoO$_{2}
    ↪ $: A combined experimental and theoretical study. Eur. Phys. J.
    ↪ B 93, 155 (2020)

_aflow_title 'Na$_{0.74}$CoO$_{2}$ Structure'
_aflow_proto 'AB2C2_hP10_194_a_bc_f'
_aflow_params 'a,c/a,z_{4}'
_aflow_params_values '2.84,3.80669014085,0.0913'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hP10'

_symmetry_space_group_name_H-M "P 63/m 2/m 2/c"
_symmetry_Int_Tables_number 194

_cell_length_a 2.84000
_cell_length_b 2.84000
_cell_length_c 10.81100
_cell_angle_alpha 90.00000
_cell_angle_beta 90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z+1/2
3 -y,x-y,z
4 -x,-y,z+1/2
5 -x+y,-x,z
6 y,-x+y,z+1/2
7 x-y,-y,-z
8 x,x-y,-z+1/2
9 y,x,-z
10 -x+y,y,-z+1/2
11 -x,-x+y,-z
12 -y,-x,-z+1/2
13 -x,-y,-z
14 -x+y,-x,-z+1/2
15 y,-x+y,-z
16 x,y,-z+1/2
17 x-y,x,-z
18 -y,x-y,-z+1/2
19 -x+y,y,z
20 -x,-x+y,z+1/2
21 -y,-x,z
22 x-y,-y,z+1/2
23 x,x-y,z
24 y,x,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Co1 Co 2 a 0.00000 0.00000 0.00000 1.00000
Na1 Na 2 b 0.00000 0.00000 0.25000 0.21000
Na2 Na 2 c 0.33333 0.66667 0.25000 0.51000
O1 O 4 f 0.33333 0.66667 0.09130 1.00000
```

$Na_{0.74}CoO_2$: AB2C2_hP10_194_a_bc_f - POSCAR

```
AB2C2_hP10_194_a_bc_f & a,c/a,z4 --params=2.84,3.80669014085,0.0913 &
    ↪ P6_{3}/mmc D_{6h}^{4} #194 (abcf) & hP10 & None & CoNa0.74O2 &
    ↪ CoNa0.74O2 & R. J. Balsys and R. L. Davis, Solid State Ion. 93,
    ↪ 279-282 (1997)
   1.00000000000000
   1.42000000000000  -2.45951214674781   0.00000000000000
   1.42000000000000   2.45951214674781   0.00000000000000
   0.00000000000000   0.00000000000000  10.81100000000000
     Co    Na    O
      2     4     4
Direct
   0.00000000000000   0.00000000000000   0.00000000000000  Co   (2a)
   0.00000000000000   0.00000000000000   0.50000000000000  Co   (2a)
   0.00000000000000   0.00000000000000   0.25000000000000  Na   (2b)
   0.00000000000000   0.00000000000000   0.75000000000000  Na   (2b)
   0.33333333333333   0.66666666666667   0.25000000000000  Na   (2c)
   0.66666666666667   0.33333333333333   0.75000000000000  Na   (2c)
   0.33333333333333   0.66666666666667   0.09130000000000   O   (4f)
   0.66666666666667   0.33333333333333   0.59130000000000   O   (4f)
   0.66666666666667   0.33333333333333  -0.09130000000000   O   (4f)
   0.33333333333333   0.66666666666667   0.40870000000000   O   (4f)
```

$EuIn_2P_2$: AB2C2_hP10_194_a_f_f - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'EuIn2P2'
_chemical_formula_sum 'Eu In2 P2'

loop_
_publ_author_name
 'J. Jiang'
 'S. M. Kauzlarich'
_journal_name_full_name
;
 Chemistry of Materials
;
_journal_volume 18
_journal_year 2006
_journal_page_first 435
_journal_page_last 441
_publ_Section_title
;
 Colossal Magnetoresistance in a Rare Earth Zintl Compound with a New
      ↪ Structure Type: EuIn$_{2}$P$_{2}$
;

_aflow_title 'EuIn$_{2}$P$_{2}$ Structure'
_aflow_proto 'AB2C2_hP10_194_a_f_f'
_aflow_params 'a,c/a,z_{2},z_{3}'
_aflow_params_values '4.0829,4.30943691984,0.82845,0.10706'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hP10'

_symmetry_space_group_name_H-M "P 63/m 2/m 2/c"
_symmetry_Int_Tables_number 194

_cell_length_a    4.08290
_cell_length_b    4.08290
_cell_length_c    17.59500
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z+1/2
3 -y,x-y,z
4 -x,-y,z+1/2
5 -x+y,-x,z
6 y,-x+y,z+1/2
7 x-y,-y,-z
8 x,x-y,-z+1/2
9 y,x,-z
10 -x+y,y,-z+1/2
11 -x,-x+y,-z
12 -y,-x,-z+1/2
13 -x,-y,-z
14 -x+y,-x,-z+1/2
15 y,-x+y,-z
16 x,y,-z+1/2
17 x-y,x,-z
18 -y,x-y,-z+1/2
19 -x+y,y,z
20 -x,-x+y,z+1/2
21 -y,-x,z
22 x-y,-y,z+1/2
23 x,x-y,z
24 y,x,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Eu1 Eu 2 a 0.00000 0.00000 0.00000 1.00000
In1 In 4 f 0.33333 0.66667 0.82845 1.00000
P1 P 4 f 0.33333 0.66667 0.10706 1.00000
```

$EuIn_2P_2$: AB2C2_hP10_194_a_f_f - POSCAR

```
AB2C2_hP10_194_a_f_f & a,c/a,z2,z3 --params=4.0829,4.30943691984,0.82845
      ↪ ,0.10706 & P6_{3}/mmc D_{6h}^{4} #194 (af^2) & hP10 & None &
      ↪ EuIn2P2 & EuIn2P2 & J. Jiang and S. M. Kauzlarich, Chem. Mater.
      ↪ 18, 435-441 (2006)
   1.00000000000000
   2.04145000000000  -3.53589512111148   0.00000000000000
   2.04145000000000   3.53589512111148   0.00000000000000
   0.00000000000000   0.00000000000000  17.59500000000000
     Eu    In    P
      2     4     4
Direct
   0.00000000000000   0.00000000000000   0.00000000000000   Eu   (2a)
   0.00000000000000   0.00000000000000   0.50000000000000   Eu   (2a)
   0.33333333333333   0.66666666666667   0.82845000000000   In   (4f)
   0.66666666666667   0.33333333333333   1.32845000000000   In   (4f)
   0.66666666666667   0.33333333333333  -0.82845000000000   In   (4f)
   0.33333333333333   0.66666666666667  -0.32845000000000   In   (4f)
   0.33333333333333   0.66666666666667   0.10706000000000    P   (4f)
   0.66666666666667   0.33333333333333   0.60706000000000    P   (4f)
   0.66666666666667   0.33333333333333  -0.10706000000000    P   (4f)
   0.33333333333333   0.66666666666667   0.39294000000000    P   (4f)
```

Lu_2CoGa_3: AB3C2_hP24_194_f_k_bh - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'CoGa3Lu2'
_chemical_formula_sum 'Co Ga3 Lu2'

loop_
_publ_author_name
 'R. E. Gladyshevskii'
 'K. Cenzual'
 'E. Parth\\'{e}'
_journal_name_full_name
;
 Journal of Alloys and Compounds
;
_journal_volume 189
_journal_year 1992
_journal_page_first 221
_journal_page_last 228
_publ_Section_title
;
 Er$_{2}$RhSi$_{3}$ and $R$$_{2}$CoGa$_{3}$ ($R$ = Y, Tb, Dy, Ho, Er, Tm
      ↪ , Yb) with Lu$_{2}$CoGa$_{3}$ type structure: new members of
      ↪ the AlB$_{2}$ structure family
;

_aflow_title 'Lu$_{2}$CoGa$_{3}$ Structure'
_aflow_proto 'AB3C2_hP24_194_f_k_bh'
_aflow_params 'a,c/a,z_{2},x_{3},x_{4},z_{4}'
_aflow_params_values '8.659,0.787966277861,0.05,0.5231,0.1692,0.0432'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hP24'

_symmetry_space_group_name_H-M "P 63/m 2/m 2/c"
_symmetry_Int_Tables_number 194

_cell_length_a    8.65900
_cell_length_b    8.65900
_cell_length_c    6.82300
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z+1/2
3 -y,x-y,z
4 -x,-y,z+1/2
5 -x+y,-x,z
6 y,-x+y,z+1/2
7 x-y,-y,-z
8 x,x-y,-z+1/2
9 y,x,-z
10 -x+y,y,-z+1/2
11 -x,-x+y,-z
12 -y,-x,-z+1/2
13 -x,-y,-z
14 -x+y,-x,-z+1/2
15 y,-x+y,-z
16 x,y,-z+1/2
17 x-y,x,-z
18 -y,x-y,-z+1/2
19 -x+y,y,z
20 -x,-x+y,z+1/2
21 -y,-x,z
22 x-y,-y,z+1/2
23 x,x-y,z
24 y,x,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Lu1 Lu 2 b 0.00000 0.00000 0.25000 1.00000
Co1 Co 4 f 0.33333 0.66667 0.05000 1.00000
Lu2 Lu 6 h 0.52310 0.04620 0.25000 1.00000
Ga1 Ga 12 k 0.16920 0.33840 0.04320 1.00000
```

Lu_2CoGa_3: AB3C2_hP24_194_f_k_bh - POSCAR

```
AB3C2_hP24_194_f_k_bh & a,c/a,z2,x3,x4,z4 --params=8.659,0.787966277861,
      ↪ 0.05,0.5231,0.1692,0.0432 & P6_{3}/mmc D_{6h}^{4} #194 (bfhk) &
      ↪ hP24 & None & CoGa3Lu2 & CoGa3Lu2 & R. E. Gladyshevskii and K.
      ↪ Cenzual and E. Parth\'{e}, J. Alloys Compd. 189, 221-228 (1992
      ↪ )
   1.00000000000000
   4.32950000000000  -7.49891397136945   0.00000000000000
   4.32950000000000   7.49891397136945   0.00000000000000
   0.00000000000000   0.00000000000000   6.82300000000000
     Co    Ga    Lu
      4    12     8
```

```
Direct
    0.33333333333333    0.66666666666667    0.05000000000000   Co    (4f)
    0.66666666666667    0.33333333333333    0.55000000000000   Co    (4f)
    0.66666666666667    0.33333333333333   -0.05000000000000   Co    (4f)
    0.33333333333333    0.66666666666667    0.45000000000000   Co    (4f)
    0.16920000000000    0.33840000000000    0.04320000000000   Ga   (12k)
   -0.33840000000000   -0.16920000000000    0.04320000000000   Ga   (12k)
    0.16920000000000   -0.16920000000000    0.04320000000000   Ga   (12k)
   -0.16920000000000   -0.33840000000000    0.54320000000000   Ga   (12k)
    0.33840000000000    0.16920000000000    0.54320000000000   Ga   (12k)
   -0.16920000000000    0.16920000000000    0.54320000000000   Ga   (12k)
    0.33840000000000    0.16920000000000   -0.04320000000000   Ga   (12k)
   -0.16920000000000   -0.33840000000000   -0.04320000000000   Ga   (12k)
   -0.16920000000000    0.16920000000000   -0.04320000000000   Ga   (12k)
   -0.33840000000000   -0.16920000000000    0.45680000000000   Ga   (12k)
    0.16920000000000    0.33840000000000    0.45680000000000   Ga   (12k)
    0.16920000000000   -0.16920000000000    0.45680000000000   Ga   (12k)
    0.00000000000000    0.00000000000000    0.25000000000000   Lu    (2b)
    0.00000000000000    0.00000000000000    0.75000000000000   Lu    (2b)
    0.52310000000000    1.04620000000000    0.25000000000000   Lu    (6h)
   -1.04620000000000   -0.52310000000000    0.25000000000000   Lu    (6h)
    0.52310000000000   -0.52310000000000    0.25000000000000   Lu    (6h)
   -0.52310000000000   -1.04620000000000    0.75000000000000   Lu    (6h)
    1.04620000000000    0.52310000000000    0.75000000000000   Lu    (6h)
   -0.52310000000000    0.52310000000000    0.75000000000000   Lu    (6h)
```

Hexagonal Delafossite ($CuAlO_2$): ABC2_hP8_194_a_c_f - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Delafossite'
_chemical_formula_sum 'Al Cu O2'

loop_
_publ_author_name
 'B. U. K\"{o}hler'
 'M. Jansen'
_journal_name_full_name
;
 Zeitschrift fur Anorganische und Allgemeine Chemie
;
_journal_volume 543
_journal_year 1986
_journal_page_first 73
_journal_page_last 80
_publ_Section_title
;
 Darstellung und Strukturdaten von, ``Delafossiten\''\ CuMO$_{2}$ ($M$ =
      ↪ Al, Ga, Sc, Y)
;

# Found in Crystal chemistry and electrical properties of the
      ↪ delafossite structure, 2006

_aflow_title 'Hexagonal Delafossite (CuAlO$_{2}$) Structure'
_aflow_proto 'ABC2_hP8_194_a_c_f'
_aflow_params 'a,c/a,z_{3}'
_aflow_params_values '2.863,3.95179881243,0.0851'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'hP8'

_symmetry_space_group_name_H-M "P 63/m 2/m 2/c"
_symmetry_Int_Tables_number 194

_cell_length_a     2.86300
_cell_length_b     2.86300
_cell_length_c     11.31400
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z+1/2
3 -y,x-y,z
4 -x,-y,z+1/2
5 -x+y,-x,z
6 y,-x+y,z+1/2
7 x-y,-y,-z
8 x,x-y,-z+1/2
9 y,x,-z
10 -x+y,y,-z+1/2
11 -x,-x+y,-z
12 -y,-x,-z+1/2
13 -x,-y,-z
14 -x+y,-x,-z+1/2
15 y,-x+y,-z
16 x,y,-z+1/2
17 x-y,x,-z
18 -y,x-y,-z+1/2
19 -x+y,y,z
20 -x,-x+y,z+1/2
21 -y,-x,z
22 x-y,-y,z+1/2
23 x,x-y,z
24 y,x,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Al1 Al 2 a 0.00000 0.00000 0.00000 1.00000
Cu1 Cu 2 c 0.33333 0.66667 0.25000 1.00000
O1 O 4 f 0.33333 0.66667 0.08510 1.00000
```

Hexagonal Delafossite ($CuAlO_2$): ABC2_hP8_194_a_c_f - POSCAR

```
ABC2_hP8_194_a_c_f & a,c/a,z3 --params=2.863,3.95179881243,0.0851 & P6_{
      ↪ 3}/mmc D_{6h}^{4} #194 (acf) & hP8 & None & AlCuO2 &
      ↪ Delafossite & B. U. K\"{o}hler and M. Jansen, Z. Anorg. Allg.
      ↪ Chem. 543, 73-80 (1986)
   1.00000000000000
   1.43150000000000   -2.47943073103485    0.00000000000000
   1.43150000000000    2.47943073103485    0.00000000000000
   0.00000000000000    0.00000000000000   11.31400000000000
     Al     Cu     O
      2      2      4
Direct
   0.00000000000000    0.00000000000000    0.00000000000000   Al    (2a)
   0.00000000000000    0.00000000000000    0.50000000000000   Al    (2a)
   0.33333333333333    0.66666666666667    0.25000000000000   Cu    (2c)
   0.66666666666667    0.33333333333333    0.75000000000000   Cu    (2c)
   0.33333333333333    0.66666666666667    0.08510000000000   O     (4f)
   0.66666666666667    0.33333333333333    0.58510000000000   O     (4f)
   0.66666666666667    0.33333333333333   -0.08510000000000   O     (4f)
   0.33333333333333    0.66666666666667    0.41490000000000   O     (4f)
```

$LiZn_2$ (C_k): AB_hP4_194_a_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'LiZn2'
_chemical_formula_sum 'Li Zn'

loop_
_publ_author_name
 'E. Zintl'
 'A. Schneider'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Elektrochemie und angewandte physikalische Chemie
;
_journal_volume 41
_journal_year 1935
_journal_page_first 764
_journal_page_last 767
_publ_Section_title
;
 R{\"o}ntgenanalyse der Lithium-Zink-Legierungen (15. Mitteilung {\"u}
      ↪ ber Metalle und Legierungen)
;

# Found in A Handbook of Lattice Spacings and Structures of Metals and
      ↪ Alloys, {N.~R.~C. No. 4303}, 1958

_aflow_title 'LiZn$_{2}$ ($C_{k}$) Structure'
_aflow_proto 'AB_hP4_194_a_c'
_aflow_params 'a,c/a'
_aflow_params_values '4.362,0.575424117377'
_aflow_Strukturbericht '$C_{k}$'
_aflow_Pearson 'hP4'

_symmetry_space_group_name_H-M "P 63/m 2/m 2/c"
_symmetry_Int_Tables_number 194

_cell_length_a     4.36200
_cell_length_b     4.36200
_cell_length_c     2.51000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z+1/2
3 -y,x-y,z
4 -x,-y,z+1/2
5 -x+y,-x,z
6 y,-x+y,z+1/2
7 x-y,-y,-z
8 x,x-y,-z+1/2
9 y,x,-z
10 -x+y,y,-z+1/2
11 -x,-x+y,-z
12 -y,-x,-z+1/2
13 -x,-y,-z
14 -x+y,-x,-z+1/2
15 y,-x+y,-z
16 x,y,-z+1/2
17 x-y,x,-z
18 -y,x-y,-z+1/2
19 -x+y,y,z
20 -x,-x+y,z+1/2
21 -y,-x,z
22 x-y,-y,z+1/2
23 x,x-y,z
24 y,x,z+1/2
```

```
loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Li1 Li 2 a 0.00000 0.00000 0.00000 0.40000
Zn1 Zn 2 c 0.33333 0.66667 0.25000 1.00000
```

$LiZn_2$ (C_k): AB_hP4_194_a_c - POSCAR

```
AB_hP4_194_a_c & a,c/a --params=4.362,0.575424117377 & P6_{3}/mmc D_{6h
    ↪ }^{4} #194 (ac) & hP4 & $C_{k}$ & LiZn2 & LiZn2 & E. Zintl and
    ↪ A. Schneider, Z. Elektrochem. 41, 764-767 (1935)
   1.00000000000000
   2.18100000000000  -3.77760281130772   0.00000000000000
   2.18100000000000   3.77760281130772   0.00000000000000
   0.00000000000000   0.00000000000000   2.51000000000000
    Li   Zn
     2    2
Direct
   0.00000000000000   0.00000000000000   0.00000000000000   Li   (2a)
   0.00000000000000   0.00000000000000   0.50000000000000   Li   (2a)
   0.33333333333333   0.66666666666667   0.25000000000000   Zn   (2c)
   0.66666666666667   0.33333333333333   0.75000000000000   Zn   (2c)
```

Fe_2N (approximate, $L'3_0$): AB_hP4_194_c_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Fe2N'
_chemical_formula_sum 'Fe N'

loop_
_publ_author_name
 'C. J. Smithells'
_journal_year 1955
_publ_Section_title
;
 Metals Reference Book
;

_aflow_title 'Fe$_{2}$N (approximate, $L'3_{0}$) Structure'
_aflow_proto 'AB_hP4_194_c_a'
_aflow_params 'a,c/a'
_aflow_params_values '2.767,1.59522949042'
_aflow_Strukturbericht '$L'3_{0}$'
_aflow_Pearson 'hP4'

_symmetry_space_group_name_H-M "P 63/m 2/m 2/c"
_symmetry_Int_Tables_number 194

_cell_length_a     2.76700
_cell_length_b     2.76700
_cell_length_c     4.41400
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  120.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x-y,x,z+1/2
3 -y,x-y,z
4 -x,-y,z+1/2
5 -x+y,-x,z
6 y,-x+y,z+1/2
7 x-y,-y,-z
8 x,x-y,-z+1/2
9 y,x,-z
10 -x+y,y,-z+1/2
11 -x,-x+y,-z
12 -y,-x,-z+1/2
13 -x,-y,-z
14 -x+y,-x,-z+1/2
15 y,-x+y,-z
16 x,y,-z+1/2
17 x-y,x,-z
18 -y,x-y,-z+1/2
19 -x+y,y,z
20 -x,-x+y,z+1/2
21 -y,-x,z
22 x-y,-y,z+1/2
23 x,x-y,z
24 y,x,z+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
N1 N 2 a 0.00000 0.00000 0.00000 0.50000
Fe1 Fe 2 c 0.33333 0.66667 0.25000 1.00000
```

Fe_2N (approximate, $L'3_0$): AB_hP4_194_c_a - POSCAR

```
AB_hP4_194_c_a & a,c/a --params=2.767,1.59522949042 & P6_{3}/mmc D_{6h
    ↪ }^{4} #194 (ac) & hP4 & $L'3_{0}$ & Fe2N & Fe2N & C. J.
    ↪ Smithells, (1955)
   1.00000000000000
   1.38350000000000  -2.39629229227154   0.00000000000000
   1.38350000000000   2.39629229227154   0.00000000000000
   0.00000000000000   0.00000000000000   4.41400000000000
    Fe   N
     2   2
Direct
   0.33333333333333   0.66666666666667   0.25000000000000   Fe   (2c)
   0.66666666666667   0.33333333333333   0.75000000000000   Fe   (2c)
   0.00000000000000   0.00000000000000   0.00000000000000   N   (2a)
   0.00000000000000   0.00000000000000   0.50000000000000   N   (2a)
```

Cubic Cu_2OSeO_3: A2B4C_cP56_198_ab_2a2b_2a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Cu2O4Se'
_chemical_formula_sum 'Cu2 O4 Se'

loop_
_publ_author_name
 'H. Effenberger'
 'F. Pertlik'
_journal_name_full_name
;
 Monatshefte f{\"u}r Chemie - Chemical Monthly
;
_journal_volume 117
_journal_year 1986
_journal_page_first 887
_journal_page_last 896
_publ_Section_title
;
 Die Kristallstrukturen der Kupfer(II)-oxo-selenite Cu$_{2}$O(SeO$_{3}$)
    ↪ (kubisch und monoklin) und Cu$_{4}$O(SeO$_{3}$)$_{3}$ (
    ↪ monoklin und triklin)
;

# Found in Magnon spectrum of the helimagnetic insulator Cu$_{2}$OSeO$_{
    ↪ 3}$, 2016

_aflow_title 'Cubic Cu$_{2}$OSeO$_{3}$ Structure'
_aflow_proto 'A2B4C_cP56_198_ab_2a2b_2a'
_aflow_params 'a,x_{1},x_{2},x_{3},x_{4},x_{5},x_{6},y_{6},z_{6},x_{7},
    ↪ y_{7},z_{7},x_{8},y_{8},z_{8}'
_aflow_params_values '8.925,0.886,0.0105,0.7621,0.459,0.2113,0.1335,
    ↪ 0.1211,0.8719,0.2699,0.4834,0.4706,0.271,0.1892,0.0313'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'cP56'

_symmetry_space_group_name_H-M "P 21 3"
_symmetry_Int_Tables_number 198

_cell_length_a     8.92500
_cell_length_b     8.92500
_cell_length_c     8.92500
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z
3 -x,y+1/2,-z+1/2
4 -x+1/2,-y,z+1/2
5 y,z,x
6 y+1/2,-z+1/2,-x
7 -y,z+1/2,-x+1/2
8 -y+1/2,-z,x+1/2
9 z,x,y
10 z+1/2,-x+1/2,-y
11 -z,x+1/2,-y+1/2
12 -z+1/2,-x,y+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cu1 Cu 4 a 0.88600 0.88600 0.88600 1.00000
O1 O 4 a 0.01050 0.01050 0.01050 1.00000
O2 O 4 a 0.76210 0.76210 0.76210 1.00000
Se1 Se 4 a 0.45900 0.45900 0.45900 1.00000
Se2 Se 4 a 0.21130 0.21130 0.21130 1.00000
Cu2 Cu 12 b 0.13350 0.12110 0.87190 1.00000
O3 O 12 b 0.26990 0.48340 0.47060 1.00000
O4 O 12 b 0.27100 0.18920 0.03130 1.00000
```

Cubic Cu_2OSeO_3: A2B4C_cP56_198_ab_2a2b_2a - POSCAR

```
A2B4C_cP56_198_ab_2a2b_2a & a,x1,x2,x3,x4,x5,x6,y6,z6,x7,y7,z7,x8,y8,z8
    ↪ --params=8.925,0.886,0.0105,0.7621,0.459,0.2113,0.1335,0.1211,
    ↪ 0.8719,0.2699,0.4834,0.4706,0.271,0.1892,0.0313 & P2_{1}3 T^{4}
    ↪ #198 (a^5b^3) & cP56 & None & Cu2O4Se & Cu2O4Se & H.
    ↪ Effenberger and F. Pertlik, Monatshefte f{"u}r Chemie -
    ↪ Chemical Monthly 117, 887-896 (1986)
   1.00000000000000
   8.92500000000000    0.00000000000000    0.00000000000000
   0.00000000000000    8.92500000000000    0.00000000000000
   0.00000000000000    0.00000000000000    8.92500000000000
    Cu     O     Se
    16    32     8
Direct
   0.88600000000000    0.88600000000000    0.88600000000000  Cu   (4a)
  -0.38600000000000   -0.88600000000000    1.38600000000000  Cu   (4a)
  -0.88600000000000    1.38600000000000   -0.38600000000000  Cu   (4a)
   1.38600000000000   -0.38600000000000   -0.88600000000000  Cu   (4a)
   0.13350000000000    0.12110000000000    0.87190000000000  Cu  (12b)
   0.36650000000000   -0.12110000000000    1.37190000000000  Cu  (12b)
  -0.13350000000000    0.62110000000000   -0.37190000000000  Cu  (12b)
   0.63350000000000    0.37890000000000   -0.87190000000000  Cu  (12b)
   0.87190000000000    0.13350000000000    0.12110000000000  Cu  (12b)
   1.37190000000000    0.36650000000000   -0.12110000000000  Cu  (12b)
  -0.37190000000000   -0.13350000000000    0.62110000000000  Cu  (12b)
  -0.87190000000000    0.63350000000000    0.37890000000000  Cu  (12b)
   0.12110000000000    0.87190000000000    0.13350000000000  Cu  (12b)
  -0.12110000000000    1.37190000000000    0.36650000000000  Cu  (12b)
   0.62110000000000   -0.37190000000000   -0.13350000000000  Cu  (12b)
   0.37890000000000   -0.87190000000000    0.63350000000000  Cu  (12b)
   0.01050000000000    0.01050000000000    0.01050000000000   O   (4a)
   0.48950000000000   -0.01050000000000    0.51050000000000   O   (4a)
  -0.01050000000000    0.51050000000000    0.48950000000000   O   (4a)
   0.51050000000000    0.48950000000000   -0.01050000000000   O   (4a)
   0.76210000000000    0.76210000000000    0.76210000000000   O   (4a)
  -0.26210000000000   -0.76210000000000    1.26210000000000   O   (4a)
  -0.76210000000000    1.26210000000000   -0.26210000000000   O   (4a)
   1.26210000000000   -0.26210000000000   -0.76210000000000   O   (4a)
   0.26990000000000    0.48340000000000    0.47060000000000   O  (12b)
   0.23010000000000   -0.48340000000000    0.97060000000000   O  (12b)
  -0.26990000000000    0.98340000000000    0.02940000000000   O  (12b)
   0.76990000000000    0.01660000000000   -0.47060000000000   O  (12b)
   0.47060000000000    0.26990000000000    0.48340000000000   O  (12b)
   0.97060000000000    0.23010000000000   -0.48340000000000   O  (12b)
   0.02940000000000   -0.26990000000000    0.98340000000000   O  (12b)
  -0.47060000000000    0.76990000000000    0.01660000000000   O  (12b)
   0.48340000000000    0.47060000000000    0.26990000000000   O  (12b)
  -0.48340000000000    0.97060000000000    0.23010000000000   O  (12b)
   0.98340000000000    0.02940000000000   -0.26990000000000   O  (12b)
   0.01660000000000   -0.47060000000000    0.76990000000000   O  (12b)
   0.27100000000000    0.18920000000000    0.03130000000000   O  (12b)
   0.22900000000000   -0.18920000000000    0.53130000000000   O  (12b)
  -0.27100000000000    0.68920000000000    0.46870000000000   O  (12b)
   0.77100000000000    0.31080000000000   -0.03130000000000   O  (12b)
   0.03130000000000    0.27100000000000    0.18920000000000   O  (12b)
   0.53130000000000    0.22900000000000   -0.18920000000000   O  (12b)
   0.46870000000000   -0.27100000000000    0.68920000000000   O  (12b)
  -0.03130000000000    0.77100000000000    0.31080000000000   O  (12b)
   0.18920000000000    0.03130000000000    0.27100000000000   O  (12b)
  -0.18920000000000    0.53130000000000    0.22900000000000   O  (12b)
   0.68920000000000    0.46870000000000   -0.27100000000000   O  (12b)
   0.31080000000000   -0.03130000000000    0.77100000000000   O  (12b)
   0.45900000000000    0.45900000000000    0.45900000000000  Se   (4a)
   0.04100000000000   -0.45900000000000    0.95900000000000  Se   (4a)
  -0.45900000000000    0.95900000000000    0.04100000000000  Se   (4a)
   0.95900000000000    0.04100000000000   -0.45900000000000  Se   (4a)
   0.21130000000000    0.21130000000000    0.21130000000000  Se   (4a)
   0.28870000000000   -0.21130000000000    0.71130000000000  Se   (4a)
  -0.21130000000000    0.71130000000000    0.28870000000000  Se   (4a)
   0.71130000000000    0.28870000000000   -0.21130000000000  Se   (4a)
```

Na_2CaSiO_4 ($S6_6$): AB2C4D_cP32_198_a_2a_ab_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'CaNa2O4Si'
_chemical_formula_sum 'Ca Na2 O4 Si'

loop_
_publ_author_name
 'T. F. W. Barth'
 'E. Posnjak'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 81
_journal_year 1932
_journal_page_first 370
_journal_page_last 375
_publ_Section_title
;
 Silicate structures of the cristobalite type: II. The crystal structure
    ↪ of Na$_{2}$CaSiO$_{4}$
;

_aflow_title 'Na$_{2}$CaSiO$_{4}$ ($S6_{6}$) Structure'
_aflow_proto 'AB2C4D_cP32_198_a_2a_ab_a'
_aflow_params 'a,x_{1},x_{2},x_{3},x_{4},x_{5},x_{6},y_{6},z_{6}'
_aflow_params_values '7.48,-0.007,0.5,0.75,0.133,0.253,0.556,0.667,0.222
    ↪ '
_aflow_Strukturbericht '$S6_{6}$'
_aflow_Pearson 'cP32'

_symmetry_space_group_name_H-M "P 21 3"
_symmetry_Int_Tables_number 198

_cell_length_a    7.48000
_cell_length_b    7.48000
_cell_length_c    7.48000
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z
3 -x,y+1/2,-z+1/2
4 -x+1/2,-y,z+1/2
5 y,z,x
6 y+1/2,-z+1/2,-x
7 -y,z+1/2,-x+1/2
8 -y+1/2,-z,x+1/2
9 z,x,y
10 z+1/2,-x+1/2,-y
11 -z,x+1/2,-y+1/2
12 -z+1/2,-x,y+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ca1 Ca 4 a -0.00700 -0.00700 -0.00700 1.00000
Na1 Na 4 a 0.50000 0.50000 0.50000 1.00000
Na2 Na 4 a 0.75000 0.75000 0.75000 1.00000
O1 O 4 a 0.13300 0.13300 0.13300 1.00000
Si1 Si 4 a 0.25300 0.25300 0.25300 1.00000
O2 O 12 b 0.55600 0.66700 0.22200 1.00000
```

Na_2CaSiO_4 ($S6_6$): AB2C4D_cP32_198_a_2a_ab_a - POSCAR

```
AB2C4D_cP32_198_a_2a_ab_a & a,x1,x2,x3,x4,x5,x6,y6,z6 --params=7.48,-
    ↪ 0.007,0.5,0.75,0.133,0.253,0.556,0.667,0.222 & P2_{1}3 T^{4} #
    ↪ 198 (a^5b) & cP32 & $S6_{6}$ & CaNa2O4Si & CaNa2O4Si & T. F. W.
    ↪ Barth and E. Posnjak, Zeitschrift f{"u}r Kristallographie -
    ↪ Crystalline Materials 81, 370-375 (1932)
   1.00000000000000
   7.48000000000000    0.00000000000000    0.00000000000000
   0.00000000000000    7.48000000000000    0.00000000000000
   0.00000000000000    0.00000000000000    7.48000000000000
    Ca     Na     O     Si
    4      8     16     4
Direct
  -0.00700000000000   -0.00700000000000   -0.00700000000000  Ca   (4a)
   0.50700000000000    0.00700000000000    0.49300000000000  Ca   (4a)
   0.00700000000000    0.49300000000000    0.50700000000000  Ca   (4a)
   0.49300000000000    0.50700000000000    0.00700000000000  Ca   (4a)
   0.50000000000000    0.50000000000000    0.50000000000000  Na   (4a)
   0.00000000000000   -0.50000000000000    1.00000000000000  Na   (4a)
  -0.50000000000000    1.00000000000000    0.00000000000000  Na   (4a)
   1.00000000000000    0.00000000000000   -0.50000000000000  Na   (4a)
   0.75000000000000    0.75000000000000    0.75000000000000  Na   (4a)
  -0.25000000000000   -0.75000000000000    1.25000000000000  Na   (4a)
  -0.75000000000000    1.25000000000000   -0.25000000000000  Na   (4a)
   1.25000000000000   -0.25000000000000   -0.75000000000000  Na   (4a)
   0.13300000000000    0.13300000000000    0.13300000000000   O   (4a)
   0.36700000000000   -0.13300000000000    0.63300000000000   O   (4a)
  -0.13300000000000    0.63300000000000    0.36700000000000   O   (4a)
   0.63300000000000    0.36700000000000   -0.13300000000000   O   (4a)
   0.55600000000000    0.66700000000000    0.22200000000000   O  (12b)
  -0.05600000000000   -0.66700000000000    0.72200000000000   O  (12b)
  -0.55600000000000    1.16700000000000    0.27800000000000   O  (12b)
   1.05600000000000   -0.16700000000000   -0.22200000000000   O  (12b)
   0.22200000000000    0.55600000000000    0.66700000000000   O  (12b)
   0.72200000000000   -0.05600000000000   -0.66700000000000   O  (12b)
   0.27800000000000   -0.55600000000000    1.16700000000000   O  (12b)
  -0.22200000000000    1.05600000000000   -0.16700000000000   O  (12b)
   0.66700000000000    0.22200000000000    0.55600000000000   O  (12b)
  -0.66700000000000    0.72200000000000   -0.05600000000000   O  (12b)
   1.16700000000000    0.27800000000000   -0.55600000000000   O  (12b)
  -0.16700000000000   -0.22200000000000    1.05600000000000   O  (12b)
   0.25300000000000    0.25300000000000    0.25300000000000  Si   (4a)
   0.24700000000000   -0.25300000000000    0.75300000000000  Si   (4a)
  -0.25300000000000    0.75300000000000    0.24700000000000  Si   (4a)
   0.75300000000000    0.24700000000000   -0.25300000000000  Si   (4a)
```

α-Carnegieite ($NaAlSiO_4$, $S6_5$): ABC4D_cP28_198_a_a_ab_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral '$\alpha$-carnegieite'
_chemical_formula_sum 'Al Na O4 Si'

loop_
_publ_author_name
 'T. F. W. Barth'
 'E. Posnjak'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
```

```
;
_journal_volume 81
_journal_year 1932
_journal_page_first 135
_journal_page_last 141
_publ_Section_title
;
 Silicate structures of the cristobalite type: I. The crystal structure
      ↪ of $\alpha$-carnegieite (NaAlSiO$_{4}$)
;

_aflow_title '$\alpha$-Carnegieite (NaAlSiO$_{4}$, $S6_{5}$) Structure'
_aflow_proto 'ABC4D_cP28_198_a_a_ab_a'
_aflow_params 'a,x_{1},x_{2},x_{3},x_{4},x_{5},y_{5},z_{5}'
_aflow_params_values '7.37,0.258,0.744,0.125,0.0,0.658,0.644,0.0556'
_aflow_Strukturbericht '$S6_{5}$'
_aflow_Pearson 'cP28'

_symmetry_space_group_name_H-M "P 21 3"
_symmetry_Int_Tables_number 198

_cell_length_a     7.37000
_cell_length_b     7.37000
_cell_length_c     7.37000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z
3 -x,y+1/2,-z+1/2
4 -x+1/2,-y,z+1/2
5 y,z,x
6 y+1/2,-z+1/2,-x
7 -y,z+1/2,-x+1/2
8 -y+1/2,-z,x+1/2
9 z,x,y
10 z+1/2,-x+1/2,-y
11 -z,x+1/2,-y+1/2
12 -z+1/2,-x,y+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Al1 Al 4 a 0.25800 0.25800 0.25800 1.00000
Na1 Na 4 a 0.74400 0.74400 0.74400 1.00000
O1 O 4 a 0.12500 0.12500 0.12500 1.00000
Si1 Si 4 a 0.00000 0.00000 0.00000 1.00000
O2 O 12 b 0.65800 0.64400 0.05560 1.00000
```

α-Carnegieite ($NaAlSiO_4$, $S6_5$): ABC4D_cP28_198_a_a_ab_a - POSCAR

```
ABC4D_cP28_198_a_a_ab_a & a,x1,x2,x3,x4,x5,y5,z5 --params=7.37,0.258,
      ↪ 0.744,0.125,0.0,0.658,0.644,0.0556 & P2_{1}3 T^{4} #198 (a^4b)
      ↪ & cP28 & $S6_{5}$ & AlNaO4Si & $\alpha$-carnegieite & T. F. W.
      ↪ Barth and E. Posnjak, Zeitschrift f{"u}r Kristallographie -
      ↪ Crystalline Materials 81, 135-141 (1932)
   1.00000000000000
   7.37000000000000    0.00000000000000    0.00000000000000
   0.00000000000000    7.37000000000000    0.00000000000000
   0.00000000000000    0.00000000000000    7.37000000000000
     Al     Na      O     Si
      4      4     16      4
Direct
   0.25800000000000    0.25800000000000    0.25800000000000   Al   (4a)
   0.24200000000000   -0.25800000000000    0.75800000000000   Al   (4a)
  -0.25800000000000    0.75800000000000    0.24200000000000   Al   (4a)
   0.75800000000000    0.24200000000000   -0.25800000000000   Al   (4a)
   0.74400000000000    0.74400000000000    0.74400000000000   Na   (4a)
  -0.24400000000000   -0.74400000000000    1.24400000000000   Na   (4a)
  -0.74400000000000    1.24400000000000   -0.24400000000000   Na   (4a)
   1.24400000000000   -0.24400000000000   -0.74400000000000   Na   (4a)
   0.12500000000000    0.12500000000000    0.12500000000000   O    (4a)
   0.37500000000000   -0.12500000000000    0.62500000000000   O    (4a)
  -0.12500000000000    0.62500000000000    0.37500000000000   O    (4a)
   0.62500000000000    0.37500000000000   -0.12500000000000   O    (4a)
   0.65800000000000    0.64400000000000    0.05560000000000   O   (12b)
  -0.15800000000000   -0.64400000000000    0.55560000000000   O   (12b)
  -0.65800000000000    1.14400000000000    0.44440000000000   O   (12b)
   1.15800000000000   -0.14400000000000   -0.05560000000000   O   (12b)
   0.05560000000000    0.65800000000000    0.64400000000000   O   (12b)
   0.55560000000000   -0.15800000000000   -0.64400000000000   O   (12b)
   0.44440000000000   -0.65800000000000    1.14400000000000   O   (12b)
  -0.05560000000000    1.15800000000000   -0.14400000000000   O   (12b)
   0.64400000000000    0.05560000000000    0.65800000000000   O   (12b)
  -0.64400000000000    0.55560000000000   -0.15800000000000   O   (12b)
   1.14400000000000    0.44440000000000   -0.65800000000000   O   (12b)
  -0.14400000000000   -0.05560000000000    1.15800000000000   O   (12b)
   0.00000000000000    0.00000000000000    0.00000000000000   Si   (4a)
   0.50000000000000    0.00000000000000    0.50000000000000   Si   (4a)
   0.00000000000000    0.50000000000000    0.50000000000000   Si   (4a)
   0.50000000000000    0.50000000000000    0.00000000000000   Si   (4a)
```

$C26_a$ (NO_2) (*obsolete*): AB2_cI36_199_b_c - CIF

```
# CIF file

data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'NO2'
_chemical_formula_sum 'N O2'

loop_
_publ_author_name
 'L. Vegard'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Physik
;
_journal_volume 68
_journal_year 1931
_journal_page_first 184
_journal_page_last 203
_publ_Section_title
;
 Die Struktur von festem N$_{2}$O$_{4}$ bei der Temperatur von fl\"{u}
      ↪ ssiger Luft
;

# Found in Strukturbericht Band II 1928-1932, 1937

_aflow_title '$C26_{a}$ (NO$_{2}$) ({\em{obsolete}}) Structure'
_aflow_proto 'AB2_cI36_199_b_c'
_aflow_params 'a,x_{1},x_{2},y_{2},z_{2}'
_aflow_params_values '7.77,0.4,0.178,0.25,0.403'
_aflow_Strukturbericht '$C26_{a}$'
_aflow_Pearson 'cI36'

_symmetry_space_group_name_H-M "I 21 3"
_symmetry_Int_Tables_number 199

_cell_length_a     7.77000
_cell_length_b     7.77000
_cell_length_c     7.77000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z+1/2
3 -x+1/2,y,-z
4 -x,-y+1/2,z
5 y,z,x
6 y,-z,-x+1/2
7 -y+1/2,z,-x
8 -y,-z+1/2,x
9 z,x,y
10 z,-x,-y+1/2
11 -z+1/2,x,-y
12 -z,-x+1/2,y
13 x+1/2,y+1/2,z+1/2
14 x+1/2,-y+1/2,-z
15 -x,y+1/2,-z+1/2
16 -x+1/2,-y,z+1/2
17 y+1/2,z+1/2,x+1/2
18 y+1/2,-z+1/2,-x
19 -y,z+1/2,-x+1/2
20 -y+1/2,-z,x+1/2
21 z+1/2,x+1/2,y+1/2
22 z+1/2,-x+1/2,-y
23 -z,x+1/2,-y+1/2
24 -z+1/2,-x,y+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
N1 N 12 b 0.40000 0.00000 0.25000 1.00000
O1 O 24 c 0.17800 0.25000 0.40300 1.00000
```

$C26_a$ (NO_2) (*obsolete*): AB2_cI36_199_b_c - POSCAR

```
AB2_cI36_199_b_c & a,x1,x2,y2,z2 --params=7.77,0.4,0.178,0.25,0.403 &
      ↪ I2_{1}3 T^{5} #199 (bc) & cI36 & $C26_{a}$ & NO2 & NO2 & L.
      ↪ Vegard, Z. Phys. 68, 184-203 (1931)
   1.00000000000000
  -3.88500000000000    3.88500000000000    3.88500000000000
   3.88500000000000   -3.88500000000000    3.88500000000000
   3.88500000000000    3.88500000000000   -3.88500000000000
     N      O
     6     12
Direct
   0.25000000000000    0.65000000000000    0.40000000000000   N   (12b)
   0.75000000000000   -0.15000000000000    0.10000000000000   N   (12b)
   0.40000000000000    0.25000000000000    0.65000000000000   N   (12b)
   0.10000000000000    0.75000000000000   -0.15000000000000   N   (12b)
   0.65000000000000    0.40000000000000    0.25000000000000   N   (12b)
  -0.15000000000000    0.10000000000000    0.75000000000000   N   (12b)
   0.65300000000000    0.58100000000000    0.42800000000000   O   (24c)
   0.65300000000000    0.22500000000000    0.07200000000000   O   (24c)
  -0.15300000000000   -0.08100000000000    0.57200000000000   O   (24c)
  -0.15300000000000    0.27500000000000   -0.07200000000000   O   (24c)
   0.42800000000000    0.65300000000000    0.58100000000000   O   (24c)
```

```
   0.07200000000000   0.65300000000000   0.22500000000000   O  (24c)
   0.57200000000000  -0.15300000000000  -0.08100000000000   O  (24c)
  -0.07200000000000  -0.15300000000000   0.27500000000000   O  (24c)
   0.58100000000000   0.42800000000000   0.65300000000000   O  (24c)
   0.22500000000000   0.07200000000000   0.65300000000000   O  (24c)
  -0.08100000000000   0.57200000000000  -0.15300000000000   O  (24c)
   0.27500000000000  -0.07200000000000  -0.15300000000000   O  (24c)
```

$Bi_3Ru_3O_{11}$: A3B11C3_cP68_201_be_efh_g - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Bi3O11Ru3'
_chemical_formula_sum 'Bi3 O11 Ru3'

loop_
_publ_author_name
 'F. Abraham'
 'D. Thomas'
 'G. Nowogrocki'
_journal_name_full_name
;
 Bulletin de la Societ{\'e} fran{\c{c}}aise de Mineralogie et de
     ↪ Crystallographie
;
_journal_volume 98
_journal_year 1975
_journal_page_first 25
_journal_page_last 29
_publ_Section_title
;
 Structure cristalline de Bi$_{3}$Ru$_{3}$O$_{11}$
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'Bi$_{3}$Ru$_{3}$O$_{11}$ Structure'
_aflow_proto 'A3B11C3_cP68_201_be_efh_g'
_aflow_params 'a,x_{2},x_{3},x_{4},x_{5},x_{6},y_{6},z_{6}'
_aflow_params_values '9.302,0.38379,0.152,0.59,0.3897,0.599,0.247,0.547'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'cP68'

_symmetry_space_group_name_H-M "P 2/n -3 (origin choice 2)"
_symmetry_Int_Tables_number 201

_cell_length_a    9.30200
_cell_length_b    9.30200
_cell_length_c    9.30200
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y+1/2,-z+1/2
3 -x+1/2,y,-z+1/2
4 -x+1/2,-y+1/2,z
5 y,z,x
6 y,-z+1/2,-x+1/2
7 -y+1/2,z,-x+1/2
8 -y+1/2,-z+1/2,x
9 z,x,y
10 z,-x+1/2,-y+1/2
11 -z+1/2,x,-y+1/2
12 -z+1/2,-x+1/2,y
13 -x,-y,-z
14 -x,y+1/2,z+1/2
15 x+1/2,-y,z+1/2
16 x+1/2,y+1/2,-z
17 -y,-z,-x
18 -y,z+1/2,x+1/2
19 y+1/2,-z,x+1/2
20 y+1/2,z+1/2,-x
21 -z,-x,-y
22 -z,x+1/2,y+1/2
23 z+1/2,-x,y+1/2
24 z+1/2,x+1/2,-y

loop_
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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Bi1 Bi 4 b 0.00000 0.00000 0.00000 1.00000
Bi2 Bi 8 e 0.38379 0.38379 0.38379 1.00000
O1 O 8 e 0.15200 0.15200 0.15200 1.00000
O2 O 12 f 0.59000 0.25000 0.25000 1.00000
Ru1 Ru 12 g 0.38970 0.75000 0.25000 1.00000
O3 O 24 h 0.59900 0.24700 0.54700 1.00000
```

$Bi_3Ru_3O_{11}$: A3B11C3_cP68_201_be_efh_g - POSCAR

```
A3B11C3_cP68_201_be_efh_g & a,x2,x3,x4,x5,x6,y6,z6 --params=9.302,
    ↪ 0.38379,0.152,0.59,0.3897,0.599,0.247,0.547 & Pn-3 T_{h}^{2} #
    ↪ 201 (be^2fgh) & cP68 & None & Bi3O11Ru3 & Bi3O11Ru3 & F.
    ↪ Abraham and D. Thomas and G. Nowogrocki, Bull. Soc. fr. Min'
    ↪ eral. Crystallogr. 98, 25-29 (1975)
  1.00000000000000
  9.30200000000000   0.00000000000000   0.00000000000000
  0.00000000000000   9.30200000000000   0.00000000000000
  0.00000000000000   0.00000000000000   9.30200000000000
   Bi    O    Ru
   12   44   12
Direct
  0.00000000000000   0.00000000000000   0.00000000000000   Bi   (4b)
  0.50000000000000   0.50000000000000   0.00000000000000   Bi   (4b)
  0.50000000000000   0.00000000000000   0.50000000000000   Bi   (4b)
  0.00000000000000   0.50000000000000   0.50000000000000   Bi   (4b)
  0.38379000000000   0.38379000000000   0.38379000000000   Bi   (8e)
  0.11621000000000   0.11621000000000   0.38379000000000   Bi   (8e)
  0.11621000000000   0.38379000000000   0.11621000000000   Bi   (8e)
  0.38379000000000   0.11621000000000   0.11621000000000   Bi   (8e)
 -0.38379000000000  -0.38379000000000  -0.38379000000000   Bi   (8e)
  0.88379000000000   0.88379000000000  -0.38379000000000   Bi   (8e)
  0.88379000000000  -0.38379000000000   0.88379000000000   Bi   (8e)
 -0.38379000000000   0.88379000000000   0.88379000000000   Bi   (8e)
  0.15200000000000   0.15200000000000   0.15200000000000   O    (8e)
  0.34800000000000   0.34800000000000   0.15200000000000   O    (8e)
  0.34800000000000   0.15200000000000   0.34800000000000   O    (8e)
  0.15200000000000   0.34800000000000   0.34800000000000   O    (8e)
 -0.15200000000000  -0.15200000000000  -0.15200000000000   O    (8e)
  0.65200000000000   0.65200000000000  -0.15200000000000   O    (8e)
  0.65200000000000  -0.15200000000000   0.65200000000000   O    (8e)
 -0.15200000000000   0.65200000000000   0.65200000000000   O    (8e)
  0.59000000000000   0.25000000000000   0.25000000000000   O   (12f)
 -0.09000000000000   0.25000000000000   0.25000000000000   O   (12f)
  0.25000000000000   0.59000000000000   0.25000000000000   O   (12f)
  0.25000000000000  -0.09000000000000   0.25000000000000   O   (12f)
  0.25000000000000   0.25000000000000   0.59000000000000   O   (12f)
  0.25000000000000   0.25000000000000  -0.09000000000000   O   (12f)
 -0.59000000000000   0.75000000000000   0.75000000000000   O   (12f)
  1.09000000000000   0.75000000000000   0.75000000000000   O   (12f)
  0.75000000000000  -0.59000000000000   0.75000000000000   O   (12f)
  0.75000000000000   1.09000000000000   0.75000000000000   O   (12f)
  0.75000000000000   0.75000000000000  -0.59000000000000   O   (12f)
  0.75000000000000   0.75000000000000   1.09000000000000   O   (12f)
  0.59900000000000   0.24700000000000   0.54700000000000   O   (24h)
 -0.09900000000000   0.25300000000000   0.54700000000000   O   (24h)
 -0.09900000000000   0.24700000000000  -0.04700000000000   O   (24h)
  0.59900000000000   0.25300000000000  -0.04700000000000   O   (24h)
  0.54700000000000   0.59900000000000   0.24700000000000   O   (24h)
  0.54700000000000  -0.09900000000000   0.25300000000000   O   (24h)
 -0.04700000000000  -0.09900000000000   0.24700000000000   O   (24h)
 -0.04700000000000   0.59900000000000   0.25300000000000   O   (24h)
  0.24700000000000   0.54700000000000   0.59900000000000   O   (24h)
  0.25300000000000   0.54700000000000  -0.09900000000000   O   (24h)
  0.24700000000000  -0.04700000000000  -0.09900000000000   O   (24h)
  0.25300000000000  -0.04700000000000   0.59900000000000   O   (24h)
 -0.59900000000000  -0.24700000000000  -0.54700000000000   O   (24h)
  1.09900000000000   0.74700000000000  -0.54700000000000   O   (24h)
  1.09900000000000  -0.24700000000000   1.04700000000000   O   (24h)
 -0.59900000000000   0.74700000000000   1.04700000000000   O   (24h)
 -0.54700000000000  -0.59900000000000  -0.24700000000000   O   (24h)
 -0.54700000000000   1.09900000000000   0.74700000000000   O   (24h)
  1.04700000000000   1.09900000000000  -0.24700000000000   O   (24h)
  1.04700000000000  -0.59900000000000   0.74700000000000   O   (24h)
 -0.24700000000000  -0.54700000000000  -0.59900000000000   O   (24h)
  0.74700000000000  -0.54700000000000   1.09900000000000   O   (24h)
 -0.24700000000000   1.04700000000000   1.09900000000000   O   (24h)
  0.74700000000000   1.04700000000000  -0.59900000000000   O   (24h)
  0.38970000000000   0.75000000000000   0.25000000000000   Ru  (12g)
  0.11030000000000   0.75000000000000   0.25000000000000   Ru  (12g)
  0.25000000000000   0.38970000000000   0.75000000000000   Ru  (12g)
  0.25000000000000   0.11030000000000   0.75000000000000   Ru  (12g)
  0.75000000000000   0.25000000000000   0.38970000000000   Ru  (12g)
  0.75000000000000   0.25000000000000   0.11030000000000   Ru  (12g)
 -0.38970000000000   0.25000000000000   0.75000000000000   Ru  (12g)
  0.88970000000000   0.25000000000000   0.75000000000000   Ru  (12g)
  0.75000000000000  -0.38970000000000   0.25000000000000   Ru  (12g)
  0.75000000000000   0.88970000000000   0.25000000000000   Ru  (12g)
  0.25000000000000   0.75000000000000  -0.38970000000000   Ru  (12g)
  0.25000000000000   0.75000000000000   0.88970000000000   Ru  (12g)
```

$K_3Co(NO_2)_6$ ($J2_4$): AB3C6D12_cF88_202_a_bc_e_h - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'CoK3N6O12'
_chemical_formula_sum 'Co K3 N6 O12'

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_publ_author_name
 'M. {van Driel}'
 'H. J. Verweel'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 95
_journal_year 1936
_journal_page_first 308
_journal_page_last 314
_publ_Section_title
;
 \"{U}ber die Struktur der Tripelnitrite
;

# Found in Strukturbericht Band IV 1936, 1938
```

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_aflow_title 'K$_{3}$Co(NO$_{2}$)$_{6}$ ($J2_{4}$) Structure'
_aflow_proto 'AB3C6D12_cF88_202_a_bc_e_h'
_aflow_params 'a,x_{4},y_{5},z_{5}'
_aflow_params_values '10.46,0.195,0.235,0.1'
_aflow_Strukturbericht '$J2_{4}$'
_aflow_Pearson 'cF88'

_symmetry_space_group_name_H-M "F 2/m -3"
_symmetry_Int_Tables_number 202

_cell_length_a      10.46000
_cell_length_b      10.46000
_cell_length_c      10.46000
_cell_angle_alpha   90.00000
_cell_angle_beta    90.00000
_cell_angle_gamma   90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 y,z,x
6 y,-z,-x
7 -y,z,-x
8 -y,-z,x
9 z,x,y
10 z,-x,-y
11 -z,x,-y
12 -z,-x,y
13 -x,-y,-z
14 -x,y,z
15 x,-y,z
16 x,y,-z
17 -y,-z,-x
18 -y,z,x
19 y,-z,x
20 y,z,-x
21 -z,-x,-y
22 -z,x,y
23 z,-x,y
24 z,x,-y
25 x,y+1/2,z+1/2
26 x,-y+1/2,-z+1/2
27 -x,y+1/2,-z+1/2
28 -x,-y+1/2,z+1/2
29 y,z+1/2,x+1/2
30 y,-z+1/2,-x+1/2
31 -y,z+1/2,-x+1/2
32 -y,-z+1/2,x+1/2
33 z,x+1/2,y+1/2
34 z,-x+1/2,-y+1/2
35 -z,x+1/2,-y+1/2
36 -z,-x+1/2,y+1/2
37 -x,-y+1/2,-z+1/2
38 -x,y+1/2,z+1/2
39 x,-y+1/2,z+1/2
40 x,y+1/2,-z+1/2
41 -y,-z+1/2,-x+1/2
42 -y,z+1/2,x+1/2
43 y,-z+1/2,x+1/2
44 y,z+1/2,-x+1/2
45 -z,-x+1/2,-y+1/2
46 -z,x+1/2,y+1/2
47 z,-x+1/2,y+1/2
48 z,x+1/2,-y+1/2
49 x+1/2,y,z+1/2
50 x+1/2,-y,-z+1/2
51 -x+1/2,y,-z+1/2
52 -x+1/2,-y,z+1/2
53 y+1/2,z,x+1/2
54 y+1/2,-z,-x+1/2
55 -y+1/2,z,-x+1/2
56 -y+1/2,-z,x+1/2
57 z+1/2,x,y+1/2
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61 -x+1/2,-y,-z+1/2
62 -x+1/2,y,z+1/2
63 x+1/2,-y,z+1/2
64 x+1/2,y,-z+1/2
65 -y+1/2,-z,-x+1/2
66 -y+1/2,z,x+1/2
67 y+1/2,-z,x+1/2
68 y+1/2,z,-x+1/2
69 -z+1/2,-x,-y+1/2
70 -z+1/2,x,y+1/2
71 z+1/2,-x,y+1/2
72 z+1/2,x,-y+1/2
73 x+1/2,y+1/2,z
74 x+1/2,-y+1/2,-z
75 -x+1/2,y+1/2,-z
76 -x+1/2,-y+1/2,z
77 y+1/2,z+1/2,x
78 y+1/2,-z+1/2,-x
79 -y+1/2,z+1/2,-x
80 -y+1/2,-z+1/2,x
81 z+1/2,x+1/2,y
82 z+1/2,-x+1/2,-y
83 -z+1/2,x+1/2,-y
84 -z+1/2,-x+1/2,y
85 -x+1/2,-y+1/2,-z
86 -x+1/2,y+1/2,z
87 x+1/2,-y+1/2,z
88 x+1/2,y+1/2,-z
89 -y+1/2,-z+1/2,-x
90 -y+1/2,z+1/2,x
91 y+1/2,-z+1/2,x
92 y+1/2,z+1/2,-x
93 -z+1/2,-x+1/2,-y
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95 z+1/2,-x+1/2,y
96 z+1/2,x+1/2,-y

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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Co1 Co 4 a 0.00000 0.00000 0.00000 1.00000
K1 K 4 b 0.50000 0.50000 0.50000 1.00000
K2 K 8 c 0.25000 0.25000 0.25000 1.00000
N1 N 24 e 0.19500 0.00000 0.00000 1.00000
O1 O 48 h 0.00000 0.23500 0.10000 1.00000
```

$K_3Co(NO_2)_6$ ($J2_4$): AB3C6D12_cF88_202_a_bc_e_h - POSCAR

```
AB3C6D12_cF88_202_a_bc_e_h & a,x4,y5,z5 --params=10.46,0.195,0.235,0.1 &
    ↪ Fm-3 T_{h}^{3} #202 (abceh) & cF88 & $J2_{4}$ & CoK3N6O12 &
    ↪ CoK3N6O12 & M. {van Driel} and H. J. Verweel, Zeitschrift f{"u}
    ↪ r Kristallographie - Crystalline Materials 95, 308-314 (1936)
   1.00000000000000
     0.00000000000000     5.23000000000000     5.23000000000000
     5.23000000000000     0.00000000000000     5.23000000000000
     5.23000000000000     5.23000000000000     0.00000000000000
    Co    K    N    O
     1    3    6   12
Direct
     0.00000000000000     0.00000000000000     0.00000000000000  Co  (4a)
     0.50000000000000     0.50000000000000     0.50000000000000  K   (4b)
     0.25000000000000     0.25000000000000     0.25000000000000  K   (8c)
     0.75000000000000     0.75000000000000     0.75000000000000  K   (8c)
    -0.19500000000000     0.19500000000000     0.19500000000000  N  (24e)
     0.19500000000000    -0.19500000000000    -0.19500000000000  N  (24e)
     0.19500000000000    -0.19500000000000     0.19500000000000  N  (24e)
    -0.19500000000000     0.19500000000000    -0.19500000000000  N  (24e)
     0.19500000000000     0.19500000000000    -0.19500000000000  N  (24e)
    -0.19500000000000    -0.19500000000000     0.19500000000000  N  (24e)
     0.33500000000000    -0.13500000000000     0.13500000000000  O  (48h)
    -0.13500000000000     0.33500000000000    -0.33500000000000  O  (48h)
     0.13500000000000    -0.33500000000000     0.33500000000000  O  (48h)
    -0.33500000000000     0.13500000000000    -0.13500000000000  O  (48h)
     0.13500000000000     0.33500000000000    -0.13500000000000  O  (48h)
    -0.33500000000000    -0.13500000000000     0.33500000000000  O  (48h)
     0.33500000000000     0.13500000000000    -0.33500000000000  O  (48h)
    -0.13500000000000    -0.33500000000000     0.13500000000000  O  (48h)
    -0.13500000000000     0.13500000000000     0.33500000000000  O  (48h)
     0.33500000000000    -0.33500000000000    -0.13500000000000  O  (48h)
    -0.33500000000000     0.33500000000000     0.13500000000000  O  (48h)
     0.13500000000000    -0.13500000000000    -0.33500000000000  O  (48h)
```

$LaFe_4P_{12}$: A4BC12_cI34_204_c_a_g - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Fe4LaP12'
_chemical_formula_sum 'Fe4 La P12'

loop_
_publ_author_name
 'W. Jeitschko'
 'D. Braun'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 33
_journal_year 1977
_journal_page_first 3401
_journal_page_last 3406
_publ_Section_title
;
 LaFe$_{4}$P$_{12}$ with filled CoAs$_{3}$-type structure and isotypic
    ↪ lanthanoid-transition metal polyphosphides
;

# Found in Novel Heavy Fermion Behavior in Praseodymium-based Materials:
    ↪ Experimental Study of PrOs$_{4}$Sb$_{12}$, {Ph. D. Thesis,
    ↪ University of Florida}, 2007

_aflow_title 'LaFe$_{4}$P$_{12}$ Structure'
_aflow_proto 'A4BC12_cI34_204_c_a_g'
_aflow_params 'a,y_{3},z_{3}'
_aflow_params_values '7.832,0.3539,0.1504'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'cI34'

_symmetry_space_group_name_H-M "I 2/m -3"
_symmetry_Int_Tables_number 204
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_cell_length_b     7.83200
_cell_length_c     7.83200
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 y,z,x
6 y,-z,-x
7 -y,z,-x
8 -y,-z,x
9 z,x,y
10 z,-x,-y
11 -z,x,-y
12 -z,-x,y
13 -x,-y,-z
14 -x,y,z
15 x,-y,z
16 x,y,-z
17 -y,-z,-x
18 -y,z,x
19 y,-z,x
20 y,z,-x
21 -z,-x,-y
22 -z,x,y
23 z,-x,y
24 z,x,-y
25 x+1/2,y+1/2,z+1/2
26 x+1/2,-y+1/2,-z+1/2
27 -x+1/2,y+1/2,-z+1/2
28 -x+1/2,-y+1/2,z+1/2
29 y+1/2,z+1/2,x+1/2
30 y+1/2,-z+1/2,-x+1/2
31 -y+1/2,z+1/2,-x+1/2
32 -y+1/2,-z+1/2,x+1/2
33 z+1/2,x+1/2,y+1/2
34 z+1/2,-x+1/2,-y+1/2
35 -z+1/2,x+1/2,-y+1/2
36 -z+1/2,-x+1/2,y+1/2
37 -x+1/2,-y+1/2,-z+1/2
38 -x+1/2,y+1/2,z+1/2
39 x+1/2,-y+1/2,z+1/2
40 x+1/2,y+1/2,-z+1/2
41 -y+1/2,-z+1/2,-x+1/2
42 -y+1/2,z+1/2,x+1/2
43 y+1/2,-z+1/2,x+1/2
44 y+1/2,z+1/2,-x+1/2
45 -z+1/2,-x+1/2,-y+1/2
46 -z+1/2,x+1/2,y+1/2
47 z+1/2,-x+1/2,y+1/2
48 z+1/2,x+1/2,-y+1/2

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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
La1 La 2 a 0.00000 0.00000 0.00000 1.00000
Fe1 Fe 8 c 0.25000 0.25000 0.25000 1.00000
P1 P 24 g 0.00000 0.35390 0.15040 1.00000
```

$LaFe_4P_{12}$: A4BC12_cI34_204_c_a_g - POSCAR

```
A4BC12_cI34_204_c_a_g & a,y3,z3 --params=7.832,0.3539,0.1504 & Im-3 T_{h
    ↪ }^{5} #204 (acg) & cI34 & None & Fe4LaP12 & Fe4LaP12 & W.
    ↪ Jeitschko and D. Braun, Acta Crystallogr. Sect. B Struct. Sci.
    ↪ 33, 3401-3406 (1977)
   1.00000000000000
  -3.91600000000000    3.91600000000000    3.91600000000000
   3.91600000000000   -3.91600000000000    3.91600000000000
   3.91600000000000    3.91600000000000   -3.91600000000000
    Fe     La     P
     4      1     12
Direct
   0.50000000000000    0.50000000000000    0.50000000000000   Fe   (8c)
   0.00000000000000    0.00000000000000    0.50000000000000   Fe   (8c)
   0.00000000000000    0.50000000000000    0.00000000000000   Fe   (8c)
   0.50000000000000    0.00000000000000    0.00000000000000   Fe   (8c)
   0.00000000000000    0.00000000000000    0.00000000000000   La   (2a)
   0.50430000000000    0.15040000000000    0.35390000000000    P   (24g)
  -0.20350000000000    0.15040000000000   -0.35390000000000    P   (24g)
   0.20350000000000   -0.15040000000000    0.35390000000000    P   (24g)
  -0.50430000000000   -0.15040000000000   -0.35390000000000    P   (24g)
   0.35390000000000    0.50430000000000    0.15040000000000    P   (24g)
  -0.35390000000000   -0.20350000000000    0.15040000000000    P   (24g)
   0.35390000000000    0.20350000000000   -0.15040000000000    P   (24g)
  -0.35390000000000   -0.50430000000000   -0.15040000000000    P   (24g)
   0.15040000000000    0.35390000000000    0.50430000000000    P   (24g)
   0.15040000000000   -0.35390000000000   -0.20350000000000    P   (24g)
  -0.15040000000000    0.35390000000000    0.20350000000000    P   (24g)
  -0.15040000000000   -0.35390000000000   -0.50430000000000    P   (24g)
```

$NaMn_7O_{12}$: A7BC12_cI40_204_bc_a_g - CIF

```
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_audit_creation_method FINDSYM

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_chemical_formula_sum 'Mn7 Na O12'

loop_
_publ_author_name
 'E. Gilioli'
 'G. Calestani'
 'F. Licci'
 'A. Gauzzi'
 'F. Bolzoni'
 'A. Prodi'
 'M. Marezio'
_journal_name_full_name
;
 Solid State Sciences
;
_journal_volume 7
_journal_year 2005
_journal_page_first 746
_journal_page_last 752
_publ_Section_title
;
 $P-T$ phase diagram and single crystal structural refinement of NaMn$_{
    ↪ 7}$O$_{12}$
;

# Found in Crystal growth and structural refinement of NaMn$_{7}$O$_{12}
    ↪ $, 2005

_aflow_title 'NaMn$_{7}$O$_{12}$ Structure'
_aflow_proto 'A7BC12_cI40_204_bc_a_g'
_aflow_params 'a,y_{4},z_{4}'
_aflow_params_values '7.312,0.3128,0.1829'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'cI40'

_symmetry_space_group_name_H-M "I 2/m -3"
_symmetry_Int_Tables_number 204

_cell_length_a     7.31200
_cell_length_b     7.31200
_cell_length_c     7.31200
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

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_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 y,z,x
6 y,-z,-x
7 -y,z,-x
8 -y,-z,x
9 z,x,y
10 z,-x,-y
11 -z,x,-y
12 -z,-x,y
13 -x,-y,-z
14 -x,y,z
15 x,-y,z
16 x,y,-z
17 -y,-z,-x
18 -y,z,x
19 y,-z,x
20 y,z,-x
21 -z,-x,-y
22 -z,x,y
23 z,-x,y
24 z,x,-y
25 x+1/2,y+1/2,z+1/2
26 x+1/2,-y+1/2,-z+1/2
27 -x+1/2,y+1/2,-z+1/2
28 -x+1/2,-y+1/2,z+1/2
29 y+1/2,z+1/2,x+1/2
30 y+1/2,-z+1/2,-x+1/2
31 -y+1/2,z+1/2,-x+1/2
32 -y+1/2,-z+1/2,x+1/2
33 z+1/2,x+1/2,y+1/2
34 z+1/2,-x+1/2,-y+1/2
35 -z+1/2,x+1/2,-y+1/2
36 -z+1/2,-x+1/2,y+1/2
37 -x+1/2,-y+1/2,-z+1/2
38 -x+1/2,y+1/2,z+1/2
39 x+1/2,-y+1/2,z+1/2
40 x+1/2,y+1/2,-z+1/2
41 -y+1/2,-z+1/2,-x+1/2
42 -y+1/2,z+1/2,x+1/2
43 y+1/2,-z+1/2,x+1/2
44 y+1/2,z+1/2,-x+1/2
45 -z+1/2,-x+1/2,-y+1/2
46 -z+1/2,x+1/2,y+1/2
47 z+1/2,-x+1/2,y+1/2
48 z+1/2,x+1/2,-y+1/2

loop_
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_atom_site_type_symbol
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_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Na1 Na 2 a 0.00000 0.00000 0.00000 1.00000
Mn1 Mn 6 b 0.00000 0.50000 0.50000 1.00000
Mn2 Mn 8 c 0.25000 0.25000 0.25000 1.00000
O1 O 24 g 0.00000 0.31280 0.18290 1.00000
```

$NaMn_7O_{12}$: A7BC12_cI40_204_bc_a_g - POSCAR

```
A7BC12_cI40_204_bc_a_g & a,y4,z4 --params=7.312,0.3128,0.1829 & Im-3 T_{
    ↪ h}^{5} #204 (abcg) & cI40 & None & Mn7NaO12 & Mn7NaO12 & E.
    ↪ Gilioli et al., Solid State Sci. 7, 746-752 (2005)
   1.00000000000000
  -3.65600000000000    3.65600000000000    3.65600000000000
   3.65600000000000   -3.65600000000000    3.65600000000000
   3.65600000000000    3.65600000000000   -3.65600000000000
    Mn     Na     O
     7      1     12
Direct
   0.00000000000000    0.50000000000000    0.50000000000000   Mn   (6b)
   0.50000000000000    0.00000000000000    0.50000000000000   Mn   (6b)
   0.50000000000000    0.50000000000000    0.00000000000000   Mn   (6b)
   0.50000000000000    0.50000000000000    0.50000000000000   Mn   (8c)
   0.00000000000000    0.00000000000000    0.50000000000000   Mn   (8c)
   0.00000000000000    0.50000000000000    0.00000000000000   Mn   (8c)
   0.50000000000000    0.00000000000000    0.00000000000000   Mn   (8c)
   0.00000000000000    0.00000000000000    0.00000000000000   Na   (2a)
   0.49570000000000    0.18290000000000    0.31280000000000   O   (24g)
  -0.12990000000000    0.18290000000000   -0.31280000000000   O   (24g)
   0.12990000000000   -0.18290000000000    0.31280000000000   O   (24g)
  -0.49570000000000   -0.18290000000000   -0.31280000000000   O   (24g)
   0.31280000000000    0.49570000000000    0.18290000000000   O   (24g)
  -0.31280000000000   -0.12990000000000    0.18290000000000   O   (24g)
   0.31280000000000    0.12990000000000   -0.18290000000000   O   (24g)
  -0.31280000000000   -0.49570000000000   -0.18290000000000   O   (24g)
   0.18290000000000    0.31280000000000    0.49570000000000   O   (24g)
   0.18290000000000   -0.31280000000000   -0.12990000000000   O   (24g)
  -0.18290000000000    0.31280000000000    0.12990000000000   O   (24g)
  -0.18290000000000   -0.31280000000000   -0.49570000000000   O   (24g)
```

NO_2 (Modern, *C*26): AB2_cI36_204_d_g - CIF

```
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_audit_creation_method FINDSYM

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_chemical_formula_sum 'N O2'

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 '{\AA}. Kvick'
 'R. K. {McMullan}'
 'M. D. Newton'
_journal_name_full_name
;
 Journal of Chemical Physics
;
_journal_volume 76
_journal_year 1982
_journal_page_first 3754
_journal_page_last 3761
_publ_Section_title
;
 The structure of dinitrogen tetroxide N$_{2}$O$_{4}$: Neutron
    ↪ diffraction study at 100, 60 and 20 K and {\em ab initio}
    ↪ theoretical calculations
;

# Found in Crystal Structure Data of Inorganic Compounds, 2005 Found in
    ↪ Crystal Structure Data of Inorganic Compounds, {
    ↪ Landolt-Bornstein Volume III 43A2},

_aflow_title 'NO$_{2}$ (Modern, $C26$) Structure'
_aflow_proto 'AB2_cI36_204_d_g'
_aflow_params 'a,x_{1},y_{2},z_{2}'
_aflow_params_values '7.6937,0.38587,0.32597,0.1425'
_aflow_Strukturbericht '$C26$'
_aflow_Pearson 'cI36'

_symmetry_space_group_name_H-M "I 2/m -3"
_symmetry_Int_Tables_number 204

_cell_length_a     7.69370
_cell_length_b     7.69370
_cell_length_c     7.69370
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 y,z,x
6 y,-z,-x
7 -y,z,-x
8 -y,-z,x
9 z,x,y
10 z,-x,-y
11 -z,x,-y
12 -z,-x,y
13 -x,-y,-z
14 -x,y,z
15 x,-y,z
16 x,y,-z
17 -y,-z,-x
18 -y,z,x
19 y,-z,x
20 y,z,-x
21 -z,-x,-y
22 -z,x,y
23 z,-x,y
24 z,x,-y
25 x+1/2,y+1/2,z+1/2
26 x+1/2,-y+1/2,-z+1/2
27 -x+1/2,y+1/2,-z+1/2
28 -x+1/2,-y+1/2,z+1/2
29 y+1/2,z+1/2,x+1/2
30 y+1/2,-z+1/2,-x+1/2
31 -y+1/2,z+1/2,-x+1/2
32 -y+1/2,-z+1/2,x+1/2
33 z+1/2,x+1/2,y+1/2
34 z+1/2,-x+1/2,-y+1/2
35 -z+1/2,x+1/2,-y+1/2
36 -z+1/2,-x+1/2,y+1/2
37 -x+1/2,-y+1/2,-z+1/2
38 -x+1/2,y+1/2,z+1/2
39 x+1/2,-y+1/2,z+1/2
40 x+1/2,y+1/2,-z+1/2
41 -y+1/2,-z+1/2,-x+1/2
42 -y+1/2,z+1/2,x+1/2
43 y+1/2,-z+1/2,x+1/2
44 y+1/2,z+1/2,-x+1/2
45 -z+1/2,-x+1/2,-y+1/2
46 -z+1/2,x+1/2,y+1/2
47 z+1/2,-x+1/2,y+1/2
48 z+1/2,x+1/2,-y+1/2

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_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
N1 N 12 d 0.38587 0.00000 0.00000 1.00000
O1 O 24 g 0.00000 0.32597 0.14250 1.00000
```

NO_2 (Modern, *C*26): AB2_cI36_204_d_g - POSCAR

```
AB2_cI36_204_d_g & a,x1,y2,z2 --params=7.6937,0.38587,0.32597,0.1425 &
    ↪ Im-3 T_{h}^{5} #204 (dg) & cI36 & $C26$ & NO2 & NO2 & {\AA}.
    ↪ Kvick and R. K. {McMullan} and M. D. Newton, J. Chem. Phys. 76,
    ↪ 3754-3761 (1982)
   1.00000000000000
  -3.84685000000000    3.84685000000000    3.84685000000000
   3.84685000000000   -3.84685000000000    3.84685000000000
   3.84685000000000    3.84685000000000   -3.84685000000000
     N      O
     6     12
Direct
   0.00000000000000    0.38587000000000    0.38587000000000   N   (12d)
   0.00000000000000   -0.38587000000000   -0.38587000000000   N   (12d)
   0.38587000000000    0.00000000000000    0.38587000000000   N   (12d)
  -0.38587000000000    0.00000000000000   -0.38587000000000   N   (12d)
   0.38587000000000    0.38587000000000    0.00000000000000   N   (12d)
  -0.38587000000000   -0.38587000000000    0.00000000000000   N   (12d)
   0.46847000000000    0.14250000000000    0.32597000000000   O   (24g)
  -0.18347000000000    0.14250000000000   -0.32597000000000   O   (24g)
   0.18347000000000   -0.14250000000000    0.32597000000000   O   (24g)
  -0.46847000000000   -0.14250000000000   -0.32597000000000   O   (24g)
   0.32597000000000    0.46847000000000    0.14250000000000   O   (24g)
  -0.32597000000000   -0.18347000000000    0.14250000000000   O   (24g)
   0.32597000000000    0.18347000000000   -0.14250000000000   O   (24g)
  -0.32597000000000   -0.46847000000000   -0.14250000000000   O   (24g)
   0.14250000000000    0.32597000000000    0.46847000000000   O   (24g)
   0.14250000000000   -0.32597000000000   -0.18347000000000   O   (24g)
  -0.14250000000000    0.32597000000000    0.18347000000000   O   (24g)
  -0.14250000000000   -0.32597000000000   -0.46847000000000   O   (24g)
```

$Zn(BrO_3)_2{\cdot}6H_2O$ ($J1_{10}$): A2B6C6D_cP60_205_c_d_d_a - CIF

```
# CIF file
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_chemical_name_mineral 'Br2(H6O)6O6Zn'
_chemical_formula_sum 'Br2 (H2O)6 O6 Zn'

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_publ_author_name
 'S. H. Y\"{u}'
 'C. A. Beevers'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 95
_journal_year 1936
_journal_page_first 426
```

```
_journal_page_last 434
_publ_Section_title
;
 The Crystal Structure of Zinc Bromate Hexahydrate [Zn(BrO$_{3}$)$_{2}$
      ↪ $\cdot$ 6H$_{2}$O]
;

# Found in Strukturbericht Band IV 1936, 1938

_aflow_title 'Zn(BrO$_{3}$)$_{2}$$\cdot$6H$_{2}$O ($J1_{10}$) Structure'
_aflow_proto 'A2B6C6D_cP60_205_c_d_d_a'
_aflow_params 'a,x_{2},x_{3},y_{3},z_{3},x_{4},y_{4},z_{4}'
_aflow_params_values '10.316,0.259,0.195,0.05,0.965,0.19,0.145,0.33'
_aflow_Strukturbericht '$J1_{10}$'
_aflow_Pearson 'cP60'

_symmetry_space_group_name_H-M "P 21/a -3"
_symmetry_Int_Tables_number 205

_cell_length_a    10.31600
_cell_length_b    10.31600
_cell_length_c    10.31600
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

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_space_group_symop_id
_space_group_symop_operation_xyz
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2 x+1/2,-y+1/2,-z
3 -x,y+1/2,-z+1/2
4 -x+1/2,-y,z+1/2
5 y,z,x
6 y+1/2,-z+1/2,-x
7 -y,z+1/2,-x+1/2
8 -y+1/2,-z,x+1/2
9 z,x,y
10 z+1/2,-x+1/2,-y
11 -z,x+1/2,-y+1/2
12 -z+1/2,-x,y+1/2
13 -x,-y,-z
14 -x+1/2,y+1/2,z
15 x,-y+1/2,z+1/2
16 x+1/2,y,-z+1/2
17 -y,-z,-x
18 -y+1/2,z+1/2,x
19 y,-z+1/2,x+1/2
20 y+1/2,z,-x+1/2
21 -z,-x,-y
22 -z+1/2,x+1/2,y
23 z,-x+1/2,y+1/2
24 z+1/2,x,-y+1/2

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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Zn1 Zn 4 a 0.00000 0.00000 0.00000 1.00000
Br1 Br 8 c 0.25900 0.25900 0.25900 1.00000
H2O1 H2O 24 d 0.19500 0.05000 0.96500 1.00000
O1 O 24 d 0.19000 0.14500 0.33000 1.00000
```

$Zn(BrO_3)_2{\cdot}6H_2O$ ($J1_{10}$): A2B6C6D_cP60_205_c_d_d_a - POSCAR

```
A2B6C6D_cP60_205_c_d_d_a & a,x2,x3,y3,z3,x4,y4,z4 --params=10.316,0.259,
      ↪ 0.195,0.05,0.965,0.19,0.145,0.33 & Pa-3 T_{h}^{6} #205 (acd^2)
      ↪ & cP60 & $J1_{10}$ & Br2(H6O)6O6Zn & Br2(H6O)6O6Zn & S. H. Y\"{
      ↪ u} and C. A. Beevers, Zeitschrift f{"u}r Kristallographie -
      ↪ Crystalline Materials 95, 426-434 (1936)
   1.00000000000000
  10.31600000000000    0.00000000000000    0.00000000000000
   0.00000000000000   10.31600000000000    0.00000000000000
   0.00000000000000    0.00000000000000   10.31600000000000
     Br    H2O      O     Zn
      8     24     24      4
Direct
   0.25900000000000    0.25900000000000    0.25900000000000   Br    (8c)
   0.24100000000000   -0.25900000000000    0.75900000000000   Br    (8c)
  -0.25900000000000    0.75900000000000    0.24100000000000   Br    (8c)
   0.75900000000000    0.24100000000000   -0.25900000000000   Br    (8c)
  -0.25900000000000   -0.25900000000000   -0.25900000000000   Br    (8c)
   0.75900000000000    0.25900000000000    0.24100000000000   Br    (8c)
   0.25900000000000    0.24100000000000    0.75900000000000   Br    (8c)
   0.24100000000000    0.75900000000000    0.25900000000000   Br    (8c)
   0.19500000000000    0.05000000000000    0.96500000000000   H2O   (24d)
   0.30500000000000   -0.05000000000000    1.46500000000000   H2O   (24d)
  -0.19500000000000    0.55000000000000   -0.46500000000000   H2O   (24d)
   0.69500000000000    0.45000000000000   -0.96500000000000   H2O   (24d)
   0.96500000000000    0.19500000000000    0.05000000000000   H2O   (24d)
   1.46500000000000    0.30500000000000   -0.05000000000000   H2O   (24d)
  -0.46500000000000   -0.19500000000000    0.55000000000000   H2O   (24d)
  -0.96500000000000    0.69500000000000    0.45000000000000   H2O   (24d)
   0.05000000000000    0.96500000000000    0.19500000000000   H2O   (24d)
  -0.05000000000000    1.46500000000000    0.30500000000000   H2O   (24d)
   0.55000000000000   -0.46500000000000   -0.19500000000000   H2O   (24d)
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  -0.19500000000000   -0.05000000000000   -0.96500000000000   H2O   (24d)
   0.69500000000000    0.05000000000000   -0.46500000000000   H2O   (24d)
   0.19500000000000    0.45000000000000    1.46500000000000   H2O   (24d)
   0.30500000000000    0.55000000000000    0.96500000000000   H2O   (24d)
  -0.96500000000000   -0.19500000000000   -0.05000000000000   H2O   (24d)
  -0.46500000000000    0.69500000000000    0.05000000000000   H2O   (24d)
   1.46500000000000    0.19500000000000    0.45000000000000   H2O   (24d)
   0.96500000000000    0.30500000000000    0.55000000000000   H2O   (24d)
  -0.05000000000000   -0.96500000000000   -0.19500000000000   H2O   (24d)
   0.05000000000000   -0.46500000000000    0.69500000000000   H2O   (24d)
   0.45000000000000    1.46500000000000    0.19500000000000   H2O   (24d)
   0.55000000000000    0.96500000000000    0.30500000000000   H2O   (24d)
   0.19000000000000    0.14500000000000    0.33000000000000   O     (24d)
   0.31000000000000   -0.14500000000000    0.83000000000000   O     (24d)
  -0.19000000000000    0.64500000000000    0.17000000000000   O     (24d)
   0.69000000000000    0.35500000000000   -0.33000000000000   O     (24d)
   0.33000000000000    0.19000000000000    0.14500000000000   O     (24d)
   0.83000000000000    0.31000000000000   -0.14500000000000   O     (24d)
   0.17000000000000   -0.19000000000000    0.64500000000000   O     (24d)
  -0.33000000000000    0.69000000000000    0.35500000000000   O     (24d)
   0.14500000000000    0.33000000000000    0.19000000000000   O     (24d)
  -0.14500000000000    0.83000000000000    0.31000000000000   O     (24d)
   0.64500000000000    0.17000000000000   -0.19000000000000   O     (24d)
   0.35500000000000   -0.33000000000000    0.69000000000000   O     (24d)
  -0.19000000000000   -0.14500000000000   -0.33000000000000   O     (24d)
   0.69000000000000    0.14500000000000    0.17000000000000   O     (24d)
   0.19000000000000    0.35500000000000    0.83000000000000   O     (24d)
   0.31000000000000    0.64500000000000    0.33000000000000   O     (24d)
  -0.33000000000000   -0.19000000000000   -0.14500000000000   O     (24d)
   0.17000000000000    0.69000000000000    0.14500000000000   O     (24d)
   0.83000000000000    0.19000000000000    0.35500000000000   O     (24d)
   0.33000000000000    0.31000000000000    0.64500000000000   O     (24d)
  -0.14500000000000   -0.33000000000000   -0.19000000000000   O     (24d)
   0.14500000000000    0.17000000000000    0.69000000000000   O     (24d)
   0.35500000000000    0.83000000000000    0.19000000000000   O     (24d)
   0.64500000000000    0.33000000000000    0.31000000000000   O     (24d)
   0.00000000000000    0.00000000000000    0.00000000000000   Zn    (4a)
   0.50000000000000    0.00000000000000    0.50000000000000   Zn    (4a)
   0.00000000000000    0.50000000000000    0.50000000000000   Zn    (4a)
   0.50000000000000    0.50000000000000    0.00000000000000   Zn    (4a)
```

$H6_4$ $[Ni(NO_3)_2(NH_3)_6]$ (*obsolete*): A2B6CD6_cP60_205_c_d_a_d - CIF

```
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_audit_creation_method FINDSYM

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_chemical_formula_sum 'N2 (NH3)6 Ni O6'

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 'R. W. G. Wyckoff'
_journal_name_full_name
;
 Journal of the American Chemical Society
;
_journal_volume 44
_journal_year 1922
_journal_page_first 1260
_journal_page_last 1266
_publ_Section_title
;
 The Composition and Crystal Structure of Nickel Nitrate Hexammoniate
;

# Found in Strukturbericht 1913-1928, 1931

_aflow_title '$H6_{4}$ [Ni(NO$_{3}$)$_{2}$(NH$_{3}$)$_{6}$] ({\em{
      ↪ obsolete}}) Structure'
_aflow_proto 'A2B6CD6_cP60_205_c_d_a_d'
_aflow_params 'a,x_{2},x_{3},y_{3},z_{3},x_{4},y_{4},z_{4}'
_aflow_params_values '10.96,0.25,0.24,0.0,0.0,0.25,0.25,0.01'
_aflow_Strukturbericht '$H6_{4}$'
_aflow_Pearson 'cP60'

_symmetry_space_group_name_H-M "P 21/a -3"
_symmetry_Int_Tables_number 205

_cell_length_a    10.96000
_cell_length_b    10.96000
_cell_length_c    10.96000
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
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2 x+1/2,-y+1/2,-z
3 -x,y+1/2,-z+1/2
4 -x+1/2,-y,z+1/2
5 y,z,x
6 y+1/2,-z+1/2,-x
7 -y,z+1/2,-x+1/2
8 -y+1/2,-z,x+1/2
9 z,x,y
10 z+1/2,-x+1/2,-y
11 -z,x+1/2,-y+1/2
12 -z+1/2,-x,y+1/2
13 -x,-y,-z
14 -x+1/2,y+1/2,z
15 x,-y+1/2,z+1/2
16 x+1/2,y,-z+1/2
17 -y,-z,-x
18 -y+1/2,z+1/2,x
19 y,-z+1/2,x+1/2
```

```
20 y+1/2,z,-x+1/2
21 -z,-x,-y
22 -z+1/2,x+1/2,y
23 z,-x+1/2,y+1/2
24 z+1/2,x,-y+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ni1 Ni 4 a 0.00000 0.00000 0.00000 1.00000
N1 N 8 c 0.25000 0.25000 0.25000 1.00000
NH31 NH3 24 d 0.24000 0.00000 0.00000 1.00000
O1 O 24 d 0.25000 0.25000 0.01000 1.00000
```

*H*6$_4$ [Ni(NO$_3$)$_2$(NH$_3$)$_6$] (*obsolete*): A2B6CD6_cP60_205_c_d_a_d - POSCAR

```
A2B6CD6_cP60_205_c_d_a_d & a,x2,x3,y3,z3,x4,y4,z4 --params=10.96,0.25,
    ↪ 0.24,0.0,0.0,0.25,0.25,0.01 & Pa-3 T_{h}^{6} #205 (acd^2) &
    ↪ cP60 & $H6_{4}$ & N2(NH3)6NiO6 & N2(NH3)6NiO6 & R. W. G.
    ↪ Wyckoff, J. Am. Chem. Soc. 44, 1260-1266 (1922)
   1.00000000000000
  10.96000000000000    0.00000000000000    0.00000000000000
   0.00000000000000   10.96000000000000    0.00000000000000
   0.00000000000000    0.00000000000000   10.96000000000000
     N    NH3     Ni      O
     8     24      4     24
Direct
   0.25000000000000    0.25000000000000    0.25000000000000     N    (8c)
   0.25000000000000   -0.25000000000000    0.75000000000000     N    (8c)
  -0.25000000000000    0.75000000000000    0.25000000000000     N    (8c)
   0.75000000000000    0.25000000000000   -0.25000000000000     N    (8c)
  -0.25000000000000   -0.25000000000000   -0.25000000000000     N    (8c)
   0.75000000000000    0.25000000000000    0.25000000000000     N    (8c)
   0.25000000000000    0.25000000000000    0.75000000000000     N    (8c)
   0.25000000000000    0.75000000000000    0.25000000000000     N    (8c)
   0.24000000000000    0.00000000000000    0.00000000000000   NH3   (24d)
   0.26000000000000    0.00000000000000    0.50000000000000   NH3   (24d)
  -0.24000000000000    0.50000000000000    0.50000000000000   NH3   (24d)
   0.74000000000000    0.50000000000000    0.00000000000000   NH3   (24d)
   0.00000000000000    0.24000000000000    0.00000000000000   NH3   (24d)
   0.50000000000000    0.26000000000000    0.00000000000000   NH3   (24d)
   0.50000000000000   -0.24000000000000    0.50000000000000   NH3   (24d)
   0.00000000000000    0.74000000000000    0.50000000000000   NH3   (24d)
   0.00000000000000    0.00000000000000    0.24000000000000   NH3   (24d)
   0.00000000000000    0.50000000000000    0.26000000000000   NH3   (24d)
   0.50000000000000    0.50000000000000   -0.24000000000000   NH3   (24d)
   0.50000000000000    0.00000000000000    0.74000000000000   NH3   (24d)
  -0.24000000000000    0.00000000000000    0.00000000000000   NH3   (24d)
   0.74000000000000    0.00000000000000    0.50000000000000   NH3   (24d)
   0.24000000000000    0.50000000000000    0.50000000000000   NH3   (24d)
   0.26000000000000    0.50000000000000    0.00000000000000   NH3   (24d)
   0.00000000000000   -0.24000000000000    0.00000000000000   NH3   (24d)
   0.50000000000000    0.74000000000000    0.00000000000000   NH3   (24d)
   0.50000000000000    0.24000000000000    0.50000000000000   NH3   (24d)
   0.00000000000000    0.26000000000000    0.50000000000000   NH3   (24d)
   0.00000000000000    0.00000000000000   -0.24000000000000   NH3   (24d)
   0.00000000000000    0.50000000000000    0.74000000000000   NH3   (24d)
   0.50000000000000    0.50000000000000    0.24000000000000   NH3   (24d)
   0.50000000000000    0.00000000000000    0.26000000000000   NH3   (24d)
   0.00000000000000    0.00000000000000    0.00000000000000    Ni    (4a)
   0.50000000000000    0.00000000000000    0.50000000000000    Ni    (4a)
   0.00000000000000    0.50000000000000    0.50000000000000    Ni    (4a)
   0.50000000000000    0.50000000000000    0.00000000000000    Ni    (4a)
   0.25000000000000    0.25000000000000    0.01000000000000     O   (24d)
   0.25000000000000   -0.25000000000000    0.51000000000000     O   (24d)
  -0.25000000000000    0.75000000000000    0.49000000000000     O   (24d)
   0.75000000000000    0.25000000000000   -0.01000000000000     O   (24d)
   0.01000000000000    0.25000000000000    0.25000000000000     O   (24d)
   0.51000000000000    0.25000000000000   -0.25000000000000     O   (24d)
   0.49000000000000   -0.25000000000000    0.75000000000000     O   (24d)
  -0.01000000000000    0.75000000000000    0.25000000000000     O   (24d)
   0.25000000000000    0.01000000000000    0.25000000000000     O   (24d)
  -0.25000000000000    0.51000000000000    0.25000000000000     O   (24d)
   0.75000000000000    0.49000000000000   -0.25000000000000     O   (24d)
   0.25000000000000   -0.01000000000000    0.75000000000000     O   (24d)
  -0.25000000000000   -0.25000000000000   -0.01000000000000     O   (24d)
   0.75000000000000    0.25000000000000    0.49000000000000     O   (24d)
   0.25000000000000    0.25000000000000    0.51000000000000     O   (24d)
   0.25000000000000    0.75000000000000    0.01000000000000     O   (24d)
  -0.01000000000000   -0.25000000000000   -0.25000000000000     O   (24d)
   0.49000000000000    0.75000000000000    0.25000000000000     O   (24d)
   0.51000000000000    0.25000000000000    0.25000000000000     O   (24d)
   0.01000000000000    0.25000000000000    0.75000000000000     O   (24d)
  -0.25000000000000   -0.01000000000000   -0.25000000000000     O   (24d)
   0.25000000000000    0.49000000000000    0.75000000000000     O   (24d)
   0.25000000000000    0.51000000000000    0.25000000000000     O   (24d)
   0.75000000000000    0.01000000000000    0.25000000000000     O   (24d)
```

Pb(NO$_3$)$_2$ (*G*2$_1$): A2B6C_cP36_205_c_d_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'N2O6Pb'
_chemical_formula_sum 'N2 O6 Pb'

loop_
_publ_author_name
 'H. Nowotny'
 'G. Heger'
_journal_name_full_name
;
 Acta Crystallographica Section C: Structural Chemistry
;
_journal_volume 42
_journal_year 1986
_journal_page_first 133
_journal_page_last 135
_publ_Section_title
;
 Structure Refinement of Lead Nitrate
;

_aflow_title 'Pb(NO$_{3}$)$_{2}$ ($G2_{1}$) Structure'
_aflow_proto 'A2B6C_cP36_205_c_d_a'
_aflow_params 'a,x_{2},x_{3},y_{3},z_{3}'
_aflow_params_values '7.8586,0.34865,0.27772,0.28869,0.4773'
_aflow_Strukturbericht '$G2_{1}$'
_aflow_Pearson 'cP36'

_symmetry_space_group_name_H-M "P 21/a -3"
_symmetry_Int_Tables_number 205

_cell_length_a     7.85860
_cell_length_b     7.85860
_cell_length_c     7.85860
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

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_space_group_symop_operation_xyz
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2 x+1/2,-y+1/2,-z
3 -x,y+1/2,-z+1/2
4 -x+1/2,-y,z+1/2
5 y,z,x
6 y+1/2,-z+1/2,-x
7 -y,z+1/2,-x+1/2
8 -y+1/2,-z,x+1/2
9 z,x,y
10 z+1/2,-x+1/2,-y
11 -z,x+1/2,-y+1/2
12 -z+1/2,-x,y+1/2
13 -x,-y,-z
14 -x+1/2,y+1/2,z
15 x,-y+1/2,z+1/2
16 x+1/2,y,-z+1/2
17 -y,-z,-x
18 -y+1/2,z+1/2,x
19 y,-z+1/2,x+1/2
20 y+1/2,z,-x+1/2
21 -z,-x,-y
22 -z+1/2,x+1/2,y
23 z,-x+1/2,y+1/2
24 z+1/2,x,-y+1/2

loop_
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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Pb1 Pb 4 a 0.00000 0.00000 0.00000 1.00000
N1 N 8 c 0.34865 0.34865 0.34865 1.00000
O1 O 24 d 0.27772 0.28869 0.47730 1.00000
```

Pb(NO$_3$)$_2$ (*G*2$_1$): A2B6C_cP36_205_c_d_a - POSCAR

```
A2B6C_cP36_205_c_d_a & a,x2,x3,y3,z3 --params=7.8586,0.34865,0.27772,
    ↪ 0.28869,0.4773 & Pa-3 T_{h}^{6} #205 (acd) & cP36 & $G2_{1}$ &
    ↪ N2O6Pb & N2O6Pb & H. Nowotny and G. Heger, Acta Crystallogr. C
    ↪ 42, 133-135 (1986)
   1.00000000000000
   7.85860000000000    0.00000000000000    0.00000000000000
   0.00000000000000    7.85860000000000    0.00000000000000
   0.00000000000000    0.00000000000000    7.85860000000000
     N      O     Pb
     8     24      4
Direct
   0.34865000000000    0.34865000000000    0.34865000000000     N    (8c)
   0.15135000000000   -0.34865000000000    0.84865000000000     N    (8c)
  -0.34865000000000    0.84865000000000    0.15135000000000     N    (8c)
   0.84865000000000    0.15135000000000   -0.34865000000000     N    (8c)
  -0.34865000000000   -0.34865000000000   -0.34865000000000     N    (8c)
   0.84865000000000    0.34865000000000    0.15135000000000     N    (8c)
   0.34865000000000    0.15135000000000    0.84865000000000     N    (8c)
   0.15135000000000    0.84865000000000    0.34865000000000     N    (8c)
   0.27772000000000    0.28869000000000    0.47730000000000     O   (24d)
   0.22228000000000   -0.28869000000000    0.97730000000000     O   (24d)
  -0.27772000000000    0.78869000000000    0.02270000000000     O   (24d)
   0.77772000000000    0.21131000000000   -0.47730000000000     O   (24d)
   0.47730000000000    0.27772000000000    0.28869000000000     O   (24d)
   0.97730000000000    0.22228000000000   -0.28869000000000     O   (24d)
   0.02270000000000   -0.27772000000000    0.78869000000000     O   (24d)
  -0.47730000000000    0.77772000000000    0.21131000000000     O   (24d)
   0.28869000000000    0.47730000000000    0.27772000000000     O   (24d)
  -0.28869000000000    0.97730000000000    0.22228000000000     O   (24d)
   0.78869000000000    0.02270000000000   -0.27772000000000     O   (24d)
```

```
   0.21131000000000  −0.47730000000000   0.77772000000000   O   (24d)
  −0.27772000000000  −0.28869000000000  −0.47730000000000   O   (24d)
   0.77772000000000   0.28869000000000   0.02270000000000   O   (24d)
   0.27772000000000   0.21131000000000   0.97730000000000   O   (24d)
   0.22228000000000   0.78869000000000   0.47730000000000   O   (24d)
  −0.47730000000000  −0.27772000000000  −0.28869000000000   O   (24d)
   0.02270000000000   0.77772000000000   0.28869000000000   O   (24d)
   0.97730000000000   0.27772000000000   0.21131000000000   O   (24d)
   0.47730000000000   0.22228000000000   0.78869000000000   O   (24d)
  −0.28869000000000  −0.47730000000000  −0.27772000000000   O   (24d)
   0.28869000000000   0.02270000000000   0.77772000000000   O   (24d)
   0.21131000000000   0.97730000000000   0.27772000000000   O   (24d)
   0.78869000000000   0.47730000000000   0.22228000000000   O   (24d)
   0.00000000000000   0.00000000000000   0.00000000000000   Pb   (4a)
   0.50000000000000   0.00000000000000   0.50000000000000   Pb   (4a)
   0.00000000000000   0.50000000000000   0.50000000000000   Pb   (4a)
   0.50000000000000   0.50000000000000   0.00000000000000   Pb   (4a)
```

CaB_2O_4 (IV): A2BC4_cP84_205_d_ac_2d - CIF

```
# CIF file
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_audit_creation_method FINDSYM

_chemical_name_mineral 'B2CaO4'
_chemical_formula_sum 'B2 Ca O4'

loop_
_publ_author_name
 'M. Marezio'
 'J. P. Remeika'
 'P. D. Dernier'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 25
_journal_year 1969
_journal_page_first 965
_journal_page_last 970
_publ_Section_title
;
 The crystal structure of the high-pressure phase CaB$_{2}$O$_{4}$(IV),
      ↪ and polymorphism in CaB$_{2}$O$_{4}$
;

_aflow_title 'CaB$_{2}$O$_{4}$ (IV) Structure'
_aflow_proto 'A2BC4_cP84_205_d_ac_2d'
_aflow_params 'a,x_{2},x_{3},y_{3},z_{3},x_{4},y_{4},z_{4},x_{5},y_{5},
      ↪ z_{5}'
_aflow_params_values '9.008,0.37305,0.1189,0.1901,0.3457,0.3336,0.2692,
      ↪ 0.1208,0.0906,0.2823,0.0064'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'cP84'

_symmetry_space_group_name_H-M "P 21/a -3"
_symmetry_Int_Tables_number 205

_cell_length_a    9.00800
_cell_length_b    9.00800
_cell_length_c    9.00800
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z
3 -x,y+1/2,-z+1/2
4 -x+1/2,-y,z+1/2
5 y,z,x
6 y+1/2,-z+1/2,-x
7 -y,z+1/2,-x+1/2
8 -y+1/2,-z,x+1/2
9 z,x,y
10 z+1/2,-x+1/2,-y
11 -z,x+1/2,-y+1/2
12 -z+1/2,-x,y+1/2
13 -x,-y,-z
14 -x+1/2,y+1/2,z
15 x,-y+1/2,z+1/2
16 x+1/2,y,-z+1/2
17 -y,-z,-x
18 -y+1/2,z+1/2,x
19 y,-z+1/2,x+1/2
20 y+1/2,z,-x+1/2
21 -z,-x,-y
22 -z+1/2,x+1/2,y
23 z,-x+1/2,y+1/2
24 z+1/2,x,-y+1/2

loop_
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_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ca1 Ca 4 a 0.00000 0.00000 0.00000 1.00000
Ca2 Ca 8 c 0.37305 0.37305 0.37305 1.00000
B1 B 24 d 0.11890 0.19010 0.34570 1.00000
O1 O 24 d 0.33360 0.26920 0.12080 1.00000
O2 O 24 d 0.09060 0.28230 0.00640 1.00000
```

CaB_2O_4 (IV): A2BC4_cP84_205_d_ac_2d - POSCAR

```
A2BC4_cP84_205_d_ac_2d & a,x2,x3,y3,z3,x4,y4,z4,x5,y5,z5 --params=9.008,
      ↪ 0.37305,0.1189,0.1901,0.3457,0.3336,0.2692,0.1208,0.0906,0.2823
      ↪ ,0.0064 & Pa-3 T_{h}^{6} #205 (acd^3) & cP84 & None & B2CaO4 &
      ↪ B2CaO4 & M. Marezio and J. P. Remeika and P. D. Dernier, Acta
      ↪ Crystallogr. Sect. B Struct. Sci. 25, 965-970 (1969)
   1.00000000000000
   9.00800000000000   0.00000000000000   0.00000000000000
   0.00000000000000   9.00800000000000   0.00000000000000
   0.00000000000000   0.00000000000000   9.00800000000000
     B    Ca    O
    24    12    48
Direct
   0.11890000000000   0.19010000000000   0.34570000000000   B   (24d)
   0.38110000000000  -0.19010000000000   0.84570000000000   B   (24d)
  -0.11890000000000   0.69010000000000   0.15430000000000   B   (24d)
   0.61890000000000   0.30990000000000  -0.34570000000000   B   (24d)
   0.34570000000000   0.11890000000000   0.19010000000000   B   (24d)
   0.84570000000000   0.38110000000000  -0.19010000000000   B   (24d)
   0.15430000000000  -0.11890000000000   0.69010000000000   B   (24d)
  -0.34570000000000   0.61890000000000   0.30990000000000   B   (24d)
   0.19010000000000   0.34570000000000   0.11890000000000   B   (24d)
  -0.19010000000000   0.84570000000000   0.38110000000000   B   (24d)
   0.69010000000000   0.15430000000000  -0.11890000000000   B   (24d)
   0.30990000000000  -0.34570000000000   0.61890000000000   B   (24d)
  -0.11890000000000  -0.19010000000000  -0.34570000000000   B   (24d)
   0.61890000000000   0.19010000000000   0.15430000000000   B   (24d)
   0.11890000000000   0.30990000000000   0.84570000000000   B   (24d)
   0.38110000000000   0.69010000000000   0.34570000000000   B   (24d)
  -0.34570000000000  -0.11890000000000  -0.19010000000000   B   (24d)
   0.15430000000000   0.61890000000000   0.19010000000000   B   (24d)
   0.84570000000000   0.11890000000000   0.30990000000000   B   (24d)
   0.34570000000000   0.38110000000000   0.69010000000000   B   (24d)
  -0.19010000000000  -0.34570000000000  -0.11890000000000   B   (24d)
   0.19010000000000   0.15430000000000   0.61890000000000   B   (24d)
   0.30990000000000   0.84570000000000   0.11890000000000   B   (24d)
   0.69010000000000   0.34570000000000   0.38110000000000   B   (24d)
   0.00000000000000   0.00000000000000   0.00000000000000   Ca   (4a)
   0.50000000000000   0.00000000000000   0.50000000000000   Ca   (4a)
   0.00000000000000   0.50000000000000   0.50000000000000   Ca   (4a)
   0.50000000000000   0.50000000000000   0.00000000000000   Ca   (4a)
   0.37305000000000   0.37305000000000   0.37305000000000   Ca   (8c)
   0.12695000000000  -0.37305000000000   0.87305000000000   Ca   (8c)
  -0.37305000000000   0.87305000000000   0.12695000000000   Ca   (8c)
   0.87305000000000   0.12695000000000  -0.37305000000000   Ca   (8c)
  -0.37305000000000  -0.37305000000000  -0.37305000000000   Ca   (8c)
   0.87305000000000   0.37305000000000   0.12695000000000   Ca   (8c)
   0.37305000000000   0.12695000000000   0.87305000000000   Ca   (8c)
   0.12695000000000   0.87305000000000   0.37305000000000   Ca   (8c)
   0.33360000000000   0.26920000000000   0.12080000000000   O   (24d)
   0.16640000000000  -0.26920000000000   0.62080000000000   O   (24d)
  -0.33360000000000   0.76920000000000   0.37920000000000   O   (24d)
   0.83360000000000   0.23080000000000  -0.12080000000000   O   (24d)
   0.12080000000000   0.33360000000000   0.26920000000000   O   (24d)
   0.62080000000000   0.16640000000000  -0.26920000000000   O   (24d)
   0.37920000000000  -0.33360000000000   0.76920000000000   O   (24d)
  -0.12080000000000   0.83360000000000   0.23080000000000   O   (24d)
   0.26920000000000   0.12080000000000   0.33360000000000   O   (24d)
  -0.26920000000000   0.62080000000000   0.16640000000000   O   (24d)
   0.76920000000000   0.37920000000000  -0.33360000000000   O   (24d)
   0.23080000000000  -0.12080000000000   0.83360000000000   O   (24d)
  -0.33360000000000  -0.26920000000000  -0.12080000000000   O   (24d)
   0.83360000000000   0.26920000000000   0.37920000000000   O   (24d)
   0.33360000000000   0.23080000000000   0.62080000000000   O   (24d)
   0.16640000000000   0.76920000000000   0.12080000000000   O   (24d)
  -0.12080000000000  -0.33360000000000  -0.26920000000000   O   (24d)
   0.37920000000000   0.83360000000000   0.26920000000000   O   (24d)
   0.62080000000000   0.33360000000000   0.23080000000000   O   (24d)
   0.12080000000000   0.16640000000000   0.76920000000000   O   (24d)
  -0.26920000000000  -0.12080000000000  -0.33360000000000   O   (24d)
   0.26920000000000   0.37920000000000   0.83360000000000   O   (24d)
   0.23080000000000   0.62080000000000   0.33360000000000   O   (24d)
   0.76920000000000   0.12080000000000   0.16640000000000   O   (24d)
   0.09060000000000   0.28230000000000   0.00640000000000   O   (24d)
   0.40940000000000  -0.28230000000000   0.50640000000000   O   (24d)
  -0.09060000000000   0.78230000000000   0.49360000000000   O   (24d)
   0.59060000000000   0.21770000000000  -0.00640000000000   O   (24d)
   0.00640000000000   0.09060000000000   0.28230000000000   O   (24d)
   0.50640000000000   0.40940000000000  -0.28230000000000   O   (24d)
   0.49360000000000  -0.09060000000000   0.78230000000000   O   (24d)
  -0.00640000000000   0.59060000000000   0.21770000000000   O   (24d)
   0.28230000000000   0.00640000000000   0.09060000000000   O   (24d)
  -0.28230000000000   0.50640000000000   0.40940000000000   O   (24d)
   0.78230000000000   0.49360000000000  -0.09060000000000   O   (24d)
   0.21770000000000  -0.00640000000000   0.59060000000000   O   (24d)
  -0.09060000000000  -0.28230000000000  -0.00640000000000   O   (24d)
   0.59060000000000   0.28230000000000   0.49360000000000   O   (24d)
   0.09060000000000   0.21770000000000   0.50640000000000   O   (24d)
   0.40940000000000   0.78230000000000   0.00640000000000   O   (24d)
  -0.00640000000000  -0.09060000000000  -0.28230000000000   O   (24d)
   0.49360000000000   0.59060000000000   0.28230000000000   O   (24d)
   0.50640000000000   0.09060000000000   0.21770000000000   O   (24d)
   0.00640000000000   0.40940000000000   0.78230000000000   O   (24d)
  -0.28230000000000  -0.00640000000000  -0.09060000000000   O   (24d)
   0.28230000000000   0.49360000000000   0.59060000000000   O   (24d)
   0.21770000000000   0.50640000000000   0.09060000000000   O   (24d)
   0.78230000000000   0.00640000000000   0.40940000000000   O   (24d)
```

SnI_4 ($D1_1$): A4B_cP40_205_cd_c - CIF

```
# CIF file
```

```
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'I4Sn'
_chemical_formula_sum 'I4 Sn'

loop_
_publ_author_name
 'F. Meller'
 'I. Fankuchen'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 8
_journal_year 1955
_journal_page_first 343
_journal_page_last 344
_publ_Section_title
;
 The crystal structure of tin tetraiodide
;

# Found in The pressure-induced metallic amorphous state of SnI${4}$.
      ↪ I. A novel crystal-to-amorphous transition studied by X-ray
      ↪ scattering, 1985

_aflow_title 'SnI${4}$ (${D1_{1}}$) Structure'
_aflow_proto 'A4B_cP40_205_cd_c'
_aflow_params 'a,x_{1},x_{2},x_{3},y_{3},z_{3}'
_aflow_params_values '12.26,0.252,0.125,-0.002,-0.002,0.252'
_aflow_Strukturbericht '${D1_{1}}$'
_aflow_Pearson 'cP40'

_symmetry_space_group_name_H-M "P 21/a -3"
_symmetry_Int_Tables_number 205

_cell_length_a    12.26000
_cell_length_b    12.26000
_cell_length_c    12.26000
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z
3 -x,y+1/2,-z+1/2
4 -x+1/2,-y,z+1/2
5 y,z,x
6 y+1/2,-z+1/2,-x
7 -y,z+1/2,-x+1/2
8 -y+1/2,-z,x+1/2
9 z,x,y
10 z+1/2,-x+1/2,-y
11 -z,x+1/2,-y+1/2
12 -z+1/2,-x,y+1/2
13 -x,-y,-z
14 -x+1/2,y+1/2,z
15 x,-y+1/2,z+1/2
16 x+1/2,y,-z+1/2
17 -y,-z,-x
18 -y+1/2,z+1/2,x
19 y,-z+1/2,x+1/2
20 y+1/2,z,-x+1/2
21 -z,-x,-y
22 -z+1/2,x+1/2,y
23 z,-x+1/2,y+1/2
24 z+1/2,x,-y+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
I1 I 8 c 0.25200 0.25200 0.25200 1.00000
Sn1 Sn 8 c 0.12500 0.12500 0.12500 1.00000
I2 I 24 d -0.00200 -0.00200 0.25200 1.00000
```

SnI_4 ($D1_1$): A4B_cP40_205_cd_c - POSCAR

```
A4B_cP40_205_cd_c & a,x1,x2,x3,y3,z3 --params=12.26,0.252,0.125,-0.002,-
      ↪ 0.002,0.252 & Pa-3 T_{h}^{6} #205 (c^2d) & cP40 & ${D1_{1}}$ &
      ↪ I4Sn & I4Sn & F. Meller and I. Fankuchen, Acta Cryst. 8,
      ↪ 343-344 (1955)
   1.00000000000000
  12.26000000000000    0.00000000000000    0.00000000000000
   0.00000000000000   12.26000000000000    0.00000000000000
   0.00000000000000    0.00000000000000   12.26000000000000
     I     Sn
    32      8
Direct
   0.25200000000000    0.25200000000000    0.25200000000000     I    (8c)
   0.24800000000000   -0.25200000000000    0.75200000000000     I    (8c)
  -0.25200000000000    0.75200000000000    0.24800000000000     I    (8c)
   0.75200000000000    0.24800000000000   -0.25200000000000     I    (8c)
  -0.25200000000000   -0.25200000000000   -0.25200000000000     I    (8c)
   0.75200000000000    0.25200000000000    0.24800000000000     I    (8c)
   0.25200000000000    0.24800000000000    0.75200000000000     I    (8c)
   0.24800000000000    0.75200000000000    0.25200000000000     I    (8c)
  -0.00200000000000   -0.00200000000000    0.25200000000000     I    (24d)
   0.50200000000000    0.00200000000000    0.75200000000000     I    (24d)
   0.00200000000000    0.49800000000000    0.24800000000000     I    (24d)
   0.49800000000000    0.50200000000000   -0.25200000000000     I    (24d)
   0.25200000000000   -0.00200000000000   -0.00200000000000     I    (24d)
   0.75200000000000    0.50200000000000    0.00200000000000     I    (24d)
   0.24800000000000    0.00200000000000    0.49800000000000     I    (24d)
  -0.25200000000000    0.49800000000000    0.50200000000000     I    (24d)
  -0.00200000000000    0.25200000000000   -0.00200000000000     I    (24d)
   0.00200000000000    0.75200000000000    0.50200000000000     I    (24d)
   0.49800000000000    0.24800000000000    0.00200000000000     I    (24d)
   0.50200000000000   -0.25200000000000    0.49800000000000     I    (24d)
   0.00200000000000    0.00200000000000   -0.25200000000000     I    (24d)
   0.49800000000000   -0.00200000000000    0.24800000000000     I    (24d)
  -0.00200000000000    0.50200000000000    0.75200000000000     I    (24d)
   0.50200000000000    0.49800000000000    0.25200000000000     I    (24d)
  -0.25200000000000    0.00200000000000    0.00200000000000     I    (24d)
   0.24800000000000    0.49800000000000   -0.00200000000000     I    (24d)
   0.75200000000000   -0.00200000000000    0.50200000000000     I    (24d)
   0.25200000000000    0.50200000000000    0.49800000000000     I    (24d)
   0.00200000000000   -0.25200000000000    0.00200000000000     I    (24d)
  -0.00200000000000    0.24800000000000    0.49800000000000     I    (24d)
   0.50200000000000    0.75200000000000   -0.00200000000000     I    (24d)
   0.49800000000000    0.25200000000000    0.50200000000000     I    (24d)
   0.12500000000000    0.12500000000000    0.12500000000000    Sn    (8c)
   0.37500000000000   -0.12500000000000    0.62500000000000    Sn    (8c)
  -0.12500000000000    0.62500000000000    0.37500000000000    Sn    (8c)
   0.62500000000000    0.37500000000000   -0.12500000000000    Sn    (8c)
  -0.12500000000000   -0.12500000000000   -0.12500000000000    Sn    (8c)
   0.62500000000000    0.12500000000000    0.37500000000000    Sn    (8c)
   0.12500000000000    0.37500000000000    0.62500000000000    Sn    (8c)
   0.37500000000000    0.62500000000000    0.12500000000000    Sn    (8c)
```

$NaSbF_6$: A6BC_cP32_205_d_b_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'F6NaSb'
_chemical_formula_sum 'F6 Na Sb'

loop_
_publ_author_name
 'N. Schrewelius'
_journal_name_full_name
;
 Zeitschrift fur Anorganische und Allgemeine Chemie
;
_journal_volume 238
_journal_year 1938
_journal_page_first 241
_journal_page_last 254
_publ_Section_title
;
 R\"{o}ntgenuntersuchung der Verbindungen NaSb(OH)$_{6}$, NaSbF$_{6}$,
      ↪ NaSbO$_{3}$ und gleichartiger Stoffe
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'NaSbF$_{6}$ Structure'
_aflow_proto 'A6BC_cP32_205_d_b_a'
_aflow_params 'a,x_{3},y_{3},z_{3}'
_aflow_params_values '8.18,-0.05,0.05,0.225'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'cP32'

_symmetry_space_group_name_H-M "P 21/a -3"
_symmetry_Int_Tables_number 205

_cell_length_a    8.18000
_cell_length_b    8.18000
_cell_length_c    8.18000
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z
3 -x,y+1/2,-z+1/2
4 -x+1/2,-y,z+1/2
5 y,z,x
6 y+1/2,-z+1/2,-x
7 -y,z+1/2,-x+1/2
8 -y+1/2,-z,x+1/2
9 z,x,y
10 z+1/2,-x+1/2,-y
11 -z,x+1/2,-y+1/2
12 -z+1/2,-x,y+1/2
13 -x,-y,-z
14 -x+1/2,y+1/2,z
15 x,-y+1/2,z+1/2
16 x+1/2,y,-z+1/2
17 -y,-z,-x
18 -y+1/2,z+1/2,x
19 y,-z+1/2,x+1/2
20 y+1/2,z,-x+1/2
21 -z,-x,-y
22 -z+1/2,x+1/2,y
23 z,-x+1/2,y+1/2
```

```
24  z+1/2,x,-y+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Sb1  Sb  4  a  0.00000  0.00000  0.00000  1.00000
Na1  Na  4  b  0.50000  0.50000  0.50000  1.00000
F1  F  24  d  -0.05000  0.05000  0.22500  1.00000
```

$NaSbF_6$: A6BC_cP32_205_d_b_a - POSCAR

```
A6BC_cP32_205_d_b_a & a,x3,y3,z3 --params=8.18,-0.05,0.05,0.225 & Pa-3
      ↪ T_{h}^{6} #205 (abd) & cP32 & None & F6NaSb & F6NaSb & N.
      ↪ Schrewelius, Z. Anorg. Allg. Chem. 238, 241-254 (1938)
   1.00000000000000
   8.18000000000000    0.00000000000000    0.00000000000000
   0.00000000000000    8.18000000000000    0.00000000000000
   0.00000000000000    0.00000000000000    8.18000000000000
     F     Na     Sb
    24      4      4
Direct
  -0.05000000000000    0.05000000000000    0.22500000000000     F   (24d)
   0.55000000000000   -0.05000000000000    0.72500000000000     F   (24d)
   0.05000000000000    0.55000000000000    0.27500000000000     F   (24d)
   0.45000000000000    0.45000000000000   -0.22500000000000     F   (24d)
   0.22500000000000   -0.05000000000000    0.05000000000000     F   (24d)
   0.72500000000000    0.55000000000000   -0.05000000000000     F   (24d)
   0.27500000000000    0.05000000000000    0.55000000000000     F   (24d)
  -0.22500000000000    0.45000000000000    0.45000000000000     F   (24d)
   0.05000000000000    0.22500000000000   -0.05000000000000     F   (24d)
  -0.05000000000000    0.72500000000000    0.55000000000000     F   (24d)
   0.55000000000000    0.27500000000000    0.05000000000000     F   (24d)
   0.45000000000000   -0.22500000000000    0.45000000000000     F   (24d)
   0.05000000000000   -0.05000000000000   -0.22500000000000     F   (24d)
   0.45000000000000    0.05000000000000    0.27500000000000     F   (24d)
  -0.05000000000000    0.45000000000000    0.72500000000000     F   (24d)
   0.55000000000000    0.55000000000000    0.22500000000000     F   (24d)
  -0.22500000000000    0.05000000000000   -0.05000000000000     F   (24d)
   0.27500000000000    0.45000000000000    0.05000000000000     F   (24d)
   0.72500000000000   -0.05000000000000    0.45000000000000     F   (24d)
   0.22500000000000    0.55000000000000    0.55000000000000     F   (24d)
  -0.05000000000000   -0.22500000000000    0.05000000000000     F   (24d)
   0.05000000000000    0.27500000000000    0.45000000000000     F   (24d)
   0.45000000000000    0.72500000000000   -0.05000000000000     F   (24d)
   0.55000000000000    0.22500000000000    0.55000000000000     F   (24d)
   0.50000000000000    0.50000000000000    0.50000000000000    Na   (4b)
   0.00000000000000    0.50000000000000    0.00000000000000    Na   (4b)
   0.50000000000000    0.00000000000000    0.00000000000000    Na   (4b)
   0.00000000000000    0.00000000000000    0.50000000000000    Na   (4b)
   0.00000000000000    0.00000000000000    0.00000000000000    Sb   (4a)
   0.50000000000000    0.00000000000000    0.50000000000000    Sb   (4a)
   0.00000000000000    0.50000000000000    0.50000000000000    Sb   (4a)
   0.50000000000000    0.50000000000000    0.00000000000000    Sb   (4a)
```

ZrP_2O_7 High-Temperature ($K6_1$): A7B2C_cP40_205_bd_c_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'P2O7Zr'
_chemical_formula_sum 'O7 P2 Zr'

loop_
_publ_author_name
 'G. R. Levi'
 'G. Peyronel'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 92
_journal_year 1935
_journal_page_first 190
_journal_page_last 209
_publ_Section_title
;
 Struttura Cristallografica del Gruppo Isomorfo (Si$^{4+}$, Ti$^{4+}$,
      ↪ Zr$^{4+}$, Sn$^{4+}$, Hf$^{4+}$) P$_{2}$O$_{7}$
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'ZrP$_{2}$O$_{7}$ High-Temperature ($K6_{1}$) Structure'
_aflow_proto 'A7B2C_cP40_205_bd_c_a'
_aflow_params 'a,x_{3},x_{4},y_{4},z_{4}'
_aflow_params_values '8.2,0.39,0.394,0.218,0.458'
_aflow_Strukturbericht '$K6_{1}$'
_aflow_Pearson 'cP40'

_symmetry_space_group_name_H-M "P 21/a -3"
_symmetry_Int_Tables_number 205

_cell_length_a     8.20000
_cell_length_b     8.20000
_cell_length_c     8.20000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1  x,y,z
2  x+1/2,-y+1/2,-z
3  -x,y+1/2,-z+1/2
4  -x+1/2,-y,z+1/2
5  y,z,x
6  y+1/2,-z+1/2,-x
7  -y,z+1/2,-x+1/2
8  -y+1/2,-z,x+1/2
9  z,x,y
10  z+1/2,-x+1/2,-y
11  -z,x+1/2,-y+1/2
12  -z+1/2,-x,y+1/2
13  -x,-y,-z
14  -x+1/2,y+1/2,z
15  x,-y+1/2,z+1/2
16  x+1/2,y,-z+1/2
17  -y,-z,-x
18  -y+1/2,z+1/2,x
19  y,-z+1/2,x+1/2
20  y+1/2,z,-x+1/2
21  -z,-x,-y
22  -z+1/2,x+1/2,y
23  z,-x+1/2,y+1/2
24  z+1/2,x,-y+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Zr1  Zr  4  a  0.00000  0.00000  0.00000  1.00000
O1  O  4  b  0.50000  0.50000  0.50000  1.00000
P1  P  8  c  0.39000  0.39000  0.39000  1.00000
O2  O  24  d  0.39400  0.21800  0.45800  1.00000
```

ZrP_2O_7 High-Temperature ($K6_1$): A7B2C_cP40_205_bd_c_a - POSCAR

```
A7B2C_cP40_205_bd_c_a & a,x3,x4,y4,z4 --params=8.2,0.39,0.394,0.218,
      ↪ 0.458 & Pa-3 T_{h}^{6} #205 (abcd) & cP40 & $K6_{1}$ & P2O7Zr &
      ↪ P2O7Zr & G. R. Levi and G. Peyronel, Zeitschrift f{"u}r
      ↪ Kristallographie - Crystalline Materials 92, 190-209 (1935)
   1.00000000000000
   8.20000000000000    0.00000000000000    0.00000000000000
   0.00000000000000    8.20000000000000    0.00000000000000
   0.00000000000000    0.00000000000000    8.20000000000000
     O      P     Zr
    28      8      4
Direct
   0.50000000000000    0.50000000000000    0.50000000000000     O   (4b)
   0.00000000000000    0.50000000000000    0.00000000000000     O   (4b)
   0.50000000000000    0.00000000000000    0.00000000000000     O   (4b)
   0.00000000000000    0.00000000000000    0.50000000000000     O   (4b)
   0.39400000000000    0.21800000000000    0.45800000000000     O  (24d)
   0.10600000000000   -0.21800000000000    0.95800000000000     O  (24d)
  -0.39400000000000    0.71800000000000    0.04200000000000     O  (24d)
   0.89400000000000    0.28200000000000   -0.45800000000000     O  (24d)
   0.45800000000000    0.39400000000000    0.21800000000000     O  (24d)
   0.95800000000000    0.10600000000000   -0.21800000000000     O  (24d)
   0.04200000000000   -0.39400000000000    0.71800000000000     O  (24d)
  -0.45800000000000    0.89400000000000    0.28200000000000     O  (24d)
   0.21800000000000    0.45800000000000    0.39400000000000     O  (24d)
  -0.21800000000000    0.95800000000000    0.10600000000000     O  (24d)
   0.71800000000000    0.04200000000000   -0.39400000000000     O  (24d)
   0.28200000000000   -0.45800000000000    0.89400000000000     O  (24d)
  -0.39400000000000   -0.21800000000000   -0.45800000000000     O  (24d)
   0.89400000000000    0.21800000000000    0.04200000000000     O  (24d)
   0.39400000000000    0.28200000000000    0.95800000000000     O  (24d)
   0.10600000000000    0.71800000000000    0.45800000000000     O  (24d)
  -0.45800000000000   -0.39400000000000   -0.21800000000000     O  (24d)
   0.04200000000000    0.89400000000000    0.21800000000000     O  (24d)
   0.95800000000000    0.39400000000000    0.28200000000000     O  (24d)
   0.45800000000000    0.10600000000000    0.71800000000000     O  (24d)
  -0.21800000000000   -0.45800000000000   -0.39400000000000     O  (24d)
   0.21800000000000    0.04200000000000    0.89400000000000     O  (24d)
   0.28200000000000    0.95800000000000    0.39400000000000     O  (24d)
   0.71800000000000    0.45800000000000    0.10600000000000     O  (24d)
   0.39000000000000    0.39000000000000    0.39000000000000     P   (8c)
   0.11000000000000   -0.39000000000000    0.89000000000000     P   (8c)
  -0.39000000000000    0.89000000000000    0.11000000000000     P   (8c)
   0.89000000000000    0.11000000000000   -0.39000000000000     P   (8c)
  -0.39000000000000   -0.39000000000000   -0.39000000000000     P   (8c)
   0.89000000000000    0.39000000000000    0.11000000000000     P   (8c)
   0.39000000000000    0.11000000000000    0.89000000000000     P   (8c)
   0.11000000000000    0.89000000000000    0.39000000000000     P   (8c)
   0.00000000000000    0.00000000000000    0.00000000000000    Zr   (4a)
   0.50000000000000    0.00000000000000    0.50000000000000    Zr   (4a)
   0.00000000000000    0.50000000000000    0.50000000000000    Zr   (4a)
   0.50000000000000    0.50000000000000    0.00000000000000    Zr   (4a)
```

$NaCr(SO_4)_2{\cdot}12H_2O$ Alum: AB12CD8E2_cP96_205_a_2d_b_cd_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Alum'
_chemical_formula_sum 'Cr (H2O)12 Na O8 S2'
```

```
loop_
_publ_author_name
 'A. H. C. Ledsham'
 'H. Steeple'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 24
_journal_year 1968
_journal_page_first 1287
_journal_page_last 1289
_publ_Section_title
;
 The crystal structure of sodium chromium alum and caesium chromium alum
;

_aflow_title 'NaCr(SO$_{4}$)$_{2}$$\cdot$12H$_{2}$O Alum Structure'
_aflow_proto 'AB12CD8E2_cP96_205_a_2d_b_cd_c'
_aflow_params 'a,x_{3},x_{4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{6},x_{7},
      ↪ y_{7},z_{7}'
_aflow_params_values '12.4,0.239,0.31,0.158,0.014,0.018,0.042,0.136,
      ↪ 0.302,0.307,0.224,-0.08'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'cP96'

_symmetry_space_group_name_H-M "P 21/a -3"
_symmetry_Int_Tables_number 205

_cell_length_a     12.40000
_cell_length_b     12.40000
_cell_length_c     12.40000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z
3 -x,y+1/2,-z+1/2
4 -x+1/2,-y,z+1/2
5 y,z,x
6 y+1/2,-z+1/2,-x
7 -y,z+1/2,-x+1/2
8 -y+1/2,-z,x+1/2
9 z,x,y
10 z+1/2,-x+1/2,-y
11 -z,x+1/2,-y+1/2
12 -z+1/2,-x,y+1/2
13 -x,-y,-z
14 -x+1/2,y+1/2,z
15 x,-y+1/2,z+1/2
16 x+1/2,y,-z+1/2
17 -y,-z,-x
18 -y+1/2,z+1/2,x
19 y,-z+1/2,x+1/2
20 y+1/2,z,-x+1/2
21 -z,-x,-y
22 -z+1/2,x+1/2,y
23 z,-x+1/2,y+1/2
24 z+1/2,x,-y+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cr1 Cr 4 a 0.00000 0.00000 0.00000 1.00000
Na1 Na 4 b 0.50000 0.50000 0.50000 1.00000
O1 O 8 c 0.23900 0.23900 0.23900 1.00000
S1 S 8 c 0.31000 0.31000 0.31000 1.00000
H2O1 H2O 24 d 0.15800 0.01400 0.01800 1.00000
H2O2 H2O 24 d 0.04200 0.13600 0.30200 1.00000
O2 O 24 d 0.30700 0.22400 -0.0800 1.00000
```

$NaCr(SO_4)_2{\cdot}12H_2O$ Alum: AB12CD8E2_cP96_205_a_2d_b_cd_c - POSCAR

```
AB12CD8E2_cP96_205_a_2d_b_cd_c & a,x3,x4,x5,y5,z5,x6,y6,z6,x7,y7,z7 --
      ↪ params=12.4,0.239,0.31,0.158,0.014,0.018,0.042,0.136,0.302,
      ↪ 0.307,0.224,-0.08 & Pa-3 T_{h}^{6} #205 (abc^2d^3) & cP96 &
      ↪ None & Cr(H2O)12NaO8S2 & Alum & A. H. C. Ledsham and H. Steeple
      ↪ , Acta Crystallogr. Sect. B Struct. Sci. 24, 1287-1289 (1968)
   1.00000000000000
  12.40000000000000     0.00000000000000     0.00000000000000
   0.00000000000000    12.40000000000000     0.00000000000000
   0.00000000000000     0.00000000000000    12.40000000000000
     Cr    H2O     Na      O      S
      4     48      4     32      8
Direct
   0.00000000000000     0.00000000000000     0.00000000000000   Cr    (4a)
   0.50000000000000     0.00000000000000     0.50000000000000   Cr    (4a)
   0.00000000000000     0.50000000000000     0.50000000000000   Cr    (4a)
   0.50000000000000     0.50000000000000     0.00000000000000   Cr    (4a)
   0.15800000000000     0.01400000000000     0.01800000000000   H2O   (24d)
   0.34200000000000    -0.01400000000000     0.51800000000000   H2O   (24d)
  -0.15800000000000     0.51400000000000     0.48200000000000   H2O   (24d)
   0.65800000000000     0.48600000000000    -0.01800000000000   H2O   (24d)
   0.01800000000000     0.15800000000000     0.01400000000000   H2O   (24d)
   0.51800000000000     0.34200000000000    -0.01400000000000   H2O   (24d)
   0.48200000000000    -0.15800000000000     0.51400000000000   H2O   (24d)
  -0.01800000000000     0.65800000000000     0.48600000000000   H2O   (24d)
   0.01400000000000     0.01800000000000     0.15800000000000   H2O   (24d)
  -0.01400000000000     0.51800000000000     0.34200000000000   H2O   (24d)
   0.51400000000000     0.48200000000000    -0.15800000000000   H2O   (24d)
   0.48600000000000    -0.01800000000000     0.65800000000000   H2O   (24d)
  -0.15800000000000    -0.01400000000000    -0.01800000000000   H2O   (24d)
   0.65800000000000     0.01400000000000     0.48200000000000   H2O   (24d)
   0.15800000000000     0.48600000000000     0.51800000000000   H2O   (24d)
   0.34200000000000     0.51400000000000     0.01800000000000   H2O   (24d)
  -0.01800000000000    -0.15800000000000    -0.01400000000000   H2O   (24d)
   0.48200000000000     0.65800000000000     0.01400000000000   H2O   (24d)
   0.51800000000000     0.15800000000000     0.48600000000000   H2O   (24d)
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   0.50000000000000     0.50000000000000     0.50000000000000   Na    (4b)
   0.00000000000000     0.50000000000000     0.00000000000000   Na    (4b)
   0.50000000000000     0.00000000000000     0.00000000000000   Na    (4b)
   0.00000000000000     0.00000000000000     0.50000000000000   Na    (4b)
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   0.26100000000000    -0.23900000000000     0.73900000000000   O     (8c)
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   0.73900000000000     0.26100000000000    -0.23900000000000   O     (8c)
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   0.23900000000000     0.26100000000000     0.73900000000000   O     (8c)
   0.26100000000000     0.73900000000000     0.23900000000000   O     (8c)
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   0.19300000000000    -0.22400000000000     0.42000000000000   O     (24d)
  -0.30700000000000     0.72400000000000     0.58000000000000   O     (24d)
   0.80700000000000     0.27600000000000     0.08000000000000   O     (24d)
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   0.08000000000000     0.80700000000000     0.27600000000000   O     (24d)
   0.22400000000000    -0.08000000000000     0.30700000000000   O     (24d)
  -0.22400000000000     0.42000000000000     0.19300000000000   O     (24d)
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  -0.30700000000000    -0.22400000000000     0.08000000000000   O     (24d)
   0.80700000000000     0.22400000000000     0.58000000000000   O     (24d)
   0.30700000000000     0.27600000000000     0.42000000000000   O     (24d)
   0.19300000000000     0.72400000000000    -0.08000000000000   O     (24d)
   0.08000000000000    -0.30700000000000    -0.22400000000000   O     (24d)
   0.58000000000000     0.80700000000000     0.22400000000000   O     (24d)
   0.42000000000000     0.30700000000000     0.27600000000000   O     (24d)
  -0.08000000000000     0.19300000000000     0.72400000000000   O     (24d)
  -0.22400000000000     0.08000000000000    -0.30700000000000   O     (24d)
   0.22400000000000     0.58000000000000     0.80700000000000   O     (24d)
   0.27600000000000     0.42000000000000     0.30700000000000   O     (24d)
   0.72400000000000    -0.08000000000000     0.19300000000000   O     (24d)
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   0.19000000000000    -0.31000000000000     0.81000000000000   S     (8c)
  -0.31000000000000     0.81000000000000     0.19000000000000   S     (8c)
   0.81000000000000     0.19000000000000    -0.31000000000000   S     (8c)
  -0.31000000000000    -0.31000000000000    -0.31000000000000   S     (8c)
   0.81000000000000     0.31000000000000     0.19000000000000   S     (8c)
   0.31000000000000     0.19000000000000     0.81000000000000   S     (8c)
   0.19000000000000     0.81000000000000     0.31000000000000   S     (8c)
```

γ-Alum [$AlNa(SO_4)_2{\cdot}12H_2O$, $H4_{15}$]: AB24CD20E2_cP192_205_a_4d_b_c3d_c - CIF

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loop_
_publ_author_name
 'D. T. Cromer'
 'M. I. Kay'
 'A. C. Larson'
_journal_name_full_name
;
 Acta Cristallographica
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;
_journal_volume 22
_journal_year 1967
_journal_page_first 182
_journal_page_last 187
_publ_Section_title
;
 Refinement of the alum structures. II. X-ray and neutron diffraction of
      ↪ NaAl(SO$_{4}$)$_{2}$$\cdot$12H$_{2}$O, $\gamma$-alum
;

# Found in The American Mineralogist Crystal Structure Database, 2003

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_aflow_params 'a,x_{3},x_{4},x_{5},y_{5},z_{5},x_{6},y_{6},z_{6},x_{7},
      ↪ y_{7},z_{7},x_{8},y_{8},z_{8},x_{9},y_{9},z_{9},x_{10},y_{10},
      ↪ z_{10},x_{11},y_{11},z_{11}'
_aflow_params_values '12.213,0.3343,0.2652,0.585,0.319,0.378,0.486,0.308
      ↪ ,0.383,0.555,0.202,0.502,0.588,0.339,0.113,0.2957,0.2783,0.1508
      ↪ ,0.0767,0.0403,0.3188,0.1371,-0.0404,0.0573'
_aflow_Strukturbericht '$H4_{15}$'
_aflow_Pearson 'cP192'

_symmetry_space_group_name_H-M "P 21/a -3"
_symmetry_Int_Tables_number 205

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_space_group_symop_operation_xyz
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2 x+1/2,-y+1/2,-z
3 -x,y+1/2,-z+1/2
4 -x+1/2,-y,z+1/2
5 y,z,x
6 y+1/2,-z+1/2,-x
7 -y,z+1/2,-x+1/2
8 -y+1/2,-z,x+1/2
9 z,x,y
10 z+1/2,-x+1/2,-y
11 -z,x+1/2,-y+1/2
12 -z+1/2,-x,y+1/2
13 -x,-y,-z
14 -x+1/2,y+1/2,z
15 x,-y+1/2,z+1/2
16 x+1/2,y,-z+1/2
17 -y,-z,-x
18 -y+1/2,z+1/2,x
19 y,-z+1/2,x+1/2
20 y+1/2,z,-x+1/2
21 -z,-x,-y
22 -z+1/2,x+1/2,y
23 z,-x+1/2,y+1/2
24 z+1/2,x,-y+1/2

loop_
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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
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_atom_site_fract_z
_atom_site_occupancy
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Na1 Na 4 b 0.5000 0.5000 0.5000 1.00000
O1 O 8 c 0.3343 0.3343 0.3343 1.00000
S1 S 8 c 0.2652 0.2652 0.2652 1.00000
H1 H 24 d 0.5850 0.3190 0.3780 1.00000
H2 H 24 d 0.4860 0.3080 0.3830 1.00000
H3 H 24 d 0.5550 0.2020 0.5020 1.00000
H4 H 24 d 0.5880 0.3390 0.1130 1.00000
O2 O 24 d 0.2957 0.2783 0.1508 1.00000
O3 O 24 d 0.0767 0.0403 0.3188 1.00000
O4 O 24 d 0.1371 -0.0404 0.0573 1.00000
```

γ-Alum [$AlNa(SO_4)_2 \cdot 12H_2O$, $H4_{15}$]: AB24CD20E2_cP192_205_a_4d_b_c3d_c - POSCAR

```
AB24CD20E2_cP192_205_a_4d_b_c3d_c & a,x3,x4,x5,y5,z5,x6,y6,z6,x7,y7,z7,
      ↪ x8,y8,z8,x9,y9,z9,x10,y10,z10,x11,y11,z11 --params=12.213,
      ↪ 0.3343,0.2652,0.585,0.319,0.378,0.486,0.308,0.383,0.555,0.202,
      ↪ 0.502,0.588,0.339,0.113,0.2957,0.2783,0.1508,0.0767,0.0403,
      ↪ 0.3188,0.1371,-0.0404,0.0573 & Pa-3 T_{h}^{6} #205 (abc^2d^7) &
      ↪  cP192 & $H4_{15}$ & AlH24NaO20S2 & $\gamma$-alum & D. T.
      ↪ Cromer and M. I. Kay and A. C. Larson, Acta Cryst. 22, 182-187
      ↪ (1967)
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  12.21300000000000     0.00000000000000     0.00000000000000
   0.00000000000000    12.21300000000000     0.00000000000000
   0.00000000000000     0.00000000000000    12.21300000000000
     Al     H    Na     O     S
      4    96     4    80     8
Direct
   0.00000000000000     0.00000000000000     0.00000000000000   Al   (4a)
   0.50000000000000     0.00000000000000     0.50000000000000   Al   (4a)
   0.00000000000000     0.50000000000000     0.50000000000000   Al   (4a)
   0.50000000000000     0.50000000000000     0.00000000000000   Al   (4a)
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  -0.08500000000000    -0.31900000000000     0.87800000000000   H   (24d)
  -0.58500000000000     0.81900000000000     0.12200000000000   H   (24d)
   1.08500000000000     0.18100000000000    -0.37800000000000   H   (24d)
   0.37800000000000     0.58500000000000     0.31900000000000   H   (24d)
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   0.12200000000000    -0.58500000000000     0.81900000000000   H   (24d)
  -0.37800000000000     1.08500000000000     0.18100000000000   H   (24d)
   0.31900000000000     0.37800000000000     0.58500000000000   H   (24d)
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   0.81900000000000     0.37800000000000    -0.08500000000000   H   (24d)
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   0.01400000000000    -0.30800000000000     0.88300000000000   H   (24d)
  -0.48600000000000     0.80800000000000     0.11700000000000   H   (24d)
   0.98600000000000     0.19200000000000    -0.38300000000000   H   (24d)
   0.38300000000000     0.48600000000000     0.30800000000000   H   (24d)
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   0.50200000000000     0.55500000000000     0.20200000000000   H   (24d)
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  -0.00200000000000    -0.55500000000000     0.70200000000000   H   (24d)
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   0.29800000000000    -0.50200000000000     1.05500000000000   H   (24d)
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   1.05500000000000     0.20200000000000    -0.00200000000000   H   (24d)
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  -0.05500000000000     0.70200000000000     0.50200000000000   H   (24d)
  -0.50200000000000    -0.55500000000000    -0.20200000000000   H   (24d)
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   0.58800000000000     0.33900000000000     0.11300000000000   H   (24d)
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  -0.58800000000000     0.83900000000000     0.38700000000000   H   (24d)
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   0.11300000000000     0.58800000000000     0.33900000000000   H   (24d)
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α-Alum [$KAl(SO_4)_2 \cdot 12H_2O$, $H4_{13}$]: AB24CD28E2_cP224_205_a_4d_b_2c4d_c - CIF

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 'J. W. Steed'
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 'V. M. Petrusevski'
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 Acta Crystallographica Section B: Structural Science
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;
 Structure of the alums. I. On the sulfate group disorder in the $\
     ↪ alpha$-alums
;

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     ↪ },y_{11},z_{11},x_{12},y_{12},z_{12},x_{13},y_{13},z_{13}'
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     ↪ 0.26302,0.41979,0.30624,0.2808,0.2,0.3614,0.15173,0.028,-
     ↪ 0.01985,0.04692,0.13071,0.30399'
_aflow_Strukturbericht '$H4_{13}$'
_aflow_Pearson 'cP224'

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4 -x+1/2,-y,z+1/2
5 y,z,x
6 y+1/2,-z+1/2,-x
7 -y,z+1/2,-x+1/2
8 -y+1/2,-z,x+1/2
9 z,x,y
10 z+1/2,-x+1/2,-y
11 -z,x+1/2,-y+1/2
12 -z+1/2,-x,y+1/2
13 -x,-y,-z
14 -x+1/2,y+1/2,z
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16 x+1/2,y,-z+1/2
17 -y,-z,-x
18 -y+1/2,z+1/2,x
19 y,-z+1/2,x+1/2
20 y+1/2,z,-x+1/2
21 -z,-x,-y
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K1 K 4 b 0.50000 0.50000 0.50000 1.00000
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O2 O 8 c 0.37460 0.37460 0.37460 0.21300
S1 S 8 c 0.30610 0.30610 0.30610 1.00000
H1 H 24 d 0.17500 0.03650 -0.07275 1.00000
H2 H 24 d 0.19820 0.02620 0.03400 1.00000
H3 H 24 d 0.10400 0.15500 0.30000 1.00000
H4 H 24 d 0.00090 0.18210 0.29140 1.00000
O3 O 24 d 0.26302 0.41979 0.30624 0.78700
O4 O 24 d 0.28080 0.20000 0.36140 0.21300
O5 O 24 d 0.15173 0.02800 -0.01985 1.00000
O6 O 24 d 0.04692 0.13071 0.30399 1.00000
```

α-Alum [$KAl(SO_4)_2 \cdot 12H_2O$, $H4_{13}$]: AB24CD28E2_cP224_205_a_4d_b_2c4d_c - POSCAR

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     ↪ z8,x9,y9,z9,x10,y10,z10,x11,y11,z11,x12,y12,z12,x13,y13,z13 --
     ↪ params=12.135,0.23606,0.3746,0.3061,0.175,0.0365,-0.07275,
     ↪ 0.1982,0.0262,0.034,0.104,0.155,0.3,0.0009,0.1821,0.2914,
     ↪ 0.26302,0.41979,0.30624,0.2808,0.2,0.3614,0.15173,0.028,-
     ↪ 0.01985,0.04692,0.13071,0.30399 & Pa-3 T_{h}^{6} #205 (abc^3d^8
     ↪ ) & cP224 & $H4_{13}$ & AlH24KO20S2 & $\alpha$-alum & S. C.
     ↪ Nyburg et al., Acta Crystallogr. Sect. B Struct. Sci. 56,
     ↪ 204-209 (2000)
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   0.00000000000000  12.13500000000000   0.00000000000000
   0.00000000000000   0.00000000000000  12.13500000000000
    Al    H    K    O    S
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Direct
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β-Alum [$Al(NH_3CH_3)_2(SO_4)_2{\cdot}12H_2O$, $H4_{14}$]: AB2C36D2E20F2_cP252_205_a_c_6d_c_c3d_c - CIF

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 'G. Will'
 'W. Sch\"{a}fer'
 'A. Kirfel'
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 X-Ray and neutron diffraction study of alums
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O1 O 8 c 0.2341 0.2341 0.2341 1.00000
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O2 O 24 d 0.6911 0.7530 0.0914 1.00000
O3 O 24 d 0.8521 -0.0190 0.0185 1.00000
O4 O 24 d 0.0419 0.1373 0.2975 1.00000
```

β-Alum [$Al(NH_3CH_3)_2(SO_4)_2{\cdot}12H_2O$, $H4_{14}$]: AB2C36D2E20F2_cP252_205_a_c_6d_c_c3d_c - POSCAR

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      ↪ AlC2H36N2O20S2 & $\beta$-alum & A. M. Abdeen et al.,
      ↪ Zeitschrift f{"u}r Kristallographie - Crystalline Materials 157
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−0.04230000000000	0.70180000000000	0.47280000000000	H	(24d)
−0.47280000000000	−0.54230000000000	−0.20180000000000	H	(24d)
0.02720000000000	1.04230000000000	0.20180000000000	H	(24d)
0.97280000000000	0.54230000000000	0.29820000000000	H	(24d)
0.47280000000000	−0.04230000000000	0.70180000000000	H	(24d)
−0.20180000000000	−0.47280000000000	−0.54230000000000	H	(24d)
0.20180000000000	0.02720000000000	1.04230000000000	H	(24d)
0.29820000000000	0.97280000000000	0.54230000000000	H	(24d)
0.70180000000000	0.47280000000000	−0.04230000000000	H	(24d)
0.41320000000000	0.18460000000000	0.45020000000000	H	(24d)
0.08680000000000	−0.18460000000000	0.95020000000000	H	(24d)
−0.41320000000000	0.68460000000000	0.04980000000000	H	(24d)
0.91320000000000	0.31540000000000	−0.45020000000000	H	(24d)
0.45020000000000	0.41320000000000	0.18460000000000	H	(24d)
0.95020000000000	0.08680000000000	−0.18460000000000	H	(24d)
0.04980000000000	−0.41320000000000	0.68460000000000	H	(24d)
−0.45020000000000	0.91320000000000	0.31540000000000	H	(24d)
0.18460000000000	0.45020000000000	0.41320000000000	H	(24d)
−0.18460000000000	0.95020000000000	0.08680000000000	H	(24d)
0.68460000000000	0.04980000000000	−0.41320000000000	H	(24d)
0.31540000000000	−0.45020000000000	0.91320000000000	H	(24d)
−0.41320000000000	−0.18460000000000	−0.45020000000000	H	(24d)
0.91320000000000	0.18460000000000	0.04980000000000	H	(24d)
0.41320000000000	0.31540000000000	0.95020000000000	H	(24d)
0.08680000000000	0.68460000000000	0.45020000000000	H	(24d)
−0.45020000000000	−0.41320000000000	−0.18460000000000	H	(24d)
0.04980000000000	0.91320000000000	0.18460000000000	H	(24d)
0.95020000000000	0.41320000000000	0.31540000000000	H	(24d)
0.45020000000000	0.08680000000000	0.68460000000000	H	(24d)
−0.18460000000000	−0.45020000000000	−0.41320000000000	H	(24d)
0.18460000000000	0.04980000000000	0.91320000000000	H	(24d)
0.31540000000000	0.95020000000000	0.41320000000000	H	(24d)
0.68460000000000	0.45020000000000	0.08680000000000	H	(24d)
0.21430000000000	0.49410000000000	0.19440000000000	H	(24d)
0.28570000000000	−0.49410000000000	0.69440000000000	H	(24d)
−0.21430000000000	0.99410000000000	0.30560000000000	H	(24d)
0.71430000000000	0.00590000000000	−0.19440000000000	H	(24d)
0.19440000000000	0.21430000000000	0.49410000000000	H	(24d)
0.69440000000000	0.28570000000000	−0.49410000000000	H	(24d)
0.30560000000000	−0.21430000000000	0.99410000000000	H	(24d)
−0.19440000000000	0.71430000000000	0.00590000000000	H	(24d)
0.49410000000000	0.19440000000000	0.21430000000000	H	(24d)
−0.49410000000000	0.69440000000000	0.28570000000000	H	(24d)
0.99410000000000	0.30560000000000	−0.21430000000000	H	(24d)
0.00590000000000	−0.19440000000000	0.71430000000000	H	(24d)
−0.21430000000000	−0.49410000000000	−0.19440000000000	H	(24d)
0.71430000000000	0.49410000000000	0.30560000000000	H	(24d)
0.21430000000000	0.00590000000000	0.69440000000000	H	(24d)
0.28570000000000	0.99410000000000	0.19440000000000	H	(24d)
−0.19440000000000	−0.21430000000000	−0.49410000000000	H	(24d)
0.30560000000000	0.71430000000000	0.49410000000000	H	(24d)
0.69440000000000	0.21430000000000	0.00590000000000	H	(24d)
0.19440000000000	0.28570000000000	0.99410000000000	H	(24d)
−0.49410000000000	−0.19440000000000	−0.21430000000000	H	(24d)
0.49410000000000	0.30560000000000	0.71430000000000	H	(24d)
0.00590000000000	0.69440000000000	0.21430000000000	H	(24d)
0.99410000000000	0.19440000000000	0.28570000000000	H	(24d)
0.21480000000000	0.61030000000000	0.16970000000000	H	(24d)
0.28520000000000	−0.61030000000000	0.66970000000000	H	(24d)
−0.21480000000000	1.11030000000000	0.33030000000000	H	(24d)
0.71480000000000	−0.11030000000000	−0.16970000000000	H	(24d)
0.16970000000000	0.21480000000000	0.61030000000000	H	(24d)
0.66970000000000	0.28520000000000	−0.61030000000000	H	(24d)
0.33030000000000	−0.21480000000000	1.11030000000000	H	(24d)
−0.16970000000000	0.71480000000000	−0.11030000000000	H	(24d)
0.61030000000000	0.16970000000000	0.21480000000000	H	(24d)
−0.61030000000000	0.66970000000000	0.28520000000000	H	(24d)
1.11030000000000	0.33030000000000	−0.21480000000000	H	(24d)
−0.11030000000000	−0.16970000000000	0.71480000000000	H	(24d)
−0.21480000000000	−0.61030000000000	−0.16970000000000	H	(24d)
0.71480000000000	0.61030000000000	0.33030000000000	H	(24d)
0.21480000000000	−0.11030000000000	0.66970000000000	H	(24d)
0.28520000000000	1.11030000000000	0.16970000000000	H	(24d)
−0.16970000000000	−0.21480000000000	−0.61030000000000	H	(24d)
0.33030000000000	0.71480000000000	0.61030000000000	H	(24d)
0.66970000000000	0.21480000000000	−0.11030000000000	H	(24d)
0.16970000000000	0.28520000000000	1.11030000000000	H	(24d)
−0.61030000000000	−0.16970000000000	−0.21480000000000	H	(24d)
0.61030000000000	0.33030000000000	0.71480000000000	H	(24d)
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1.11030000000000	0.16970000000000	0.28520000000000	H	(24d)
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−0.51560000000000	1.01560000000000	−0.01560000000000	N	(8c)
1.01560000000000	−0.01560000000000	−0.51560000000000	N	(8c)
−0.51560000000000	−0.51560000000000	−0.51560000000000	N	(8c)
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0.26590000000000	−0.23410000000000	0.73410000000000	O	(8c)
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0.73410000000000	0.26590000000000	−0.23410000000000	O	(8c)
−0.23410000000000	−0.23410000000000	−0.23410000000000	O	(8c)
0.73410000000000	0.23410000000000	0.26590000000000	O	(8c)
0.23410000000000	0.26590000000000	0.73410000000000	O	(8c)
0.26590000000000	0.73410000000000	0.23410000000000	O	(8c)
0.69110000000000	0.75300000000000	0.09140000000000	O	(24d)
−0.19110000000000	−0.75300000000000	0.59140000000000	O	(24d)
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1.19110000000000	−0.25300000000000	−0.09140000000000	O	(24d)
0.09140000000000	0.69110000000000	0.75300000000000	O	(24d)
0.59140000000000	−0.19110000000000	−0.75300000000000	O	(24d)
0.40860000000000	−0.69110000000000	1.25300000000000	O	(24d)
−0.09140000000000	1.19110000000000	−0.25300000000000	O	(24d)
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1.25300000000000	0.40860000000000	−0.69110000000000	O	(24d)
−0.25300000000000	−0.09140000000000	1.19110000000000	O	(24d)
−0.69110000000000	−0.75300000000000	−0.09140000000000	O	(24d)
1.19110000000000	0.75300000000000	0.40860000000000	O	(24d)
0.69110000000000	−0.25300000000000	0.59140000000000	O	(24d)
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0.40860000000000	1.19110000000000	0.75300000000000	O	(24d)
0.59140000000000	0.69110000000000	−0.25300000000000	O	(24d)
0.09140000000000	−0.19110000000000	1.25300000000000	O	(24d)
−0.75300000000000	−0.09140000000000	−0.69110000000000	O	(24d)
0.75300000000000	0.40860000000000	1.19110000000000	O	(24d)
−0.25300000000000	0.59140000000000	0.69110000000000	O	(24d)
1.25300000000000	0.09140000000000	−0.19110000000000	O	(24d)
0.85210000000000	−0.01900000000000	0.01850000000000	O	(24d)
−0.35210000000000	0.01900000000000	0.51850000000000	O	(24d)
−0.85210000000000	0.48100000000000	0.48150000000000	O	(24d)
1.35210000000000	0.51900000000000	−0.01850000000000	O	(24d)
0.01850000000000	0.85210000000000	−0.01900000000000	O	(24d)
0.51850000000000	−0.35210000000000	0.01900000000000	O	(24d)
0.48150000000000	−0.85210000000000	0.48100000000000	O	(24d)
−0.01850000000000	1.35210000000000	0.51900000000000	O	(24d)
−0.01900000000000	0.01850000000000	0.85210000000000	O	(24d)
0.01900000000000	0.51850000000000	−0.35210000000000	O	(24d)
0.48100000000000	0.48150000000000	−0.85210000000000	O	(24d)
0.51900000000000	−0.01850000000000	1.35210000000000	O	(24d)
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1.35210000000000	−0.01900000000000	0.48150000000000	O	(24d)
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0.51900000000000	0.51850000000000	0.85210000000000	O	(24d)
0.48100000000000	0.01850000000000	−0.35210000000000	O	(24d)
0.04190000000000	0.13730000000000	0.29750000000000	O	(24d)
0.45810000000000	−0.13730000000000	0.79750000000000	O	(24d)
−0.04190000000000	0.63730000000000	0.20250000000000	O	(24d)
0.54190000000000	0.36270000000000	−0.29750000000000	O	(24d)
0.29750000000000	0.04190000000000	0.13730000000000	O	(24d)
0.79750000000000	0.45810000000000	−0.13730000000000	O	(24d)
0.20250000000000	−0.04190000000000	0.63730000000000	O	(24d)
−0.29750000000000	0.54190000000000	0.36270000000000	O	(24d)
0.13730000000000	0.29750000000000	0.04190000000000	O	(24d)
−0.13730000000000	0.79750000000000	0.45810000000000	O	(24d)
0.63730000000000	0.20250000000000	−0.04190000000000	O	(24d)
0.36270000000000	−0.29750000000000	0.54190000000000	O	(24d)
−0.04190000000000	−0.13730000000000	−0.29750000000000	O	(24d)
0.54190000000000	0.13730000000000	0.20250000000000	O	(24d)
0.04190000000000	0.36270000000000	0.79750000000000	O	(24d)
0.45810000000000	0.63730000000000	0.29750000000000	O	(24d)
−0.29750000000000	−0.04190000000000	−0.13730000000000	O	(24d)
0.20250000000000	0.54190000000000	0.13730000000000	O	(24d)
0.79750000000000	0.04190000000000	0.36270000000000	O	(24d)
0.29750000000000	0.45810000000000	0.63730000000000	O	(24d)
−0.13730000000000	−0.29750000000000	−0.04190000000000	O	(24d)
0.13730000000000	0.20250000000000	0.54190000000000	O	(24d)
0.36270000000000	0.79750000000000	0.04190000000000	O	(24d)
0.63730000000000	0.29750000000000	0.45810000000000	O	(24d)
0.30220000000000	0.30220000000000	0.30220000000000	S	(8c)
0.19780000000000	−0.30220000000000	0.80220000000000	S	(8c)
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0.80220000000000	0.30220000000000	0.19780000000000	S	(8c)
0.30220000000000	0.19780000000000	0.80220000000000	S	(8c)
0.19780000000000	0.80220000000000	0.30220000000000	S	(8c)

Maghemite (γ-Fe_2O_3, $D5_7$): A2B3_cP60_212_bcd_ace - CIF

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_chemical_name_mineral 'Maghemite or $\gamma$-corundum'
_chemical_formula_sum 'Fe2 O3'

loop_
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 'J. Thewlis'
_journal_name_full_name
;
 Philosophical Magazine
;
_journal_volume 12
_journal_year 1931
_journal_page_first 1089
_journal_page_last 1106
_publ_Section_title
;
 The structure of ferromagnetic ferric oxide
;

# Found in Strukturbericht Band II 1928-1932, 1937

_aflow_title 'Maghemite ($\gamma$-Fe$_{2}$O$_{3}$, $D5_{7}$) Structure'
_aflow_proto 'A2B3_cP60_212_bcd_ace'
_aflow_params 'a,x_{3},x_{4},y_{5},x_{6},y_{6},z_{6}'
_aflow_params_values '8.4,0.0,0.871,0.375,0.378,0.129,0.629'
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_aflow_Strukturbericht '$D5_{7}$'
_aflow_Pearson 'cP60'

_symmetry_space_group_name_H-M "P 43 3 2"
_symmetry_Int_Tables_number 212

_cell_length_a      8.40000
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_cell_length_c      8.40000
_cell_angle_alpha  90.00000
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_cell_angle_gamma  90.00000

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_space_group_symop_operation_xyz
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2 x+1/2,-y+1/2,-z
3 -x,y+1/2,-z+1/2
4 -x+1/2,-y,z+1/2
5 y,z,x
6 y+1/2,-z+1/2,-x
7 -y,z+1/2,-x+1/2
8 -y+1/2,-z,x+1/2
9 z,x,y
10 z+1/2,-x+1/2,-y
11 -z,x+1/2,-y+1/2
12 -z+1/2,-x,y+1/2
13 -y+1/4,-x+1/4,-z+1/4
14 -y+3/4,x+1/4,z+3/4
15 y+3/4,-x+3/4,z+1/4
16 y+1/4,x+3/4,-z+3/4
17 -x+1/4,-z+1/4,-y+1/4
18 -x+3/4,z+1/4,y+3/4
19 x+3/4,-z+3/4,y+1/4
20 x+1/4,z+3/4,-y+3/4
21 -z+1/4,-y+1/4,-x+1/4
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_atom_site_Wyckoff_label
_atom_site_fract_x
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_atom_site_fract_z
_atom_site_occupancy
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Fe1 Fe 4 b 0.62500 0.62500 0.62500 1.00000
Fe2 Fe 8 c 0.00000 0.00000 0.00000 1.00000
O2 O 8 c 0.87100 0.87100 0.87100 1.00000
Fe3 Fe 12 d 0.12500 0.37500 0.87500 1.00000
O3 O 24 e 0.37800 0.12900 0.62900 1.00000
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Maghemite (γ-Fe_2O_3, $D5_7$): A2B3_cP60_212_bcd_ace - POSCAR

```
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      ↪ ,0.378,0.129,0.629 & P4_{3}32 O^{6} #212 (abc^2de) & cP60 &
      ↪ $D5_{7}$ & Fe2O3 & Maghemite or $\gamma$-corundum & J. Thewlis,
      ↪  Philos. Mag. 12, 1089-1106 (1931)
   1.00000000000000
     8.40000000000000     0.00000000000000     0.00000000000000
     0.00000000000000     8.40000000000000     0.00000000000000
     0.00000000000000     0.00000000000000     8.40000000000000
      Fe    O
      24    36
Direct
     0.62500000000000     0.62500000000000     0.62500000000000  Fe   (4b)
     0.87500000000000     0.37500000000000     0.12500000000000  Fe   (4b)
     0.37500000000000     0.12500000000000     0.87500000000000  Fe   (4b)
     0.12500000000000     0.87500000000000     0.37500000000000  Fe   (4b)
     0.00000000000000     0.00000000000000     0.00000000000000  Fe   (8c)
     0.50000000000000     0.00000000000000     0.50000000000000  Fe   (8c)
     0.00000000000000     0.50000000000000     0.50000000000000  Fe   (8c)
     0.50000000000000     0.50000000000000     0.00000000000000  Fe   (8c)
     0.25000000000000     0.75000000000000     0.75000000000000  Fe   (8c)
     0.25000000000000     0.25000000000000     0.25000000000000  Fe   (8c)
     0.75000000000000     0.75000000000000     0.25000000000000  Fe   (8c)
     0.75000000000000     0.25000000000000     0.75000000000000  Fe   (8c)
     0.12500000000000     0.37500000000000    -0.12500000000000  Fe  (12d)
     0.37500000000000    -0.37500000000000     0.37500000000000  Fe  (12d)
     0.87500000000000     0.87500000000000     0.62500000000000  Fe  (12d)
     0.62500000000000     0.12500000000000     1.12500000000000  Fe  (12d)
    -0.12500000000000     0.12500000000000     0.37500000000000  Fe  (12d)
     0.37500000000000     0.37500000000000    -0.37500000000000  Fe  (12d)
     0.62500000000000     0.87500000000000     0.87500000000000  Fe  (12d)
     1.12500000000000     0.62500000000000     0.12500000000000  Fe  (12d)
     0.37500000000000    -0.12500000000000     0.12500000000000  Fe  (12d)
    -0.37500000000000     0.37500000000000     0.37500000000000  Fe  (12d)
     0.87500000000000     0.62500000000000     0.87500000000000  Fe  (12d)
     0.12500000000000     1.12500000000000     0.62500000000000  Fe  (12d)
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     0.37500000000000     0.87500000000000     0.62500000000000  O   (4a)
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     0.62500000000000     0.37500000000000     0.87500000000000  O   (4a)
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     1.12100000000000     1.62100000000000    -0.12100000000000  O   (8c)
    -0.62100000000000    -0.62100000000000    -0.62100000000000  O   (8c)
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    -0.12100000000000     1.12100000000000     1.62100000000000  O   (8c)
     0.37800000000000     0.12900000000000     0.62900000000000  O  (24e)
     0.12200000000000    -0.12900000000000     1.12900000000000  O  (24e)
    -0.37800000000000     0.62900000000000    -0.12900000000000  O  (24e)
     0.87800000000000     0.37100000000000    -0.62900000000000  O  (24e)
     0.62900000000000     0.37800000000000     0.12900000000000  O  (24e)
     1.12900000000000     0.12200000000000    -0.12900000000000  O  (24e)
    -0.12900000000000    -0.37800000000000     0.62900000000000  O  (24e)
    -0.62900000000000     0.87800000000000     0.37100000000000  O  (24e)
     0.12900000000000     0.62900000000000     0.37800000000000  O  (24e)
    -0.12900000000000     1.12900000000000     0.12200000000000  O  (24e)
     0.62900000000000    -0.12900000000000    -0.37800000000000  O  (24e)
     0.37100000000000    -0.62900000000000     0.87800000000000  O  (24e)
     0.37900000000000     1.12800000000000     0.12100000000000  O  (24e)
     0.12100000000000    -0.12800000000000    -0.37900000000000  O  (24e)
     0.87900000000000     0.37200000000000     0.87900000000000  O  (24e)
     0.62100000000000     0.62800000000000     1.37900000000000  O  (24e)
     0.62800000000000     1.37900000000000     0.62100000000000  O  (24e)
     0.37200000000000     0.87900000000000     0.87900000000000  O  (24e)
    -0.12800000000000    -0.37900000000000     0.12100000000000  O  (24e)
     1.12800000000000     0.12100000000000     0.37900000000000  O  (24e)
     0.87900000000000     0.87900000000000     0.37200000000000  O  (24e)
     1.37900000000000     0.62100000000000     0.62800000000000  O  (24e)
     0.12100000000000     0.37900000000000     1.12800000000000  O  (24e)
    -0.37900000000000     0.12100000000000    -0.12800000000000  O  (24e)
```

Al_2Mo_3C: A2BC3_cP24_213_c_a_d - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Al2CMo3'
_chemical_formula_sum 'Al2 C Mo3'

loop_
_publ_author_name
 'W. Jeitschko'
 'H. Nowotny'
 'F. Benesovsky'
_journal_name_full_name
;
 Monatshefte f{\"u}r Chemie und verwandte Teile anderer Wissenschaften
;
_journal_volume 94
_journal_year 1963
_journal_page_first 247
_journal_page_last 251
_publ_Section_title
;
 Ein Beitrag zum Dreistoff: Molybd\"{a}n-Aluminium-Kohlenstoff
;

# Found in A. Iyo et al., Superconductivity induced by Mg deficiency in
      ↪ noncentrosymmetric phosphide Mg${2}$Rh${3}$P, Phys Rev.
      ↪ Mater. 3, 124802 (2019)

_aflow_title 'Al${2}$Mo${3}$C Structure'
_aflow_proto 'A2BC3_cP24_213_c_a_d'
_aflow_params 'a,x_{2},y_{3}'
_aflow_params_values '6.84,0.061,0.206'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'cP24'

_symmetry_space_group_name_H-M "P 41 3 2"
_symmetry_Int_Tables_number 213

_cell_length_a      6.84000
_cell_length_b      6.84000
_cell_length_c      6.84000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z
3 -x,y+1/2,-z+1/2
4 -x+1/2,-y,z+1/2
5 y,z,x
6 y+1/2,-z+1/2,-x
7 -y,z+1/2,-x+1/2
8 -y+1/2,-z,x+1/2
9 z,x,y
10 z+1/2,-x+1/2,-y
11 -z,x+1/2,-y+1/2
12 -z+1/2,-x,y+1/2
13 -y+3/4,-x+3/4,-z+3/4
14 -y+1/4,x+3/4,z+1/4
15 y+1/4,-x+1/4,z+3/4
16 y+3/4,x+1/4,-z+1/4
17 -x+3/4,-z+3/4,-y+3/4
18 -x+1/4,z+3/4,y+1/4
19 x+1/4,-z+1/4,y+3/4
20 x+3/4,z+1/4,-y+1/4
21 -z+3/4,-y+3/4,-x+3/4
22 -z+1/4,y+3/4,x+1/4
23 z+1/4,-y+1/4,x+3/4
24 z+3/4,y+1/4,-x+1/4

loop_
_atom_site_label
_atom_site_type_symbol
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_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cl1 C 4 a 0.37500 0.37500 0.37500 1.00000
Al1 Al 8 c 0.06100 0.06100 0.06100 1.00000
Mo1 Mo 12 d 0.12500 0.20600 0.45600 1.00000
```

Al_2Mo_3C: A2BC3_cP24_213_c_a_d - POSCAR

```
A2BC3_cP24_213_c_a_d & a,x2,y3 --params=6.84,0.061,0.206 & P4_{1}32 O^{7
    ↪ } #213 (acd) & cP24 & None & Al2CMo3 & Al2CMo3 & W. Jeitschko
    ↪ and H. Nowotny and F. Benesovsky, Monatsh. Chem. Verw. Teile
    ↪ Anderer Wiss. 94, 247-251 (1963)
   1.00000000000000
   6.84000000000000   0.00000000000000   0.00000000000000
   0.00000000000000   6.84000000000000   0.00000000000000
   0.00000000000000   0.00000000000000   6.84000000000000
    Al    C    Mo
     8    4    12
Direct
   0.06100000000000   0.06100000000000   0.06100000000000   Al   (8c)
   0.43900000000000  -0.06100000000000   0.56100000000000   Al   (8c)
  -0.06100000000000   0.56100000000000   0.43900000000000   Al   (8c)
   0.56100000000000   0.43900000000000  -0.06100000000000   Al   (8c)
   0.81100000000000   0.31100000000000   0.18900000000000   Al   (8c)
   0.68900000000000   0.68900000000000   0.68900000000000   Al   (8c)
   0.31100000000000   0.18900000000000   0.81100000000000   Al   (8c)
   0.18900000000000   0.81100000000000   0.31100000000000   Al   (8c)
   0.37500000000000   0.37500000000000   0.37500000000000   C    (4a)
   0.12500000000000   0.62500000000000   0.87500000000000   C    (4a)
   0.62500000000000   0.87500000000000   0.12500000000000   C    (4a)
   0.87500000000000   0.12500000000000   0.62500000000000   C    (4a)
   0.12500000000000   0.20600000000000   0.45600000000000   Mo   (12d)
   0.37500000000000  -0.20600000000000   0.95600000000000   Mo   (12d)
   0.87500000000000   0.70600000000000   0.04400000000000   Mo   (12d)
   0.62500000000000   0.29400000000000   0.54400000000000   Mo   (12d)
   0.45600000000000   0.12500000000000   0.20600000000000   Mo   (12d)
   0.95600000000000   0.37500000000000  -0.20600000000000   Mo   (12d)
   0.04400000000000   0.87500000000000   0.70600000000000   Mo   (12d)
   0.54400000000000   0.62500000000000   0.29400000000000   Mo   (12d)
   0.20600000000000   0.45600000000000   0.12500000000000   Mo   (12d)
  -0.20600000000000   0.95600000000000   0.37500000000000   Mo   (12d)
   0.70600000000000   0.04400000000000   0.87500000000000   Mo   (12d)
   0.29400000000000   0.54400000000000   0.62500000000000   Mo   (12d)
```

Mg_3Ru_2: A3B2_cP20_213_d_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Mg3Ru2'
_chemical_formula_sum 'Mg3 Ru2'

loop_
_publ_author_name
 'R. P\"{o}ttgen'
 'V. Hlukhyy'
 'A. Baranov'
 'Y. Grin'
_journal_name_full_name
;
 Inorganic Chemistry
;
_journal_volume 47
_journal_year 2008
_journal_page_first 6051
_journal_page_last 6055
_publ_Section_title
;
 Crystal Structure and Chemical Bonding of Mg$_{3}$Ru$_{2}$
;

_aflow_title 'Mg$_{3}$Ru$_{2}$ Structure'
_aflow_proto 'A3B2_cP20_213_d_c'
_aflow_params 'a,x_{1},y_{2}'
_aflow_params_values '6.9352,0.07378,0.2051'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'cP20'

_symmetry_space_group_name_H-M "P 41 3 2"
_symmetry_Int_Tables_number 213

_cell_length_a    6.93520
_cell_length_b    6.93520
_cell_length_c    6.93520
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x+1/2,-y+1/2,-z
3 -x,y+1/2,-z+1/2
4 -x+1/2,-y,z+1/2
5 y,z,x
6 y+1/2,-z+1/2,-x
7 -y,z+1/2,-x+1/2
8 -y+1/2,-z,x+1/2
9 z,x,y
10 z+1/2,-x+1/2,-y
11 -z,x+1/2,-y+1/2
12 -z+1/2,-x,y+1/2
13 -y+3/4,-x+3/4,-z+3/4
14 -y+1/4,x+3/4,z+1/4
15 y+1/4,-x+1/4,z+3/4
16 y+3/4,x+1/4,-z+1/4
17 -x+3/4,-z+3/4,-y+3/4
18 -x+1/4,z+3/4,y+1/4
19 x+1/4,-z+1/4,y+3/4
20 x+3/4,z+1/4,-y+1/4
21 -z+3/4,-y+3/4,-x+3/4
22 -z+1/4,y+3/4,x+1/4
23 z+1/4,-y+1/4,x+3/4
24 z+3/4,y+1/4,-x+1/4

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ru1 Ru 8 c 0.07378 0.07378 0.07378 1.00000
Mg1 Mg 12 d 0.12500 0.20510 0.45510 1.00000
```

Mg_3Ru_2: A3B2_cP20_213_d_c - POSCAR

```
A3B2_cP20_213_d_c & a,x1,y2 --params=6.9352,0.07378,0.2051 & P4_{1}32 O
    ↪ ^{7} #213 (cd) & cP20 & None & Mg3Ru2 & Mg3Ru2 & R. P\"{o}ttgen
    ↪ et al., Inorg. Chem. 47, 6051-6055 (2008)
   1.00000000000000
   6.93520000000000   0.00000000000000   0.00000000000000
   0.00000000000000   6.93520000000000   0.00000000000000
   0.00000000000000   0.00000000000000   6.93520000000000
    Mg   Ru
    12    8
Direct
   0.12500000000000   0.20510000000000   0.45510000000000   Mg   (12d)
   0.37500000000000  -0.20510000000000   0.95510000000000   Mg   (12d)
   0.87500000000000   0.70510000000000   0.04490000000000   Mg   (12d)
   0.62500000000000   0.29490000000000   0.54490000000000   Mg   (12d)
   0.45510000000000   0.12500000000000   0.20510000000000   Mg   (12d)
   0.95510000000000   0.37500000000000  -0.20510000000000   Mg   (12d)
   0.04490000000000   0.87500000000000   0.70510000000000   Mg   (12d)
   0.54490000000000   0.62500000000000   0.29490000000000   Mg   (12d)
   0.20510000000000   0.45510000000000   0.12500000000000   Mg   (12d)
  -0.20510000000000   0.95510000000000   0.37500000000000   Mg   (12d)
   0.70510000000000   0.04490000000000   0.87500000000000   Mg   (12d)
   0.29490000000000   0.54490000000000   0.62500000000000   Mg   (12d)
   0.07378000000000   0.07378000000000   0.07378000000000   Ru   (8c)
   0.42622000000000  -0.07378000000000   0.57378000000000   Ru   (8c)
  -0.07378000000000   0.57378000000000   0.42622000000000   Ru   (8c)
   0.57378000000000   0.42622000000000  -0.07378000000000   Ru   (8c)
   0.82378000000000   0.32378000000000   0.17622000000000   Ru   (8c)
   0.67622000000000   0.67622000000000   0.67622000000000   Ru   (8c)
   0.32378000000000   0.17622000000000   0.82378000000000   Ru   (8c)
   0.17622000000000   0.82378000000000   0.32378000000000   Ru   (8c)
```

Zunyite $[Al_{13}(OH,F)_{18}Si_5O_{20}Cl, S0_8]$: A13BC18D20E5_cF228_216_dh_b_fh_2eh_ce - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Zunyite'
_chemical_formula_sum 'Al13 Cl F18 O20 Si5'

loop_
_publ_author_name
 'W. B. Kamb'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 13
_journal_year 1960
_journal_page_first 15
_journal_page_last 24
_publ_Section_title
;
 The Crystal Structure of Zunyite
;

_aflow_title 'Zunyite [Al$_{13}$(OH,F)$_{18}$Si$_{5}$O$_{20}$Cl, $S0_{8}
    ↪ $] Structure'
_aflow_proto 'A13BC18D20E5_cF228_216_dh_b_fh_2eh_ce'
_aflow_params 'a,x_{4},x_{5},x_{6},x_{7},x_{8},z_{8},x_{9},z_{9},x_{10},
    ↪ z_{10}'
_aflow_params_values '13.87,0.825,0.1818,0.1143,0.278,0.0853,0.7667,
    ↪ 0.1793,0.5466,0.1385,0.0003'
_aflow_Strukturbericht '$S0_{8}$'
_aflow_Pearson 'cF228'

_symmetry_space_group_name_H-M "F -4 3 m"
_symmetry_Int_Tables_number 216

_cell_length_a    13.87000
_cell_length_b    13.87000
_cell_length_c    13.87000
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000
```

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loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 y,z,x
6 y,-z,-x
7 -y,z,-x
8 -y,-z,x
9 z,x,y
10 z,-x,-y
11 -z,x,-y
12 -z,-x,y
13 y,x,z
14 y,-x,-z
15 -y,x,-z
16 -y,-x,z
17 x,z,y
18 x,-z,-y
19 -x,z,-y
20 -x,-z,y
21 z,y,x
22 z,-y,-x
23 -z,y,-x
24 -z,-y,x
25 x,y+1/2,z+1/2
26 x,-y+1/2,-z+1/2
27 -x,y+1/2,-z+1/2
28 -x,-y+1/2,z+1/2
29 y,z+1/2,x+1/2
30 y,-z+1/2,-x+1/2
31 -y,z+1/2,-x+1/2
32 -y,-z+1/2,x+1/2
33 z,x+1/2,y+1/2
34 z,-x+1/2,-y+1/2
35 -z,x+1/2,-y+1/2
36 -z,-x+1/2,y+1/2
37 y,x+1/2,z+1/2
38 y,-x+1/2,-z+1/2
39 -y,x+1/2,-z+1/2
40 -y,-x+1/2,z+1/2
41 x,z+1/2,y+1/2
42 x,-z+1/2,-y+1/2
43 -x,z+1/2,-y+1/2
44 -x,-z+1/2,y+1/2
45 z,y+1/2,x+1/2
46 z,-y+1/2,-x+1/2
47 -z,y+1/2,-x+1/2
48 -z,-y+1/2,x+1/2
49 x+1/2,y,z+1/2
50 x+1/2,-y,-z+1/2
51 -x+1/2,y,-z+1/2
52 -x+1/2,-y,z+1/2
53 y+1/2,z,x+1/2
54 y+1/2,-z,-x+1/2
55 -y+1/2,z,-x+1/2
56 -y+1/2,-z,x+1/2
57 z+1/2,x,y+1/2
58 z+1/2,-x,-y+1/2
59 -z+1/2,x,-y+1/2
60 -z+1/2,-x,y+1/2
61 y+1/2,x,z+1/2
62 y+1/2,-x,-z+1/2
63 -y+1/2,x,-z+1/2
64 -y+1/2,-x,z+1/2
65 x+1/2,z,y+1/2
66 x+1/2,-z,-y+1/2
67 -x+1/2,z,-y+1/2
68 -x+1/2,-z,y+1/2
69 z+1/2,y,x+1/2
70 z+1/2,-y,-x+1/2
71 -z+1/2,y,-x+1/2
72 -z+1/2,-y,x+1/2
73 x+1/2,y+1/2,z
74 x+1/2,-y+1/2,-z
75 -x+1/2,y+1/2,-z
76 -x+1/2,-y+1/2,z
77 y+1/2,z+1/2,x
78 y+1/2,-z+1/2,-x
79 -y+1/2,z+1/2,-x
80 -y+1/2,-z+1/2,x
81 z+1/2,x+1/2,y
82 z+1/2,-x+1/2,-y
83 -z+1/2,x+1/2,-y
84 -z+1/2,-x+1/2,y
85 y+1/2,x+1/2,z
86 y+1/2,-x+1/2,-z
87 -y+1/2,x+1/2,-z
88 -y+1/2,-x+1/2,z
89 x+1/2,z+1/2,y
90 x+1/2,-z+1/2,-y
91 -x+1/2,z+1/2,-y
92 -x+1/2,-z+1/2,y
93 z+1/2,y+1/2,x
94 z+1/2,-y+1/2,-x
95 -z+1/2,y+1/2,-x
96 -z+1/2,-y+1/2,x

loop_
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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cl1 Cl 4 b 0.50000 0.50000 0.50000 1.00000
Si1 Si 4 c 0.25000 0.25000 0.25000 1.00000
Al1 Al 4 d 0.75000 0.75000 0.75000 1.00000
O1 O 16 e 0.82500 0.82500 0.82500 1.00000
O2 O 16 e 0.18180 0.18180 0.18180 1.00000
Si2 Si 16 e 0.11430 0.11430 0.11430 1.00000
F1 F 24 f 0.27800 0.00000 0.00000 1.00000
Al2 Al 48 h 0.08530 0.08530 0.76670 1.00000
F2 F 48 h 0.17930 0.17930 0.54660 1.00000
O3 O 48 h 0.13850 0.13850 0.00030 1.00000
```

Zunyite [$Al_{13}(OH,F)_{18}Si_5O_{20}Cl$, $S0_8$]: A13BC18D20E5_cF228_216_dh_b_fh_2eh_ce - POSCAR

```
A13BC18D20E5_cF228_216_dh_b_fh_2eh_ce & a,x4,x5,x6,x7,x8,z8,x9,z9,x10,
    ↪ z10 --params=13.87,0.825,0.1818,0.1143,0.278,0.0853,0.7667,
    ↪ 0.1793,0.5466,0.1385,0.0003 & F-43m T_{d}^{2} #216 (bcde^3fh^3)
    ↪ & cF228 & $S0_{8}$ & A13ClF18O20Si5 & Zunyite & W. B. Kamb,
    ↪ Acta Cryst. 13, 15-24 (1960)
   1.00000000000000
   0.00000000000000   6.93500000000000   6.93500000000000
   6.93500000000000   0.00000000000000   6.93500000000000
   6.93500000000000   6.93500000000000   0.00000000000000
    Al   Cl   F   O   Si
    13    1   18   20    5
Direct
   0.75000000000000   0.75000000000000   0.75000000000000   Al   (4d)
   0.76670000000000   0.76670000000000  -0.59610000000000   Al   (48h)
   0.76670000000000   0.76670000000000  -0.93730000000000   Al   (48h)
  -0.59610000000000  -0.93730000000000   0.76670000000000   Al   (48h)
  -0.93730000000000  -0.59610000000000   0.76670000000000   Al   (48h)
  -0.59610000000000   0.76670000000000   0.76670000000000   Al   (48h)
  -0.93730000000000   0.76670000000000   0.76670000000000   Al   (48h)
   0.76670000000000  -0.59610000000000  -0.93730000000000   Al   (48h)
   0.76670000000000  -0.93730000000000  -0.59610000000000   Al   (48h)
   0.76670000000000  -0.59610000000000   0.76670000000000   Al   (48h)
   0.76670000000000  -0.93730000000000   0.76670000000000   Al   (48h)
  -0.93730000000000   0.76670000000000  -0.59610000000000   Al   (48h)
  -0.59610000000000   0.76670000000000  -0.93730000000000   Al   (48h)
   0.50000000000000   0.50000000000000   0.50000000000000   Cl   (4b)
  -0.27800000000000   0.27800000000000   0.27800000000000   F   (24f)
   0.27800000000000  -0.27800000000000  -0.27800000000000   F   (24f)
   0.27800000000000  -0.27800000000000   0.27800000000000   F   (24f)
  -0.27800000000000   0.27800000000000  -0.27800000000000   F   (24f)
   0.27800000000000   0.27800000000000  -0.27800000000000   F   (24f)
  -0.27800000000000  -0.27800000000000   0.27800000000000   F   (24f)
   0.54660000000000   0.54660000000000  -0.18800000000000   F   (48h)
   0.54660000000000   0.54660000000000  -0.90520000000000   F   (48h)
  -0.18800000000000  -0.90520000000000   0.54660000000000   F   (48h)
  -0.90520000000000  -0.18800000000000   0.54660000000000   F   (48h)
  -0.18800000000000   0.54660000000000   0.54660000000000   F   (48h)
  -0.90520000000000   0.54660000000000   0.54660000000000   F   (48h)
   0.54660000000000  -0.18800000000000  -0.90520000000000   F   (48h)
   0.54660000000000  -0.90520000000000  -0.18800000000000   F   (48h)
   0.54660000000000  -0.18800000000000   0.54660000000000   F   (48h)
   0.54660000000000  -0.90520000000000   0.54660000000000   F   (48h)
  -0.90520000000000   0.54660000000000  -0.18800000000000   F   (48h)
  -0.18800000000000   0.54660000000000  -0.90520000000000   F   (48h)
   0.82500000000000   0.82500000000000   0.82500000000000   O   (16e)
   0.82500000000000   0.82500000000000  -2.47500000000000   O   (16e)
   0.82500000000000  -2.47500000000000   0.82500000000000   O   (16e)
  -2.47500000000000   0.82500000000000   0.82500000000000   O   (16e)
   0.18180000000000   0.18180000000000   0.18180000000000   O   (16e)
   0.18180000000000   0.18180000000000  -0.54540000000000   O   (16e)
   0.18180000000000  -0.54540000000000   0.18180000000000   O   (16e)
  -0.54540000000000   0.18180000000000   0.18180000000000   O   (16e)
   0.00030000000000   0.00030000000000   0.27670000000000   O   (48h)
   0.00030000000000   0.00030000000000  -0.27730000000000   O   (48h)
   0.27670000000000  -0.27730000000000   0.00030000000000   O   (48h)
  -0.27730000000000   0.27670000000000   0.00030000000000   O   (48h)
   0.27670000000000   0.00030000000000   0.00030000000000   O   (48h)
  -0.27730000000000   0.00030000000000   0.00030000000000   O   (48h)
   0.00030000000000   0.27670000000000  -0.27730000000000   O   (48h)
   0.00030000000000  -0.27730000000000   0.27670000000000   O   (48h)
   0.00030000000000   0.27670000000000   0.00030000000000   O   (48h)
   0.00030000000000  -0.27730000000000   0.00030000000000   O   (48h)
  -0.27730000000000   0.00030000000000   0.27670000000000   O   (48h)
   0.27670000000000   0.00030000000000  -0.27730000000000   O   (48h)
   0.25000000000000   0.25000000000000   0.25000000000000   Si   (4c)
   0.11430000000000   0.11430000000000   0.11430000000000   Si   (16e)
   0.11430000000000   0.11430000000000  -0.34290000000000   Si   (16e)
   0.11430000000000  -0.34290000000000   0.11430000000000   Si   (16e)
  -0.34290000000000   0.11430000000000   0.11430000000000   Si   (16e)
```

Muratatite [$(Y,Na)_6(Zn,Fe)_5Ti_{12}O_{29}(O,F)_{10}F_4$]: A16B40C12D6E5_cF316_216_eh_e2g2h_h_f_be - CIF

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_audit_creation_method FINDSYM

_chemical_name_mineral 'Murataite'
_chemical_formula_sum 'F16 O40 Ti12 Y6 Zn5'

loop_
_publ_author_name
 'T. S. Ercit'
 'F. C. Hawthorne'
_journal_name_full_name
;
 Canadian Mineralogist
;
```

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_journal_volume 33
_journal_year 1995
_journal_page_first 1233
_journal_page_last 1229
_publ_Section_title
;
 Muraitaite, A UB$_{12}$ derivative structure with condensed Keggin
      ↪ molecules
;

_aflow_title 'Muraitaite [(Y,Na)$_{6}$(Zn,Fe)$_{5}$Ti$_{12}$O$_{29}$(O,F)
      ↪ $_{10}$F$_{4}$] Structure'
_aflow_proto 'A16B40C12D6E5_cF316_216_eh_e2g2h_h_f_be'
_aflow_params 'a,x_{2},x_{3},x_{4},x_{5},x_{6},x_{7},x_{8},z_{8},x_{9},
      ↪ z_{9},x_{10},z_{10},x_{11},z_{11}'
_aflow_params_values '14.886,0.079,0.4262,0.1722,0.1812,0.267,0.515,
      ↪ 0.894,-0.085,-0.0729,0.7024,0.6038,0.258,0.8371,0.4913'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'cF316'

_symmetry_space_group_name_H-M "F -4 3 m"
_symmetry_Int_Tables_number 216

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_cell_length_b      14.88600
_cell_length_c      14.88600
_cell_angle_alpha   90.00000
_cell_angle_beta    90.00000
_cell_angle_gamma   90.00000

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_space_group_symop_operation_xyz
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2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 y,z,x
6 y,-z,-x
7 -y,z,-x
8 -y,-z,x
9 z,x,y
10 z,-x,-y
11 -z,x,-y
12 -z,-x,y
13 y,x,z
14 y,-x,-z
15 -y,x,-z
16 -y,-x,z
17 x,z,y
18 x,-z,-y
19 -x,z,-y
20 -x,-z,y
21 z,y,x
22 z,-y,-x
23 -z,y,-x
24 -z,-y,x
25 x,y+1/2,z+1/2
26 x,-y+1/2,-z+1/2
27 -x,y+1/2,-z+1/2
28 -x,-y+1/2,z+1/2
29 y,z+1/2,x+1/2
30 y,-z+1/2,-x+1/2
31 -y,z+1/2,-x+1/2
32 -y,-z+1/2,x+1/2
33 z,x+1/2,y+1/2
34 z,-x+1/2,-y+1/2
35 -z,x+1/2,-y+1/2
36 -z,-x+1/2,y+1/2
37 y,x+1/2,z+1/2
38 y,-x+1/2,-z+1/2
39 -y,x+1/2,-z+1/2
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42 x,-z+1/2,-y+1/2
43 -x,z+1/2,-y+1/2
44 -x,-z+1/2,y+1/2
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46 z,-y+1/2,-x+1/2
47 -z,y+1/2,-x+1/2
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50 x+1/2,-y,-z+1/2
51 -x+1/2,y,-z+1/2
52 -x+1/2,-y,z+1/2
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54 y+1/2,-z,-x+1/2
55 -y+1/2,z,-x+1/2
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63 -y+1/2,x,-z+1/2
64 -y+1/2,-x,z+1/2
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66 x+1/2,-z,-y+1/2
67 -x+1/2,z,-y+1/2
68 -x+1/2,-z,y+1/2
69 z+1/2,y,x+1/2
70 z+1/2,-y,-x+1/2
71 -z+1/2,y,-x+1/2
72 -z+1/2,-y,x+1/2
73 x+1/2,y+1/2,z
74 x+1/2,-y+1/2,-z
75 -x+1/2,y+1/2,-z
76 -x+1/2,-y+1/2,z
77 y+1/2,z+1/2,x
78 y+1/2,-z+1/2,-x
79 -y+1/2,z+1/2,-x
80 -y+1/2,-z+1/2,x
81 z+1/2,x+1/2,y
82 z+1/2,-x+1/2,-y
83 -z+1/2,x+1/2,-y
84 -z+1/2,-x+1/2,y
85 y+1/2,x+1/2,z
86 y+1/2,-x+1/2,-z
87 -y+1/2,x+1/2,-z
88 -y+1/2,-x+1/2,z
89 x+1/2,z+1/2,y
90 x+1/2,-z+1/2,-y
91 -x+1/2,z+1/2,-y
92 -x+1/2,-z+1/2,y
93 z+1/2,y+1/2,x
94 z+1/2,-y+1/2,-x
95 -z+1/2,y+1/2,-x
96 -z+1/2,-y+1/2,x

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_atom_site_fract_x
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_atom_site_occupancy
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F1 F 16 e 0.07900 0.07900 0.07900 1.00000
O1 O 16 e 0.42620 0.42620 0.42620 1.00000
Zn2 Zn 16 e 0.17220 0.17220 0.17220 1.00000
Y1 Y 24 f 0.18120 0.00000 0.00000 1.00000
O2 O 24 g 0.26700 0.25000 0.25000 1.00000
O3 O 24 g 0.51500 0.25000 0.25000 1.00000
F2 F 48 h 0.89400 0.89400 -0.08500 1.00000
O4 O 48 h -0.07290 -0.07290 0.70240 1.00000
O5 O 48 h 0.60380 0.60380 0.25800 1.00000
Ti1 Ti 48 h 0.83710 0.83710 0.49130 1.00000
```

Muraitaite [(Y,Na)$_6$(Zn,Fe)$_5$Ti$_{12}$O$_{29}$(O,F)$_{10}$F$_4$]: A16B40C12D6E5_cF316_216_eh_e2g2h_h_f_be - POSCAR

```
A16B40C12D6E5_cF316_216_eh_e2g2h_h_f_be & a,x2,x3,x4,x5,x6,x7,x8,z8,x9,
      ↪ z9,x10,z10,x11,z11 --params=14.886,0.079,0.4262,0.1722,0.1812,
      ↪ 0.267,0.515,0.894,-0.085,-0.0729,0.7024,0.6038,0.258,0.8371,
      ↪ 0.4913 & F-43m T_{d}^{2} #216 (be^3fg^2h^4) & cF316 & None &
      ↪ F16O40Ti12Y6Zn5 & Muraitaite & T. S. Ercit and F. C. Hawthorne,
      ↪ Can. Mineral. 33, 1233-1229 (1995)
   1.00000000000000
     0.00000000000000     7.44300000000000     7.44300000000000
     7.44300000000000     0.00000000000000     7.44300000000000
     7.44300000000000     7.44300000000000     0.00000000000000
       F     O     Ti     Y     Zn
      16    40    12     6     5
Direct
     0.07900000000000     0.07900000000000     0.07900000000000   F   (16e)
     0.07900000000000     0.07900000000000    -0.23700000000000   F   (16e)
     0.07900000000000    -0.23700000000000     0.07900000000000   F   (16e)
    -0.23700000000000     0.07900000000000     0.07900000000000   F   (16e)
    -0.08500000000000    -0.08500000000000     1.87300000000000   F   (48h)
    -0.08500000000000    -0.08500000000000    -1.70300000000000   F   (48h)
     1.87300000000000    -1.70300000000000    -0.08500000000000   F   (48h)
    -1.70300000000000     1.87300000000000    -0.08500000000000   F   (48h)
     1.87300000000000    -0.08500000000000    -0.08500000000000   F   (48h)
    -1.70300000000000    -0.08500000000000    -0.08500000000000   F   (48h)
    -0.08500000000000     1.87300000000000    -1.70300000000000   F   (48h)
    -0.08500000000000    -1.70300000000000     1.87300000000000   F   (48h)
    -0.08500000000000     1.87300000000000    -0.08500000000000   F   (48h)
    -0.08500000000000    -1.70300000000000    -0.08500000000000   F   (48h)
    -1.70300000000000    -0.08500000000000     1.87300000000000   F   (48h)
     1.87300000000000    -0.08500000000000    -1.70300000000000   F   (48h)
     0.42620000000000     0.42620000000000     0.42620000000000   O   (16e)
     0.42620000000000     0.42620000000000    -1.27860000000000   O   (16e)
     0.42620000000000    -1.27860000000000     0.42620000000000   O   (16e)
    -1.27860000000000     0.42620000000000     0.42620000000000   O   (16e)
     0.23300000000000     0.26700000000000     0.26700000000000   O   (24g)
     0.26700000000000     0.23300000000000     0.23300000000000   O   (24g)
     0.26700000000000     0.23300000000000     0.26700000000000   O   (24g)
     0.23300000000000     0.26700000000000     0.23300000000000   O   (24g)
     0.26700000000000     0.26700000000000     0.23300000000000   O   (24g)
     0.23300000000000     0.23300000000000     0.26700000000000   O   (24g)
    -0.01500000000000     0.51500000000000     0.51500000000000   O   (24g)
     0.51500000000000    -0.01500000000000    -0.01500000000000   O   (24g)
     0.51500000000000    -0.01500000000000     0.51500000000000   O   (24g)
    -0.01500000000000     0.51500000000000    -0.01500000000000   O   (24g)
     0.51500000000000     0.51500000000000    -0.01500000000000   O   (24g)
    -0.01500000000000    -0.01500000000000     0.51500000000000   O   (24g)
     0.70240000000000     0.70240000000000    -0.84820000000000   O   (48h)
     0.70240000000000     0.70240000000000    -0.55660000000000   O   (48h)
    -0.84820000000000    -0.55660000000000     0.70240000000000   O   (48h)
    -0.55660000000000    -0.84820000000000     0.70240000000000   O   (48h)
    -0.84820000000000     0.70240000000000     0.70240000000000   O   (48h)
    -0.55660000000000     0.70240000000000     0.70240000000000   O   (48h)
     0.70240000000000    -0.84820000000000    -0.55660000000000   O   (48h)
     0.70240000000000    -0.55660000000000    -0.84820000000000   O   (48h)
     0.70240000000000    -0.84820000000000     0.70240000000000   O   (48h)
     0.70240000000000    -0.55660000000000     0.70240000000000   O   (48h)
    -0.55660000000000     0.70240000000000    -0.84820000000000   O   (48h)
```

−0.84820000000000	0.70240000000000	−0.55660000000000	O	(48h)
0.25800000000000	0.25800000000000	0.94960000000000	O	(48h)
0.25800000000000	0.25800000000000	−1.46560000000000	O	(48h)
0.94960000000000	−1.46560000000000	0.25800000000000	O	(48h)
−1.46560000000000	0.94960000000000	0.25800000000000	O	(48h)
0.94960000000000	0.25800000000000	0.25800000000000	O	(48h)
−1.46560000000000	0.25800000000000	0.25800000000000	O	(48h)
0.25800000000000	0.94960000000000	−1.46560000000000	O	(48h)
0.25800000000000	−1.46560000000000	0.94960000000000	O	(48h)
0.25800000000000	0.94960000000000	0.25800000000000	O	(48h)
0.25800000000000	−1.46560000000000	0.25800000000000	O	(48h)
−1.46560000000000	0.25800000000000	0.94960000000000	O	(48h)
0.94960000000000	0.25800000000000	−1.46560000000000	O	(48h)
0.49130000000000	0.49130000000000	1.18290000000000	Ti	(48h)
0.49130000000000	0.49130000000000	−2.16550000000000	Ti	(48h)
1.18290000000000	−2.16550000000000	0.49130000000000	Ti	(48h)
−2.16550000000000	1.18290000000000	0.49130000000000	Ti	(48h)
1.18290000000000	0.49130000000000	0.49130000000000	Ti	(48h)
−2.16550000000000	0.49130000000000	0.49130000000000	Ti	(48h)
0.49130000000000	1.18290000000000	−2.16550000000000	Ti	(48h)
0.49130000000000	−2.16550000000000	1.18290000000000	Ti	(48h)
0.49130000000000	1.18290000000000	0.49130000000000	Ti	(48h)
0.49130000000000	−2.16550000000000	0.49130000000000	Ti	(48h)
−2.16550000000000	0.49130000000000	1.18290000000000	Ti	(48h)
1.18290000000000	0.49130000000000	−2.16550000000000	Ti	(48h)
−0.18120000000000	0.18120000000000	0.18120000000000	Y	(24f)
0.18120000000000	−0.18120000000000	−0.18120000000000	Y	(24f)
0.18120000000000	−0.18120000000000	0.18120000000000	Y	(24f)
−0.18120000000000	0.18120000000000	−0.18120000000000	Y	(24f)
0.18120000000000	0.18120000000000	−0.18120000000000	Y	(24f)
−0.18120000000000	−0.18120000000000	0.18120000000000	Y	(24f)
0.50000000000000	0.50000000000000	0.50000000000000	Zn	(4b)
0.17220000000000	0.17220000000000	0.17220000000000	Zn	(16e)
0.17220000000000	0.17220000000000	−0.51660000000000	Zn	(16e)
0.17220000000000	−0.51660000000000	0.17220000000000	Zn	(16e)
−0.51660000000000	0.17220000000000	0.17220000000000	Zn	(16e)

$Sm_{11}Cd_{45}$: A45B11_cF448_216_bd4efg5h_ac2eh - CIF

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_audit_creation_method FINDSYM

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 'M. L. Fornasini'
 'B. Chabot'
 'E. Parth\\'{e}'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 34
_journal_year 1978
_journal_page_first 2093
_journal_page_last 2099
_publ_Section_title
;
 The crystal structure of Sm$_{11}$Cd$_{45}$ with $\gamma$-brass and $\
      ↪ alpha$-Mn clusters
;

_aflow_title 'Sm$_{11}$Cd$_{45}$ Structure'
_aflow_proto 'A45B11_cF448_216_bd4efg5h_ac2eh'
_aflow_params 'a,x_{5},x_{6},x_{7},x_{8},x_{9},x_{10},x_{11},x_{12},x_{
      ↪ 13},z_{13},x_{14},z_{14},x_{15},z_{15},x_{16},z_{16},x_{17},z_{
      ↪ 17},x_{18},z_{18}'
_aflow_params_values '21.699,0.0834,0.9126,0.1636,0.8297,0.4059,0.6618,
      ↪ 0.1573,0.0895,0.2958,0.3904,0.4377,0.2627,0.5455,0.6403,0.6728,
      ↪ 0.5123,0.9161,0.7637,0.1735,0.0142'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'cF448'

_symmetry_space_group_name_H-M "F -4 3 m"
_symmetry_Int_Tables_number 216

_cell_length_a    21.69900
_cell_length_b    21.69900
_cell_length_c    21.69900
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

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4 -x,-y,z
5 y,z,x
6 y,-z,-x
7 -y,z,-x
8 -y,-z,x
9 z,x,y
10 z,-x,-y
11 -z,x,-y
12 -z,-x,y
13 y,x,z
14 y,-x,-z
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18 x,-z,-y
19 -x,z,-y
20 -x,-z,y
21 z,y,x
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35 -z,x+1/2,-y+1/2
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71 -z+1/2,y,-x+1/2
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78 y+1/2,-z+1/2,-x
79 -y+1/2,z+1/2,-x
80 -y+1/2,-z+1/2,x
81 z+1/2,x+1/2,y
82 z+1/2,-x+1/2,-y
83 -z+1/2,x+1/2,-y
84 -z+1/2,-x+1/2,y
85 y+1/2,x+1/2,z
86 y+1/2,-x+1/2,-z
87 -y+1/2,x+1/2,-z
88 -y+1/2,-x+1/2,z
89 x+1/2,z+1/2,y
90 x+1/2,-z+1/2,-y
91 -x+1/2,z+1/2,-y
92 -x+1/2,-z+1/2,y
93 z+1/2,y+1/2,x
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96 -z+1/2,-y+1/2,x

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_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
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Cd1 Cd 4 b 0.50000 0.50000 0.50000 1.00000
Sm2 Sm 4 c 0.25000 0.25000 0.25000 1.00000
Cd2 Cd 4 d 0.75000 0.75000 0.75000 1.00000
Cd3 Cd 16 e 0.08340 0.08340 0.08340 1.00000
Cd4 Cd 16 e 0.91260 0.91260 0.91260 1.00000
Cd5 Cd 16 e 0.16360 0.16360 0.16360 1.00000
Cd6 Cd 16 e 0.82970 0.82970 0.82970 1.00000
Sm3 Sm 16 e 0.40590 0.40590 0.40590 1.00000
Sm4 Sm 16 e 0.66180 0.66180 0.66180 1.00000
Cd7 Cd 24 f 0.15730 0.00000 0.00000 1.00000
Cd8 Cd 24 g 0.08950 0.25000 0.25000 1.00000
Cd9 Cd 48 h 0.29580 0.29580 0.39040 1.00000
Cd10 Cd 48 h 0.43770 0.43770 0.26270 1.00000
Cd11 Cd 48 h 0.54550 0.54550 0.64030 1.00000
```

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Cd12 Cd 48 h 0.67280 0.67280 0.51230 1.00000
Cd13 Cd 48 h 0.91610 0.91610 0.76370 1.00000
Sm5 Sm 48 h 0.17350 0.17350 0.01420 1.00000
```

$Sm_{11}Cd_{45}$: A45B11_cF448_216_bd4efg5h_ac2eh - POSCAR

```
A45B11_cF448_216_bd4efg5h_ac2eh & a,x5,x6,x7,x8,x9,x10,x11,x12,x13,z13,
    ↪ x14,z14,x15,z15,x16,z16,x17,z17,x18,z18 --params=21.699,0.0834,
    ↪ 0.9126,0.1636,0.8297,0.4059,0.6618,0.1573,0.0895,0.2958,0.3904,
    ↪ 0.4377,0.2627,0.5455,0.6403,0.6728,0.5123,0.9161,0.7637,0.1735,
    ↪ 0.0142 & F-43m T_{d}^{2} #216 (abcde^6fgh^6) & cF448 & None &
    ↪ Cd45Sm11 & Cd45Sm11 & M. L. Fornasini and B. Chabot and E.
    ↪ Parth\'{e}, Acta Crystallogr. Sect. B Struct. Sci. 34,
    ↪ 2093-2099 (1978)
   1.00000000000000
   0.00000000000000  10.84950000000000  10.84950000000000
  10.84950000000000   0.00000000000000  10.84950000000000
  10.84950000000000  10.84950000000000   0.00000000000000
    Cd    Sm
    90    22
Direct
   0.50000000000000   0.50000000000000   0.50000000000000  Cd  (4b)
   0.75000000000000   0.75000000000000   0.75000000000000  Cd  (4d)
   0.08340000000000   0.08340000000000   0.08340000000000  Cd  (16e)
   0.08340000000000   0.08340000000000  -0.25020000000000  Cd  (16e)
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   0.82970000000000   0.82970000000000  -2.48910000000000  Cd  (16e)
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   0.15730000000000  -0.15730000000000  -0.15730000000000  Cd  (24f)
   0.15730000000000  -0.15730000000000   0.15730000000000  Cd  (24f)
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   0.66180000000000   0.66180000000000  -1.98540000000000  Sm  (16e)
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   0.01420000000000   0.01420000000000  -0.36120000000000  Sm  (48h)
   0.33280000000000  -0.36120000000000   0.01420000000000  Sm  (48h)
  -0.36120000000000   0.33280000000000   0.01420000000000  Sm  (48h)
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  -0.36120000000000   0.01420000000000   0.01420000000000  Sm  (48h)
   0.01420000000000   0.33280000000000  -0.36120000000000  Sm  (48h)
   0.01420000000000  -0.36120000000000   0.33280000000000  Sm  (48h)
   0.01420000000000   0.33280000000000   0.01420000000000  Sm  (48h)
   0.01420000000000  -0.36120000000000   0.01420000000000  Sm  (48h)
  -0.36120000000000   0.01420000000000   0.33280000000000  Sm  (48h)
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```

Hg_2TiCu Inverse Heusler: AB2C_cF16_216_b_ad_c - CIF

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 'M. Pu\v{s}elj'
 'Z. Ban'
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;
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;
_journal_volume 41
_journal_year 1969
_journal_page_first 79
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;
 The Crystal Structure of TiCuHg$_{2}$
;

# Found in TiCuHg$_{2}$ (CuHg$_{2}$Ti) Crystal Structure, {PAULING FILE
    ↪ in: Inorganic Solid Phases, SpringerMaterials (online database)
    ↪ , Springer, Heidelberg (ed.)}, 2016

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4 -x,-y,z
5 y,z,x
6 y,-z,-x
7 -y,z,-x
8 -y,-z,x
9 z,x,y
10 z,-x,-y
11 -z,x,-y
12 -z,-x,y
13 y,x,z
14 y,-x,-z
15 -y,x,-z
16 -y,-x,z
17 x,z,y
18 x,-z,-y
19 -x,z,-y
20 -x,-z,y
21 z,y,x
22 z,-y,-x
23 -z,y,-x
24 -z,-y,x
25 x,y+1/2,z+1/2
26 x,-y+1/2,-z+1/2
```

```
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31 -y,z+1/2,-x+1/2
32 -y,-z+1/2,x+1/2
33 z,x+1/2,y+1/2
34 z,-x+1/2,-y+1/2
35 -z,x+1/2,-y+1/2
36 -z,-x+1/2,y+1/2
37 y,x+1/2,z+1/2
38 y,-x+1/2,-z+1/2
39 -y,x+1/2,-z+1/2
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82 z+1/2,-x+1/2,-y
83 -z+1/2,x+1/2,-y
84 -z+1/2,-x+1/2,y
85 y+1/2,x+1/2,z
86 y+1/2,-x+1/2,-z
87 -y+1/2,x+1/2,-z
88 -y+1/2,-x+1/2,z
89 x+1/2,z+1/2,y
90 x+1/2,-z+1/2,-y
91 -x+1/2,z+1/2,-y
92 -x+1/2,-z+1/2,y
93 z+1/2,y+1/2,x
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95 -z+1/2,y+1/2,-x
96 -z+1/2,-y+1/2,x

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_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
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Cu1 Cu 4 b 0.50000 0.50000 0.50000 1.00000
Ti1 Ti 4 c 0.25000 0.25000 0.25000 1.00000
Hg2 Hg 4 d 0.75000 0.75000 0.75000 1.00000
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Hg_2TiCu Inverse Heusler: AB2C_cF16_216_b_ad_c - POSCAR

```
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    ↪ Z. Ban, Croat. Chem. Acta 41, 79-83 (1969)
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   3.07747000000000   3.07747000000000   0.00000000000000
    Cu    Hg    Ti
     1     2     1
Direct
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   0.00000000000000   0.00000000000000   0.00000000000000   Hg   (4a)
   0.75000000000000   0.75000000000000   0.75000000000000   Hg   (4d)
   0.25000000000000   0.25000000000000   0.25000000000000   Ti   (4c)
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$GaMo_4S_8$: AB4C8_cF52_216_a_e_2e - CIF

```
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 'J. C. Jegaden'
 'M. Potel'
 'R. Chevrel'
 'M. Sergent'
 'A. Berton'
 'J. Chaussy'
 'A. K. Rastogi'
 'R. Tournier'
_journal_name_full_name
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 Journal of Solid State Chemistry
;
_journal_volume 51
_journal_year 1984
_journal_page_first 212
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 Nouveaux chalcogenures mixtes GaMo$_{4}$($XX$\')$_{8}$ ($X$ = S, Se, Te
    ↪ ) \'{a} clusters tetraedriques Mo$_{4}$
;

# Found in The American Mineralogist Crystal Structure Database, 2003

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_symmetry_Int_Tables_number 216

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_space_group_symop_operation_xyz
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2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
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6 y,-z,-x
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8 -y,-z,x
9 z,x,y
10 z,-x,-y
11 -z,x,-y
12 -z,-x,y
13 y,x,z
14 y,-x,-z
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16 -y,-x,z
17 x,z,y
18 x,-z,-y
19 -x,z,-y
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21 z,y,x
22 z,-y,-x
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36 -z,-x+1/2,y+1/2
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38 y,-x+1/2,-z+1/2
39 -y,x+1/2,-z+1/2
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43 -x,z+1/2,-y+1/2
44 -x,-z+1/2,y+1/2
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46 z,-y+1/2,-x+1/2
47 -z,y+1/2,-x+1/2
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49 x+1/2,y,z+1/2
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51 -x+1/2,y,-z+1/2
52 -x+1/2,-y,z+1/2
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58 z+1/2,-x,-y+1/2
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92 -x+1/2,-z+1/2,y
93 z+1/2,y+1/2,x
94 z+1/2,-y+1/2,-x
95 -z+1/2,y+1/2,-x
96 -z+1/2,-y+1/2,x

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ga1 Ga 4 a 0.00000 0.00000 0.00000 1.00000
Mo1 Mo 16 e 0.39740 0.39740 0.39740 1.00000
S1 S 16 e 0.63430 0.63430 0.63430 1.00000
S2 S 16 e 0.13500 0.13500 0.13500 1.00000
```

$GaMo_4S_8$: AB4C8_cF52_216_a_e_2e - POSCAR

```
AB4C8_cF52_216_a_e_2e & a,x2,x3,x4 --params=9.7294,0.3974,0.6343,0.135 &
    ↪ F-43m T_{d}^{2} #216 (ae^3) & cF52 & None & GaMo4S8 & GaMo4S8
    ↪ & H. {Ben Yaich} et al., J. Solid State Chem. 51, 212-217 (1984
    ↪ )
   1.00000000000000
   0.00000000000000   4.86470000000000   4.86470000000000
   4.86470000000000   0.00000000000000   4.86470000000000
   4.86470000000000   4.86470000000000   0.00000000000000
    Ga    Mo    S
     1     4     8
Direct
   0.00000000000000   0.00000000000000   0.00000000000000   Ga   (4a)
   0.39740000000000   0.39740000000000   0.39740000000000   Mo   (16e)
   0.39740000000000   0.39740000000000  -1.19220000000000   Mo   (16e)
   0.39740000000000  -1.19220000000000   0.39740000000000   Mo   (16e)
  -1.19220000000000   0.39740000000000   0.39740000000000   Mo   (16e)
   0.63430000000000   0.63430000000000   0.63430000000000   S    (16e)
   0.63430000000000   0.63430000000000  -1.90290000000000   S    (16e)
   0.63430000000000  -1.90290000000000   0.63430000000000   S    (16e)
  -1.90290000000000   0.63430000000000   0.63430000000000   S    (16e)
   0.13500000000000   0.13500000000000   0.13500000000000   S    (16e)
   0.13500000000000   0.13500000000000  -0.40500000000000   S    (16e)
   0.13500000000000  -0.40500000000000   0.13500000000000   S    (16e)
  -0.40500000000000   0.13500000000000   0.13500000000000   S    (16e)
```

High-Temperature Cubic $KClO_4$ ($H0_5$): ABC4_cF24_216_b_a_e - CIF

```
# CIF file
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_audit_creation_method FINDSYM

_chemical_name_mineral 'ClKO4'
_chemical_formula_sum 'Cl K O4'

loop_
_publ_author_name
 'K. Hermann'
 'W. Ilge'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 71
_journal_year 1930
_journal_page_first 41
_journal_page_last 66
_publ_Section_title
;
 R\"{o}ntgenographische Strukturerforschung der kubischen Modifikation
    ↪ der Perchlorate
;

# Found in Strukturbericht Band II 1928-1932, 1937

_aflow_title 'High-Temperature Cubic KClO$_{4}$ ($H0_{5}$) Structure'
_aflow_proto 'ABC4_cF24_216_b_a_e'
_aflow_params 'a,x_{3}'
_aflow_params_values '7.5,0.125'
_aflow_Strukturbericht '$H0_{5}$'
_aflow_Pearson 'cF24'

_symmetry_space_group_name_H-M "F -4 3 m"
_symmetry_Int_Tables_number 216

_cell_length_a     7.50000
_cell_length_b     7.50000
_cell_length_c     7.50000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 y,z,x
6 y,-z,-x
7 -y,z,-x
8 -y,-z,x
9 z,x,y
10 z,-x,-y
11 -z,x,-y
12 -z,-x,y
13 y,x,z
14 y,-x,-z
15 -y,x,-z
16 -y,-x,z
17 x,z,y
18 x,-z,-y
19 -x,z,-y
20 -x,-z,y
21 z,y,x
22 z,-y,-x
23 -z,y,-x
24 -z,-y,x
25 x,y+1/2,z+1/2
26 x,-y+1/2,-z+1/2
27 -x,y+1/2,-z+1/2
28 -x,-y+1/2,z+1/2
29 y,z+1/2,x+1/2
30 y,-z+1/2,-x+1/2
31 -y,z+1/2,-x+1/2
32 -y,-z+1/2,x+1/2
33 z,x+1/2,y+1/2
34 z,-x+1/2,-y+1/2
35 -z,x+1/2,-y+1/2
36 -z,-x+1/2,y+1/2
37 y,x+1/2,z+1/2
38 y,-x+1/2,-z+1/2
39 -y,x+1/2,-z+1/2
40 -y,-x+1/2,z+1/2
41 x,z+1/2,y+1/2
42 x,-z+1/2,-y+1/2
43 -x,z+1/2,-y+1/2
44 -x,-z+1/2,y+1/2
45 z,y+1/2,x+1/2
46 z,-y+1/2,-x+1/2
47 -z,y+1/2,-x+1/2
48 -z,-y+1/2,x+1/2
49 x+1/2,y,z+1/2
50 x+1/2,-y,-z+1/2
51 -x+1/2,y,-z+1/2
52 -x+1/2,-y,z+1/2
53 y+1/2,z,x+1/2
54 y+1/2,-z,-x+1/2
55 -y+1/2,z,-x+1/2
56 -y+1/2,-z,x+1/2
57 z+1/2,x,y+1/2
58 z+1/2,-x,-y+1/2
59 -z+1/2,x,-y+1/2
60 -z+1/2,-x,y+1/2
61 y+1/2,x,z+1/2
62 y+1/2,-x,-z+1/2
63 -y+1/2,x,-z+1/2
64 -y+1/2,-x,z+1/2
65 x+1/2,z,y+1/2
66 x+1/2,-z,-y+1/2
67 -x+1/2,z,-y+1/2
68 -x+1/2,-z,y+1/2
69 z+1/2,y,x+1/2
70 z+1/2,-y,-x+1/2
71 -z+1/2,y,-x+1/2
72 -z+1/2,-y,x+1/2
73 x+1/2,y+1/2,z
74 x+1/2,-y+1/2,-z
75 -x+1/2,y+1/2,-z
```

```
76 -x+1/2,-y+1/2,z
77 y+1/2,z+1/2,x
78 y+1/2,-z+1/2,-x
79 -y+1/2,z+1/2,-x
80 -y+1/2,-z+1/2,x
81 z+1/2,x+1/2,y
82 z+1/2,-x+1/2,-y
83 -z+1/2,x+1/2,-y
84 -z+1/2,-x+1/2,y
85 y+1/2,x+1/2,z
86 y+1/2,-x+1/2,-z
87 -y+1/2,x+1/2,-z
88 -y+1/2,-x+1/2,z
89 x+1/2,z+1/2,y
90 x+1/2,-z+1/2,-y
91 -x+1/2,z+1/2,-y
92 -x+1/2,-z+1/2,y
93 z+1/2,y+1/2,x
94 z+1/2,-y+1/2,-x
95 -z+1/2,y+1/2,-x
96 -z+1/2,-y+1/2,x

loop_
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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
K1 K 4 a 0.00000 0.00000 0.00000 1.00000
Cl1 Cl 4 b 0.50000 0.50000 0.50000 1.00000
O1 O 16 e 0.12500 0.12500 0.12500 1.00000
```

High-Temperature Cubic $KClO_4$ ($H0_5$): ABC4_cF24_216_b_a_e - POSCAR

```
ABC4_cF24_216_b_a_e & a,x3 --params=7.5,0.125 & F-43m T_{d}^{2} #216 (
    ↪ abe) & cF24 & $H0_{5}$ & ClKO4 & ClKO4 & K. Hermann and W. Ilge
    ↪ , Zeitschrift f{"u}r Kristallographie - Crystalline Materials
    ↪ 71, 41-66 (1930)
   1.00000000000000
   0.00000000000000   3.75000000000000   3.75000000000000
   3.75000000000000   0.00000000000000   3.75000000000000
   3.75000000000000   3.75000000000000   0.00000000000000
    Cl   K   O
     1   1   4
Direct
   0.50000000000000   0.50000000000000   0.50000000000000   Cl   (4b)
   0.00000000000000   0.00000000000000   0.00000000000000   K    (4a)
   0.12500000000000   0.12500000000000   0.12500000000000   O    (16e)
   0.12500000000000   0.12500000000000  -0.37500000000000   O    (16e)
   0.12500000000000  -0.37500000000000   0.12500000000000   O    (16e)
  -0.37500000000000   0.12500000000000   0.12500000000000   O    (16e)
```

AlN (cF40): AB_cF40_216_de_ce - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'AlN'
_chemical_formula_sum 'Al N'

loop_
_publ_author_name
 'C. Liu'
 'M. Chen'
 'J. Li'
 'L. Liu'
 'P. Li'
 'M. Ma'
 'C. Shao'
 'J. He'
 'T. Liang'
_journal_name_full_name
;
 Journal of Physics and Chemistry of Solids
;
_journal_volume 130
_journal_year 2019
_journal_page_first 58
_journal_page_last 66
_publ_Section_title
;
 A first-principles study of novel cubic AlN phases
;

_aflow_title 'AlN (cF40) Structure'
_aflow_proto 'AB_cF40_216_de_ce'
_aflow_params 'a,x_{3},x_{4}'
_aflow_params_values '8.1926,0.1137,0.8804'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'cF40'

_symmetry_space_group_name_H-M "F -4 3 m"
_symmetry_Int_Tables_number 216

_cell_length_a     8.19260
_cell_length_b     8.19260
_cell_length_c     8.19260
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
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2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 y,z,x
6 y,-z,-x
7 -y,z,-x
8 -y,-z,x
9 z,x,y
10 z,-x,-y
11 -z,x,-y
12 -z,-x,y
13 y,x,z
14 y,-x,-z
15 -y,x,-z
16 -y,-x,z
17 x,z,y
18 x,-z,-y
19 -x,z,-y
20 -x,-z,y
21 z,y,x
22 z,-y,-x
23 -z,y,-x
24 -z,-y,x
25 x,y+1/2,z+1/2
26 x,-y+1/2,-z+1/2
27 -x,y+1/2,-z+1/2
28 -x,-y+1/2,z+1/2
29 y,z+1/2,x+1/2
30 y,-z+1/2,-x+1/2
31 -y,z+1/2,-x+1/2
32 -y,-z+1/2,x+1/2
33 z,x+1/2,y+1/2
34 z,-x+1/2,-y+1/2
35 -z,x+1/2,-y+1/2
36 -z,-x+1/2,y+1/2
37 y,x+1/2,z+1/2
38 y,-x+1/2,-z+1/2
39 -y,x+1/2,-z+1/2
40 -y,-x+1/2,z+1/2
41 x,z+1/2,y+1/2
42 x,-z+1/2,-y+1/2
43 -x,z+1/2,-y+1/2
44 -x,-z+1/2,y+1/2
45 z,y+1/2,x+1/2
46 z,-y+1/2,-x+1/2
47 -z,y+1/2,-x+1/2
48 -z,-y+1/2,x+1/2
49 x+1/2,y,z+1/2
50 x+1/2,-y,-z+1/2
51 -x+1/2,y,-z+1/2
52 -x+1/2,-y,z+1/2
53 y+1/2,z,x+1/2
54 y+1/2,-z,-x+1/2
55 -y+1/2,z,-x+1/2
56 -y+1/2,-z,x+1/2
57 z+1/2,x,y+1/2
58 z+1/2,-x,-y+1/2
59 -z+1/2,x,-y+1/2
60 -z+1/2,-x,y+1/2
61 y+1/2,x,z+1/2
62 y+1/2,-x,-z+1/2
63 -y+1/2,x,-z+1/2
64 -y+1/2,-x,z+1/2
65 x+1/2,z,y+1/2
66 x+1/2,-z,-y+1/2
67 -x+1/2,z,-y+1/2
68 -x+1/2,-z,y+1/2
69 z+1/2,y,x+1/2
70 z+1/2,-y,-x+1/2
71 -z+1/2,y,-x+1/2
72 -z+1/2,-y,x+1/2
73 x+1/2,y+1/2,z
74 x+1/2,-y+1/2,-z
75 -x+1/2,y+1/2,-z
76 -x+1/2,-y+1/2,z
77 y+1/2,z+1/2,x
78 y+1/2,-z+1/2,-x
79 -y+1/2,z+1/2,-x
80 -y+1/2,-z+1/2,x
81 z+1/2,x+1/2,y
82 z+1/2,-x+1/2,-y
83 -z+1/2,x+1/2,-y
84 -z+1/2,-x+1/2,y
85 y+1/2,x+1/2,z
86 y+1/2,-x+1/2,-z
87 -y+1/2,x+1/2,-z
88 -y+1/2,-x+1/2,z
89 x+1/2,z+1/2,y
90 x+1/2,-z+1/2,-y
91 -x+1/2,z+1/2,-y
92 -x+1/2,-z+1/2,y
93 z+1/2,y+1/2,x
94 z+1/2,-y+1/2,-x
95 -z+1/2,y+1/2,-x
96 -z+1/2,-y+1/2,x

loop_
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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
```

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_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
N1 N 4 c 0.25000 0.25000 0.25000 1.00000
Al1 Al 4 d 0.75000 0.75000 0.75000 1.00000
Al2 Al 16 e 0.11370 0.11370 0.11370 1.00000
N2 N 16 e 0.88040 0.88040 0.88040 1.00000
```

AlN (cF40): AB_cF40_216_de_ce - POSCAR

```
AB_cF40_216_de_ce & a,x3,x4 --params=8.1926,0.1137,0.8804 & F-43m T_{d
    ↪ }^{2} #216 (cde^2) & cF40 & None & AlN & AlN & C. Liu et al.,
    ↪ J. Phys. Chem. Solids 130, 58-66 (2019)
   1.00000000000000
   0.00000000000000   4.09630000000000   4.09630000000000
   4.09630000000000   0.00000000000000   4.09630000000000
   4.09630000000000   4.09630000000000   0.00000000000000
    Al    N
     5    5
Direct
   0.75000000000000   0.75000000000000   0.75000000000000   Al   (4d)
   0.11370000000000   0.11370000000000   0.11370000000000   Al   (16e)
   0.11370000000000   0.11370000000000  -0.34110000000000   Al   (16e)
   0.11370000000000  -0.34110000000000   0.11370000000000   Al   (16e)
  -0.34110000000000   0.11370000000000   0.11370000000000   Al   (16e)
   0.25000000000000   0.25000000000000   0.25000000000000   N    (4c)
   0.88040000000000   0.88040000000000   0.88040000000000   N    (16e)
   0.88040000000000   0.88040000000000  -2.64120000000000   N    (16e)
   0.88040000000000  -2.64120000000000   0.88040000000000   N    (16e)
  -2.64120000000000   0.88040000000000   0.88040000000000   N    (16e)
```

Tennantite ($Cu_{12}As_4S_{13}$): A4B24C13_cI82_217_c_deg_ag - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

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_chemical_formula_sum 'As4 Cu24 S13'

loop_
_publ_author_name
 'A. A. Yaroslavzev'
 'A. V. Mironov'
 'A. N. Kuznetsov'
 'A. P. Dudka'
 'O. N. Khrykina'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 75
_journal_year 2019
_journal_page_first 634
_journal_page_last 642
_publ_Section_title
;
 Tennantite: multi-temperature crystal structure, phase transition and
    ↪ electronic structure of synthetic Cu$_{12}$As$_{4}$S$_{13}$
;

_aflow_title 'Tennantite (Cu$_{12}$As$_{4}$S$_{13}$) Structure'
_aflow_proto 'A4B24C13_cI82_217_c_deg_ag'
_aflow_params 'a,x_{2},x_{4},x_{5},z_{5},x_{6},z_{6}'
_aflow_params_values '10.1439,0.24271,0.2173,0.0777,0.2123,0.61822,
    ↪ 0.14266'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'cI82'

_symmetry_space_group_name_H-M "I -4 3 m"
_symmetry_Int_Tables_number 217

_cell_length_a     10.14390
_cell_length_b     10.14390
_cell_length_c     10.14390
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 y,z,x
6 y,-z,-x
7 -y,z,-x
8 -y,-z,x
9 z,x,y
10 z,-x,-y
11 -z,x,-y
12 -z,-x,y
13 y,x,z
14 y,-x,-z
15 -y,x,-z
16 -y,-x,z
17 x,z,y
18 x,-z,-y
19 -x,z,-y
20 -x,-z,y
21 z,y,x
22 z,-y,-x
23 -z,y,-x
24 -z,-y,x
25 x+1/2,y+1/2,z+1/2
26 x+1/2,-y+1/2,-z+1/2
27 -x+1/2,y+1/2,-z+1/2
28 -x+1/2,-y+1/2,z+1/2
29 y+1/2,z+1/2,x+1/2
30 y+1/2,-z+1/2,-x+1/2
31 -y+1/2,z+1/2,-x+1/2
32 -y+1/2,-z+1/2,x+1/2
33 z+1/2,x+1/2,y+1/2
34 z+1/2,-x+1/2,-y+1/2
35 -z+1/2,x+1/2,-y+1/2
36 -z+1/2,-x+1/2,y+1/2
37 y+1/2,x+1/2,z+1/2
38 y+1/2,-x+1/2,-z+1/2
39 -y+1/2,x+1/2,-z+1/2
40 -y+1/2,-x+1/2,z+1/2
41 x+1/2,z+1/2,y+1/2
42 x+1/2,-z+1/2,-y+1/2
43 -x+1/2,z+1/2,-y+1/2
44 -x+1/2,-z+1/2,y+1/2
45 z+1/2,y+1/2,x+1/2
46 z+1/2,-y+1/2,-x+1/2
47 -z+1/2,y+1/2,-x+1/2
48 -z+1/2,-y+1/2,x+1/2

loop_
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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
S1 S 2 a 0.00000 0.00000 0.00000 1.00000
As1 As 8 c 0.24271 0.24271 0.24271 1.00000
Cu1 Cu 12 d 0.25000 0.50000 0.00000 1.00000
Cu2 Cu 12 e 0.21730 0.00000 0.00000 0.75800
Cu3 Cu 24 g 0.07770 0.07770 0.21230 0.12100
S2 S 24 g 0.61822 0.61822 0.14266 1.00000
```

Tennantite ($Cu_{12}As_4S_{13}$): A4B24C13_cI82_217_c_deg_ag - POSCAR

```
A4B24C13_cI82_217_c_deg_ag & a,x2,x4,x5,z5,x6,z6 --params=10.1439,
    ↪ 0.24271,0.2173,0.0777,0.2123,0.61822,0.14266 & I-43m T_{d}^{3}
    ↪ #217 (acdeg^2) & cI82 & None & AsCu12S13 & Tennantite & A. A.
    ↪ Yaroslavzev et al., Acta Crystallogr. Sect. B Struct. Sci. 75,
    ↪ 634-642 (2019)
   1.00000000000000
  -5.07195000000000   5.07195000000000   5.07195000000000
   5.07195000000000  -5.07195000000000   5.07195000000000
   5.07195000000000   5.07195000000000  -5.07195000000000
    As    Cu    S
     4    24    13
Direct
   0.48542000000000   0.48542000000000   0.48542000000000   As   (8c)
   0.00000000000000   0.00000000000000  -0.48542000000000   As   (8c)
   0.00000000000000  -0.48542000000000   0.00000000000000   As   (8c)
  -0.48542000000000   0.00000000000000   0.00000000000000   As   (8c)
   0.50000000000000   0.25000000000000   0.75000000000000   Cu   (12d)
   0.50000000000000   0.75000000000000   0.25000000000000   Cu   (12d)
   0.75000000000000   0.50000000000000   0.25000000000000   Cu   (12d)
   0.25000000000000   0.50000000000000   0.75000000000000   Cu   (12d)
   0.25000000000000   0.75000000000000   0.50000000000000   Cu   (12d)
   0.75000000000000   0.25000000000000   0.50000000000000   Cu   (12d)
   0.00000000000000   0.21730000000000   0.21730000000000   Cu   (12e)
   0.00000000000000  -0.21730000000000  -0.21730000000000   Cu   (12e)
   0.21730000000000   0.00000000000000   0.21730000000000   Cu   (12e)
  -0.21730000000000   0.00000000000000  -0.21730000000000   Cu   (12e)
   0.21730000000000   0.21730000000000   0.00000000000000   Cu   (12e)
  -0.21730000000000  -0.21730000000000   0.00000000000000   Cu   (12e)
   0.29000000000000   0.29000000000000   0.15540000000000   Cu   (24g)
   0.13460000000000   0.13460000000000  -0.15540000000000   Cu   (24g)
  -0.13460000000000  -0.29000000000000   0.00000000000000   Cu   (24g)
  -0.29000000000000  -0.13460000000000   0.00000000000000   Cu   (24g)
   0.15540000000000   0.29000000000000   0.29000000000000   Cu   (24g)
  -0.15540000000000   0.13460000000000   0.13460000000000   Cu   (24g)
   0.00000000000000  -0.13460000000000  -0.29000000000000   Cu   (24g)
   0.00000000000000  -0.29000000000000  -0.13460000000000   Cu   (24g)
   0.29000000000000   0.15540000000000   0.29000000000000   Cu   (24g)
   0.13460000000000  -0.15540000000000   0.13460000000000   Cu   (24g)
  -0.29000000000000   0.00000000000000  -0.13460000000000   Cu   (24g)
  -0.13460000000000   0.00000000000000  -0.29000000000000   Cu   (24g)
   0.00000000000000   0.00000000000000   0.00000000000000   S    (2a)
   0.76088000000000   0.76088000000000   1.23644000000000   S    (24g)
  -0.47556000000000  -0.47556000000000  -1.23644000000000   S    (24g)
   0.47556000000000  -0.76088000000000   0.00000000000000   S    (24g)
  -0.76088000000000   0.47556000000000   0.00000000000000   S    (24g)
   1.23644000000000   0.76088000000000   0.76088000000000   S    (24g)
  -1.23644000000000  -0.47556000000000  -0.47556000000000   S    (24g)
   0.00000000000000   0.47556000000000  -0.76088000000000   S    (24g)
   0.00000000000000  -0.76088000000000   0.47556000000000   S    (24g)
   0.76088000000000   1.23644000000000   0.76088000000000   S    (24g)
  -0.47556000000000  -1.23644000000000  -0.47556000000000   S    (24g)
  -0.76088000000000   0.00000000000000   0.47556000000000   S    (24g)
   0.47556000000000   0.00000000000000  -0.76088000000000   S    (24g)
```

AlN (cI16): AB_cI16_217_c_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM
```

```
_chemical_name_mineral 'AlN'
_chemical_formula_sum 'Al N'

loop_
_publ_author_name
 'C. Liu'
 'M. Chen'
 'J. Li'
 'L. Liu'
 'P. Li'
 'M. Ma'
 'C. Shao'
 'J. He'
 'T. Liang'
_journal_name_full_name
;
 Journal of Physics and Chemistry of Solids
;
_journal_volume 130
_journal_year 2019
_journal_page_first 58
_journal_page_last 66
_publ_Section_title
;
 A first-principles study of novel cubic AlN phases
;

_aflow_title 'AlN (cI16) Structure'
_aflow_proto 'AB_cI16_217_c_c'
_aflow_params 'a,x_{1},x_{2}'
_aflow_params_values '5.9487,0.3409,0.1657'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'cI16'

_symmetry_space_group_name_H-M "I -4 3 m"
_symmetry_Int_Tables_number 217

_cell_length_a     5.94870
_cell_length_b     5.94870
_cell_length_c     5.94870
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 y,z,x
6 y,-z,-x
7 -y,z,-x
8 -y,-z,x
9 z,x,y
10 z,-x,-y
11 -z,x,-y
12 -z,-x,y
13 y,x,z
14 y,-x,-z
15 -y,x,-z
16 -y,-x,z
17 x,z,y
18 x,-z,-y
19 -x,z,-y
20 -x,-z,y
21 z,y,x
22 z,-y,-x
23 -z,y,-x
24 -z,-y,x
25 x+1/2,y+1/2,z+1/2
26 x+1/2,-y+1/2,-z+1/2
27 -x+1/2,y+1/2,-z+1/2
28 -x+1/2,-y+1/2,z+1/2
29 y+1/2,z+1/2,x+1/2
30 y+1/2,-z+1/2,-x+1/2
31 -y+1/2,z+1/2,-x+1/2
32 -y+1/2,-z+1/2,x+1/2
33 z+1/2,x+1/2,y+1/2
34 z+1/2,-x+1/2,-y+1/2
35 -z+1/2,x+1/2,-y+1/2
36 -z+1/2,-x+1/2,y+1/2
37 y+1/2,x+1/2,z+1/2
38 y+1/2,-x+1/2,-z+1/2
39 -y+1/2,x+1/2,-z+1/2
40 -y+1/2,-x+1/2,z+1/2
41 x+1/2,z+1/2,y+1/2
42 x+1/2,-z+1/2,-y+1/2
43 -x+1/2,z+1/2,-y+1/2
44 -x+1/2,-z+1/2,y+1/2
45 z+1/2,y+1/2,x+1/2
46 z+1/2,-y+1/2,-x+1/2
47 -z+1/2,y+1/2,-x+1/2
48 -z+1/2,-y+1/2,x+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Al1 Al 8 c 0.34090 0.34090 0.34090 1.00000
N1 N 8 c 0.16570 0.16570 0.16570 1.00000
```

AlN (cI16): AB_cI16_217_c_c - POSCAR

```
AB_cI16_217_c_c & a,x1,x2 --params=5.9487,0.3409,0.1657 & I-43m T_{d}^{3
    ↪ } #217 (c^2) & cI16 & None & AlN & AlN & C. Liu et al., J.
    ↪ Phys. Chem. Solids 130, 58-66 (2019)
   1.00000000000000
  -2.97435000000000    2.97435000000000    2.97435000000000
   2.97435000000000   -2.97435000000000    2.97435000000000
   2.97435000000000    2.97435000000000   -2.97435000000000
    Al      N
     4      4
Direct
   0.68180000000000    0.68180000000000    0.68180000000000    Al    (8c)
   0.00000000000000    0.00000000000000   -0.68180000000000    Al    (8c)
   0.00000000000000   -0.68180000000000    0.00000000000000    Al    (8c)
  -0.68180000000000    0.00000000000000    0.00000000000000    Al    (8c)
   0.33140000000000    0.33140000000000    0.33140000000000    N     (8c)
   0.00000000000000    0.00000000000000   -0.33140000000000    N     (8c)
   0.00000000000000   -0.33140000000000    0.00000000000000    N     (8c)
  -0.33140000000000    0.00000000000000    0.00000000000000    N     (8c)
```

Hauyne [$(Na_{0.5}Ca_{0.3}K_{0.2})_8(Al_6Si_6O_{24})(SO_4)_{1.5}$, $S6_9$]: A3B4C4D4E16F4G3_cP76_218_c_e_e_e_ei_e_d - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Hauyne'
_chemical_formula_sum 'Al3 Ca4 K4 Na4 O16 S4 Si3'

loop_
_publ_author_name
 'I. Hassan'
 'H. D. Grundy'
_journal_name_full_name
;
 Canadian Mineralogist
;
_journal_volume 29
_journal_year 1991
_journal_page_first 123
_journal_page_last 130
_publ_Section_title
;
 The Crystal Structure of Hauyne at 293 and 153 K
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'Hauyne [(Na$_{0.5}$Ca$_{0.3}$K$_{0.2}$)$_{8}$(Al$_{6}$Si$_
    ↪ {6}$O$_{24}$)(SO$_{4}$)$_{1.5}$, $S6_{9}$] Structure'
_aflow_proto 'A3B4C4D4E16F4G3_cP76_218_c_e_e_e_ei_e_d'
_aflow_params 'a,x_{3},x_{4},x_{5},x_{6},x_{7},x_{8},y_{8},z_{8}'
_aflow_params_values '9.1097,0.2009,0.164,0.2392,0.0995,-0.0331,0.6443,
    ↪ 0.6558,-0.0331'
_aflow_Strukturbericht '$S6_{9}$'
_aflow_Pearson 'cP76'

_symmetry_space_group_name_H-M "P -4 3 n"
_symmetry_Int_Tables_number 218

_cell_length_a     9.10970
_cell_length_b     9.10970
_cell_length_c     9.10970
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 y,z,x
6 y,-z,-x
7 -y,z,-x
8 -y,-z,x
9 z,x,y
10 z,-x,-y
11 -z,x,-y
12 -z,-x,y
13 y+1/2,x+1/2,z+1/2
14 y+1/2,-x+1/2,-z+1/2
15 -y+1/2,x+1/2,-z+1/2
16 -y+1/2,-x+1/2,z+1/2
17 x+1/2,z+1/2,y+1/2
18 x+1/2,-z+1/2,-y+1/2
19 -x+1/2,z+1/2,-y+1/2
20 -x+1/2,-z+1/2,y+1/2
21 z+1/2,y+1/2,x+1/2
22 z+1/2,-y+1/2,-x+1/2
23 -z+1/2,y+1/2,-x+1/2
24 -z+1/2,-y+1/2,x+1/2

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
```

```
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Al1 Al 6 c 0.25000 0.50000 0.00000 1.00000
Si1 Si 6 d 0.25000 0.00000 0.50000 1.00000
Ca1 Ca 8 e 0.20090 0.20090 0.20090 0.30000
K1 K 8 e 0.16400 0.16400 0.16400 0.20000
Na1 Na 8 e 0.23920 0.23920 0.23920 0.54000
O1 O 8 e 0.09950 0.09950 0.09950 0.75000
S1 S 8 e -0.03310 -0.03310 -0.03310 0.19000
O2 O 24 i 0.64430 0.65580 -0.03310 1.00000
```

Hauyne $[(Na_{0.5}Ca_{0.3}K_{0.2})_8(Al_6Si_6O_{24})(SO_4)_{1.5}, S6_9]$: A3B4C4D4E16F4G3_cP76_218_c_e_e_e_ei_e_d - POSCAR

```
A3B4C4D4E16F4G3_cP76_218_c_e_e_e_ei_e_d & a,x3,x4,x5,x6,x7,x8,y8,z8 --
    ↪ params=9.1097,0.2009,0.164,0.2392,0.0995,-0.0331,0.6443,0.6558
    ↪ ,-0.0331 & P-43n T_{d}^{4} #218 (cde^5i) & cP76 & $S6_{9}$ &
    ↪ Al6Ca2.4K1.6Na4O30S1.5Si6 & Hauyne & I. Hassan and H. D. Grundy
    ↪ , Can. Mineral. 29, 123-130 (1991)
   1.00000000000000
   9.10970000000000   0.00000000000000   0.00000000000000
   0.00000000000000   9.10970000000000   0.00000000000000
   0.00000000000000   0.00000000000000   9.10970000000000
    Al  Ca  K  Na  O  S  Si
     6   8  8   8 32  8   6
Direct
   0.25000000000000   0.50000000000000   0.00000000000000  Al  (6c)
   0.75000000000000   0.50000000000000   0.00000000000000  Al  (6c)
   0.00000000000000   0.25000000000000   0.50000000000000  Al  (6c)
   0.00000000000000   0.75000000000000   0.50000000000000  Al  (6c)
   0.50000000000000   0.00000000000000   0.25000000000000  Al  (6c)
   0.50000000000000   0.00000000000000   0.75000000000000  Al  (6c)
   0.20090000000000   0.20090000000000   0.20090000000000  Ca  (8e)
  -0.20090000000000  -0.20090000000000   0.20090000000000  Ca  (8e)
  -0.20090000000000   0.20090000000000  -0.20090000000000  Ca  (8e)
   0.20090000000000  -0.20090000000000  -0.20090000000000  Ca  (8e)
   0.70090000000000   0.70090000000000   0.70090000000000  Ca  (8e)
   0.29910000000000   0.29910000000000   0.70090000000000  Ca  (8e)
   0.70090000000000   0.29910000000000   0.29910000000000  Ca  (8e)
   0.29910000000000   0.70090000000000   0.29910000000000  Ca  (8e)
   0.16400000000000   0.16400000000000   0.16400000000000  K  (8e)
  -0.16400000000000  -0.16400000000000   0.16400000000000  K  (8e)
  -0.16400000000000   0.16400000000000  -0.16400000000000  K  (8e)
   0.16400000000000  -0.16400000000000  -0.16400000000000  K  (8e)
   0.66400000000000   0.66400000000000   0.66400000000000  K  (8e)
   0.33600000000000   0.33600000000000   0.66400000000000  K  (8e)
   0.66400000000000   0.33600000000000   0.33600000000000  K  (8e)
   0.33600000000000   0.66400000000000   0.33600000000000  K  (8e)
   0.23920000000000   0.23920000000000   0.23920000000000  Na  (8e)
  -0.23920000000000  -0.23920000000000   0.23920000000000  Na  (8e)
  -0.23920000000000   0.23920000000000  -0.23920000000000  Na  (8e)
   0.23920000000000  -0.23920000000000  -0.23920000000000  Na  (8e)
   0.73920000000000   0.73920000000000   0.73920000000000  Na  (8e)
   0.26080000000000   0.26080000000000   0.73920000000000  Na  (8e)
   0.73920000000000   0.26080000000000   0.26080000000000  Na  (8e)
   0.26080000000000   0.73920000000000   0.26080000000000  Na  (8e)
   0.09950000000000   0.09950000000000   0.09950000000000  O  (8e)
  -0.09950000000000  -0.09950000000000   0.09950000000000  O  (8e)
  -0.09950000000000   0.09950000000000  -0.09950000000000  O  (8e)
   0.09950000000000  -0.09950000000000  -0.09950000000000  O  (8e)
   0.59950000000000   0.59950000000000   0.59950000000000  O  (8e)
   0.40050000000000   0.40050000000000   0.59950000000000  O  (8e)
   0.59950000000000   0.40050000000000   0.40050000000000  O  (8e)
   0.40050000000000   0.59950000000000   0.40050000000000  O  (8e)
   0.64430000000000   0.65580000000000  -0.03310000000000  O  (24i)
  -0.64430000000000  -0.65580000000000  -0.03310000000000  O  (24i)
  -0.64430000000000   0.65580000000000   0.03310000000000  O  (24i)
   0.64430000000000  -0.65580000000000   0.03310000000000  O  (24i)
  -0.03310000000000   0.64430000000000   0.65580000000000  O  (24i)
  -0.03310000000000  -0.64430000000000  -0.65580000000000  O  (24i)
   0.03310000000000  -0.64430000000000   0.65580000000000  O  (24i)
   0.03310000000000   0.64430000000000  -0.65580000000000  O  (24i)
   0.65580000000000  -0.03310000000000   0.64430000000000  O  (24i)
  -0.65580000000000  -0.03310000000000  -0.64430000000000  O  (24i)
   0.65580000000000   0.03310000000000  -0.64430000000000  O  (24i)
  -0.65580000000000   0.03310000000000   0.64430000000000  O  (24i)
   1.15580000000000   1.14430000000000   0.46690000000000  O  (24i)
  -0.15580000000000  -0.14430000000000   0.46690000000000  O  (24i)
   1.15580000000000  -0.14430000000000   0.53310000000000  O  (24i)
  -0.15580000000000   1.14430000000000   0.53310000000000  O  (24i)
   1.14430000000000   0.46690000000000   1.15580000000000  O  (24i)
  -0.14430000000000   0.46690000000000  -0.15580000000000  O  (24i)
  -0.14430000000000   0.53310000000000   1.15580000000000  O  (24i)
   1.14430000000000   0.53310000000000  -0.15580000000000  O  (24i)
   0.46690000000000   1.15580000000000   1.14430000000000  O  (24i)
   0.46690000000000  -0.15580000000000  -0.14430000000000  O  (24i)
   0.53310000000000   1.15580000000000  -0.14430000000000  O  (24i)
   0.53310000000000  -0.15580000000000   1.14430000000000  O  (24i)
  -0.03310000000000  -0.03310000000000  -0.03310000000000  S  (8e)
   0.03310000000000   0.03310000000000  -0.03310000000000  S  (8e)
   0.03310000000000  -0.03310000000000   0.03310000000000  S  (8e)
  -0.03310000000000   0.03310000000000   0.03310000000000  S  (8e)
   0.46690000000000   0.46690000000000   0.46690000000000  S  (8e)
   0.53310000000000   0.53310000000000   0.46690000000000  S  (8e)
   0.46690000000000   0.53310000000000   0.53310000000000  S  (8e)
   0.53310000000000   0.46690000000000   0.53310000000000  S  (8e)
   0.25000000000000   0.00000000000000   0.50000000000000  Si  (6d)
   0.75000000000000   0.00000000000000   0.50000000000000  Si  (6d)
   0.50000000000000   0.25000000000000   0.00000000000000  Si  (6d)
   0.50000000000000   0.75000000000000   0.00000000000000  Si  (6d)
   0.00000000000000   0.50000000000000   0.25000000000000  Si  (6d)
   0.00000000000000   0.50000000000000   0.75000000000000  Si  (6d)
```

Sodalite $[Na_4(AlSiO_4)_3Cl, S6_2]$: A3BC4D12E3_cP46_218_d_a_e_i_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Sodalite'
_chemical_formula_sum 'Al3 Cl Na4 O12 Si3'

loop_
_publ_author_name
 'I. Hassan'
 'H. D. Grundy'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 40
_journal_year 1984
_journal_page_first 6
_journal_page_last 13
_publ_Section_title
;
 The Crystal Structures of Sodalite-Group Minerals
;

_aflow_title 'Sodalite [Na$_{4}$(AlSiO$_{4}$)$_{3}$Cl, $S6_{2}$]
    ↪ Structure'
_aflow_proto 'A3BC4D12E3_cP46_218_d_a_e_i_c'
_aflow_params 'a,x_{4},x_{5},y_{5},z_{5}'
_aflow_params_values '8.882,0.1778,0.139,0.1494,0.4383'
_aflow_Strukturbericht '$S6_{2}$'
_aflow_Pearson 'cP46'

_symmetry_space_group_name_H-M "P -4 3 n"
_symmetry_Int_Tables_number 218

_cell_length_a    8.88200
_cell_length_b    8.88200
_cell_length_c    8.88200
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 y,z,x
6 y,-z,-x
7 -y,z,-x
8 -y,-z,x
9 z,x,y
10 z,-x,-y
11 -z,x,-y
12 -z,-x,y
13 y+1/2,x+1/2,z+1/2
14 y+1/2,-x+1/2,-z+1/2
15 -y+1/2,x+1/2,-z+1/2
16 -y+1/2,-x+1/2,z+1/2
17 x+1/2,z+1/2,y+1/2
18 x+1/2,-z+1/2,-y+1/2
19 -x+1/2,z+1/2,-y+1/2
20 -x+1/2,-z+1/2,y+1/2
21 z+1/2,y+1/2,x+1/2
22 z+1/2,-y+1/2,-x+1/2
23 -z+1/2,y+1/2,-x+1/2
24 -z+1/2,-y+1/2,x+1/2

loop_
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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cl1 Cl 2 a 0.00000 0.00000 0.00000 1.00000
Si1 Si 6 c 0.25000 0.50000 0.00000 1.00000
Al1 Al 6 d 0.25000 0.00000 0.50000 1.00000
Na1 Na 8 e 0.17780 0.17780 0.17780 1.00000
O1 O 24 i 0.13900 0.14940 0.43830 1.00000
```

Sodalite $[Na_4(AlSiO_4)_3Cl, S6_2]$: A3BC4D12E3_cP46_218_d_a_e_i_c - POSCAR

```
A3BC4D12E3_cP46_218_d_a_e_i_c & a,x4,x5,y5,z5 --params=8.882,0.1778,
    ↪ 0.139,0.1494,0.4383 & P-43n T_{d}^{4} #218 (acdei) & cP46 &
    ↪ $S6_{2}$ & Al3ClNa4O12Si3 & Sodalite & I. Hassan and H. D.
    ↪ Grundy, Acta Crystallogr. Sect. B Struct. Sci. 40, 6-13 (1984)
   1.00000000000000
   8.88200000000000   0.00000000000000   0.00000000000000
   0.00000000000000   8.88200000000000   0.00000000000000
   0.00000000000000   0.00000000000000   8.88200000000000
    Al  Cl  Na  O  Si
     6   2   8 24   6
Direct
   0.25000000000000   0.00000000000000   0.50000000000000  Al  (6d)
   0.75000000000000   0.00000000000000   0.50000000000000  Al  (6d)
   0.50000000000000   0.25000000000000   0.00000000000000  Al  (6d)
   0.50000000000000   0.75000000000000   0.00000000000000  Al  (6d)
   0.00000000000000   0.50000000000000   0.25000000000000  Al  (6d)
```

```
   0.00000000000000   0.50000000000000   0.75000000000000   Al   (6d)
   0.00000000000000   0.00000000000000   0.00000000000000   Cl   (2a)
   0.50000000000000   0.50000000000000   0.50000000000000   Cl   (2a)
   0.17780000000000   0.17780000000000   0.17780000000000   Na   (8e)
  -0.17780000000000  -0.17780000000000   0.17780000000000   Na   (8e)
  -0.17780000000000   0.17780000000000  -0.17780000000000   Na   (8e)
   0.17780000000000  -0.17780000000000  -0.17780000000000   Na   (8e)
   0.67780000000000   0.67780000000000   0.67780000000000   Na   (8e)
   0.32220000000000   0.32220000000000   0.67780000000000   Na   (8e)
   0.67780000000000   0.32220000000000   0.32220000000000   Na   (8e)
   0.32220000000000   0.67780000000000   0.32220000000000   Na   (8e)
   0.13900000000000   0.14940000000000   0.43830000000000   O   (24i)
  -0.13900000000000  -0.14940000000000   0.43830000000000   O   (24i)
  -0.13900000000000   0.14940000000000  -0.43830000000000   O   (24i)
   0.13900000000000  -0.14940000000000  -0.43830000000000   O   (24i)
   0.43830000000000   0.13900000000000   0.14940000000000   O   (24i)
   0.43830000000000  -0.13900000000000  -0.14940000000000   O   (24i)
  -0.43830000000000  -0.13900000000000   0.14940000000000   O   (24i)
  -0.43830000000000   0.13900000000000  -0.14940000000000   O   (24i)
   0.14940000000000   0.43830000000000   0.13900000000000   O   (24i)
  -0.14940000000000   0.43830000000000  -0.13900000000000   O   (24i)
   0.14940000000000  -0.43830000000000  -0.13900000000000   O   (24i)
  -0.14940000000000  -0.43830000000000   0.13900000000000   O   (24i)
   0.64940000000000   0.63900000000000   0.93830000000000   O   (24i)
   0.35060000000000   0.36100000000000   0.93830000000000   O   (24i)
   0.64940000000000   0.36100000000000   0.06170000000000   O   (24i)
   0.35060000000000   0.63900000000000   0.06170000000000   O   (24i)
   0.63900000000000   0.93830000000000   0.64940000000000   O   (24i)
   0.36100000000000   0.93830000000000   0.35060000000000   O   (24i)
   0.36100000000000   0.06170000000000   0.64940000000000   O   (24i)
   0.63900000000000   0.06170000000000   0.35060000000000   O   (24i)
   0.93830000000000   0.64940000000000   0.63900000000000   O   (24i)
   0.93830000000000   0.35060000000000   0.36100000000000   O   (24i)
   0.06170000000000   0.64940000000000   0.36100000000000   O   (24i)
   0.06170000000000   0.35060000000000   0.63900000000000   O   (24i)
   0.25000000000000   0.50000000000000   0.00000000000000   Si   (6c)
   0.75000000000000   0.50000000000000   0.00000000000000   Si   (6c)
   0.00000000000000   0.25000000000000   0.50000000000000   Si   (6c)
   0.00000000000000   0.75000000000000   0.50000000000000   Si   (6c)
   0.50000000000000   0.00000000000000   0.25000000000000   Si   (6c)
   0.50000000000000   0.00000000000000   0.75000000000000   Si   (6c)
```

Eulytine ($Bi_4(SiO_4)_3$, $S1_5$): A4B12C3_cI76_220_c_e_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Eulytine'
_chemical_formula_sum 'Bi4 O12 Si3'

loop_
_publ_author_name
 'H. Liu'
 'C. Kuo'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 212
_journal_year 1997
_journal_page_first 48
_journal_page_last 48
_publ_Section_title
;
 Crystal structure of bismuth(III) silicate, Bi$_{4}$(SiO$_{4}$)$_{3}$
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'Eulytine (Bi$_{4}$(SiO$_{4}$)$_{3}$, $S1_{5}$) Structure'
_aflow_proto 'A4B12C3_cI76_220_c_e_a'
_aflow_params 'a,x_{2},x_{3},y_{3},z_{3}'
_aflow_params_values '10.2867,0.0849,0.059,0.131,0.2889'
_aflow_Strukturbericht '$S1_{5}$'
_aflow_Pearson 'cI76'

_symmetry_space_group_name_H-M "I -4 3 d"
_symmetry_Int_Tables_number 220

_cell_length_a    10.28670
_cell_length_b    10.28670
_cell_length_c    10.28670
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

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_space_group_symop_id
_space_group_symop_operation_xyz
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2 x,-y,-z+1/2
3 -x+1/2,y,-z
4 -x,-y+1/2,z
5 y,z,x
6 y,-z,-x+1/2
7 -y+1/2,z,-x
8 -y,-z+1/2,x
9 z,x,y
10 z,-x,-y+1/2
11 -z+1/2,x,-y
12 -z,-x+1/2,y
13 y+1/4,x+1/4,z+1/4
14 y+1/4,-x+3/4,-z+1/4
15 -y+1/4,x+1/4,-z+3/4
16 -y+3/4,-x+1/4,z+1/4
17 x+1/4,z+1/4,y+1/4
18 x+1/4,-z+3/4,-y+1/4
19 -x+1/4,z+1/4,-y+3/4
20 -x+3/4,-z+1/4,y+1/4
21 z+1/4,y+1/4,x+1/4
22 z+1/4,-y+3/4,-x+1/4
23 -z+1/4,y+1/4,-x+3/4
24 -z+3/4,-y+1/4,x+1/4
25 x+1/2,y+1/2,z+1/2
26 x+1/2,-y+1/2,-z
27 -x,y+1/2,-z+1/2
28 -x+1/2,-y,z+1/2
29 y+1/2,z+1/2,x+1/2
30 y+1/2,-z+1/2,-x
31 -y,z+1/2,-x+1/2
32 -y+1/2,-z,x+1/2
33 z+1/2,x+1/2,y+1/2
34 z+1/2,-x+1/2,-y
35 -z,x+1/2,-y+1/2
36 -z+1/2,-x,y+1/2
37 y+3/4,x+3/4,z+3/4
38 y+3/4,-x+1/4,-z+3/4
39 -y+3/4,x+3/4,-z+1/4
40 -y+1/4,-x+3/4,z+3/4
41 x+3/4,z+3/4,y+3/4
42 x+3/4,-z+1/4,-y+3/4
43 -x+3/4,z+3/4,-y+1/4
44 -x+1/4,-z+3/4,y+3/4
45 z+3/4,y+3/4,x+3/4
46 z+3/4,-y+1/4,-x+3/4
47 -z+3/4,y+3/4,-x+1/4
48 -z+1/4,-y+3/4,x+3/4

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_atom_site_fract_x
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Bi1 Bi 16 c 0.08490 0.08490 0.08490 1.00000
O1 O 48 e 0.05900 0.13100 0.28890 1.00000
```

Eulytine ($Bi_4(SiO_4)_3$, $S1_5$): A4B12C3_cI76_220_c_e_a - POSCAR

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    ↪ 0.131,0.2889 & I-43d T_{d}^{6} #220 (ace) & cI76 & $S1_{5}$ &
    ↪ Bi4O12Si3 & Eulytine & H. Liu and C. Kuo, Zeitschrift f{"u}r
    ↪ Kristallographie - Crystalline Materials 212, 48(1997)
   1.00000000000000
  -5.14335000000000   5.14335000000000   5.14335000000000
   5.14335000000000  -5.14335000000000   5.14335000000000
   5.14335000000000   5.14335000000000  -5.14335000000000
    Bi    O    Si
     8   24    6
Direct
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   0.00000000000000   0.33020000000000   0.50000000000000   Bi   (16c)
   0.33020000000000   0.50000000000000   0.00000000000000   Bi   (16c)
   0.66980000000000   0.66980000000000   0.66980000000000   Bi   (16c)
   0.50000000000000   0.00000000000000  -0.16980000000000   Bi   (16c)
  -0.16980000000000   0.50000000000000   0.00000000000000   Bi   (16c)
   0.00000000000000  -0.16980000000000   0.50000000000000   Bi   (16c)
   0.41990000000000   0.34790000000000   0.19000000000000    O   (48e)
   0.65790000000000   0.22990000000000   0.31000000000000    O   (48e)
  -0.15790000000000   0.15210000000000   0.57200000000000    O   (48e)
   0.08010000000000   0.27010000000000  -0.07200000000000    O   (48e)
   0.19000000000000   0.41990000000000   0.34790000000000    O   (48e)
   0.31000000000000   0.65790000000000   0.22990000000000    O   (48e)
   0.57200000000000  -0.15790000000000   0.15210000000000    O   (48e)
  -0.07200000000000   0.08010000000000   0.27010000000000    O   (48e)
   0.34790000000000   0.19000000000000   0.41990000000000    O   (48e)
   0.22990000000000   0.31000000000000   0.65790000000000    O   (48e)
   0.15210000000000   0.57200000000000  -0.15790000000000    O   (48e)
   0.27010000000000  -0.07200000000000   0.08010000000000    O   (48e)
   0.84790000000000   0.91990000000000   0.69000000000000    O   (48e)
   0.72990000000000   0.15790000000000  -0.19000000000000    O   (48e)
  -0.34790000000000   0.34210000000000   0.07200000000000    O   (48e)
  -0.22990000000000  -0.41990000000000   0.42800000000000    O   (48e)
   0.91990000000000   0.69000000000000   0.84790000000000    O   (48e)
   0.15790000000000  -0.19000000000000   0.72990000000000    O   (48e)
   0.34210000000000   0.07200000000000  -0.34790000000000    O   (48e)
  -0.41990000000000   0.42800000000000  -0.22990000000000    O   (48e)
   0.69000000000000   0.84790000000000   0.91990000000000    O   (48e)
  -0.19000000000000   0.72990000000000   0.15790000000000    O   (48e)
   0.07200000000000  -0.34790000000000   0.34210000000000    O   (48e)
   0.42800000000000  -0.22990000000000  -0.41990000000000    O   (48e)
   0.25000000000000   0.62500000000000   0.37500000000000   Si   (12a)
   0.75000000000000   0.87500000000000   0.12500000000000   Si   (12a)
   0.37500000000000   0.25000000000000   0.62500000000000   Si   (12a)
   0.12500000000000   0.75000000000000   0.87500000000000   Si   (12a)
   0.62500000000000   0.37500000000000   0.25000000000000   Si   (12a)
   0.87500000000000   0.12500000000000   0.75000000000000   Si   (12a)
```

Mayenite ($12CaO{\cdot}7Al_2O_3$, $K7_4$, C12A7): A7B12C19_cI152_220_bc_2d_ace - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM
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_chemical_name_mineral 'Mayenite'
_chemical_formula_sum 'Al7 Ca12 O19'

loop_
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 'H. Boysen'
 'M. Lerch'
 'A. Stys'
 'A. Senyshyn'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 63
_journal_year 2007
_journal_page_first 675
_journal_page_last 682
_publ_Section_title
;
 Structure and oxygen mobility in mayenite (Ca$_{12}$Al$_{14}$O$_{33}$):
      ↪ a high-temperature neutron powder diffraction study
;

# Found in The American Mineralogist Crystal Structure Database, 2003

_aflow_title 'Mayenite (12CaO\cdot$7Al$_{2}$O$_{3}$, $K7_{4}$, C12A7)
      ↪ Structure'
_aflow_proto 'A7B12C19_cI152_220_bc_2d_ace'
_aflow_params 'a,x_{3},x_{4},x_{5},x_{6},x_{7},y_{7},z_{7}'
_aflow_params_values '11.9794,0.0188,0.43519,0.1432,0.1867,0.78672,
      ↪ 0.09946,0.30708'
_aflow_Strukturbericht '$K7_{4}$'
_aflow_Pearson 'cI152'

_symmetry_space_group_name_H-M "I -4 3 d"
_symmetry_Int_Tables_number 220

_cell_length_a    11.97940
_cell_length_b    11.97940
_cell_length_c    11.97940
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
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2 x,-y,-z+1/2
3 -x+1/2,y,-z
4 -x,-y+1/2,z
5 y,z,x
6 y,-z,-x+1/2
7 -y+1/2,z,-x
8 -y,-z+1/2,x
9 z,x,y
10 z,-x,-y+1/2
11 -z+1/2,x,-y
12 -z,-x+1/2,y
13 y+1/4,x+1/4,z+1/4
14 y+1/4,-x+3/4,-z+1/4
15 -y+1/4,x+1/4,-z+3/4
16 -y+3/4,-x+1/4,z+1/4
17 x+1/4,z+1/4,y+1/4
18 x+1/4,-z+3/4,-y+1/4
19 -x+1/4,z+1/4,-y+3/4
20 -x+3/4,-z+1/4,y+1/4
21 z+1/4,y+1/4,x+1/4
22 z+1/4,-y+3/4,-x+1/4
23 -z+1/4,y+1/4,-x+3/4
24 -z+3/4,-y+1/4,x+1/4
25 x+1/2,y+1/2,z+1/2
26 x+1/2,-y+1/2,-z
27 -x,y+1/2,-z+1/2
28 -x+1/2,-y,z+1/2
29 y+1/2,z+1/2,x+1/2
30 y+1/2,-z+1/2,-x
31 -y,z+1/2,-x+1/2
32 -y+1/2,-z,x+1/2
33 z+1/2,x+1/2,y+1/2
34 z+1/2,-x+1/2,-y
35 -z,x+1/2,-y+1/2
36 -z+1/2,-x,y+1/2
37 y+3/4,x+3/4,z+3/4
38 y+3/4,-x+1/4,-z+3/4
39 -y+3/4,x+3/4,-z+1/4
40 -y+1/4,-x+3/4,z+3/4
41 x+3/4,z+3/4,y+3/4
42 x+3/4,-z+1/4,-y+3/4
43 -x+3/4,z+3/4,-y+1/4
44 -x+1/4,-z+3/4,y+3/4
45 z+3/4,y+3/4,x+3/4
46 z+3/4,-y+1/4,-x+3/4
47 -z+3/4,y+3/4,-x+1/4
48 -z+1/4,-y+3/4,x+3/4

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_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
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Al1 Al 12 b 0.87500 0.00000 0.25000 1.00000
Al2 Al 16 c 0.01880 0.01880 0.01880 1.00000
O2 O 16 c 0.43519 0.43519 0.43519 1.00000
Ca1 Ca 24 d 0.14320 0.00000 0.25000 0.87500
Ca2 Ca 24 d 0.18670 0.00000 0.25000 0.12500
O3 O 48 e 0.78672 0.09946 0.30708 1.00000
```

Mayenite (12CaO·7Al_2O_3, $K7_4$, C12A7): A7B12C19_cI152_220_bc_2d_ace - POSCAR

```
A7B12C19_cI152_220_bc_2d_ace & a,x3,x4,x5,x6,x7,y7,z7 --params=11.9794,
      ↪ 0.0188,0.43519,0.1432,0.1867,0.78672,0.09946,0.30708 & I-43d T_
      ↪ {d}^{6} #220 (abc^2d^2e) & cI152 & $K7_{4}$ & Al14Ca12O33 &
      ↪ Mayenite & H. Boysen et al., Acta Crystallogr. Sect. B Struct.
      ↪ Sci. 63, 675-682 (2007)
   1.00000000000000
  -5.98970000000000    5.98970000000000    5.98970000000000
   5.98970000000000   -5.98970000000000    5.98970000000000
   5.98970000000000    5.98970000000000   -5.98970000000000
    Al     Ca     O
    14     24     38
Direct
   0.25000000000000    0.12500000000000    0.87500000000000   Al   (12b)
   0.75000000000000    0.37500000000000    0.62500000000000   Al   (12b)
   0.87500000000000    0.25000000000000    0.12500000000000   Al   (12b)
   0.62500000000000    0.75000000000000    0.37500000000000   Al   (12b)
   0.12500000000000    0.87500000000000    0.25000000000000   Al   (12b)
   0.37500000000000    0.62500000000000    0.75000000000000   Al   (12b)
   0.03760000000000    0.03760000000000    0.03760000000000   Al   (16c)
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   0.46240000000000    0.50000000000000    0.00000000000000   Al   (16c)
   0.53760000000000    0.53760000000000    0.53760000000000   Al   (16c)
   0.50000000000000    0.00000000000000   -0.03760000000000   Al   (16c)
  -0.03760000000000    0.50000000000000    0.00000000000000   Al   (16c)
   0.00000000000000   -0.03760000000000    0.50000000000000   Al   (16c)
   0.25000000000000    0.39320000000000    0.14320000000000   Ca   (24d)
   0.75000000000000    0.10680000000000    0.35680000000000   Ca   (24d)
   0.14320000000000    0.25000000000000    0.39320000000000   Ca   (24d)
   0.35680000000000    0.75000000000000    0.10680000000000   Ca   (24d)
   0.39320000000000    0.14320000000000    0.25000000000000   Ca   (24d)
   0.10680000000000    0.35680000000000    0.75000000000000   Ca   (24d)
   0.89320000000000    0.75000000000000    0.64320000000000   Ca   (24d)
   0.60680000000000    0.25000000000000   -0.14320000000000   Ca   (24d)
   0.75000000000000    0.64320000000000    0.89320000000000   Ca   (24d)
   0.25000000000000   -0.14320000000000    0.60680000000000   Ca   (24d)
   0.64320000000000    0.89320000000000    0.75000000000000   Ca   (24d)
  -0.14320000000000    0.60680000000000    0.25000000000000   Ca   (24d)
   0.25000000000000    0.43670000000000    0.18670000000000   Ca   (24d)
   0.75000000000000    0.06330000000000    0.31330000000000   Ca   (24d)
   0.18670000000000    0.25000000000000    0.43670000000000   Ca   (24d)
   0.31330000000000    0.75000000000000    0.06330000000000   Ca   (24d)
   0.43670000000000    0.18670000000000    0.25000000000000   Ca   (24d)
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   0.93670000000000    0.75000000000000    0.68670000000000   Ca   (24d)
   0.56330000000000    0.25000000000000   -0.18670000000000   Ca   (24d)
   0.75000000000000    0.68670000000000    0.93670000000000   Ca   (24d)
   0.25000000000000   -0.18670000000000    0.56330000000000   Ca   (24d)
   0.68670000000000    0.93670000000000    0.75000000000000   Ca   (24d)
  -0.18670000000000    0.56330000000000    0.25000000000000   Ca   (24d)
   0.25000000000000    0.62500000000000    0.37500000000000   O    (12a)
   0.75000000000000    0.87500000000000    0.12500000000000   O    (12a)
   0.37500000000000    0.25000000000000    0.62500000000000   O    (12a)
   0.12500000000000    0.75000000000000    0.87500000000000   O    (12a)
   0.62500000000000    0.37500000000000    0.25000000000000   O    (12a)
   0.87500000000000    0.12500000000000    0.75000000000000   O    (12a)
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   0.50000000000000    0.00000000000000   -0.37038000000000   O    (16c)
   0.00000000000000   -0.37038000000000    0.50000000000000   O    (16c)
  -0.37038000000000    0.50000000000000    0.00000000000000   O    (16c)
   1.37038000000000    1.37038000000000    1.37038000000000   O    (16c)
   0.50000000000000    0.00000000000000   -0.87038000000000   O    (16c)
  -0.87038000000000    0.50000000000000    0.00000000000000   O    (16c)
   0.00000000000000   -0.87038000000000    0.50000000000000   O    (16c)
   0.40654000000000    1.09380000000000    0.88618000000000   O    (48e)
   0.70762000000000   -0.47964000000000   -0.38618000000000   O    (48e)
  -0.20762000000000   -0.59380000000000   -0.18726000000000   O    (48e)
   0.09346000000000    0.97964000000000    0.68726000000000   O    (48e)
   0.88618000000000    0.40654000000000    1.09380000000000   O    (48e)
  -0.38618000000000    0.70762000000000   -0.47964000000000   O    (48e)
  -0.18726000000000   -0.20762000000000   -0.59380000000000   O    (48e)
   0.68726000000000    0.09346000000000    0.97964000000000   O    (48e)
   1.09380000000000    0.88618000000000    0.40654000000000   O    (48e)
  -0.47964000000000   -0.38618000000000    0.70762000000000   O    (48e)
  -0.59380000000000   -0.18726000000000   -0.20762000000000   O    (48e)
   0.97964000000000    0.68726000000000    0.09346000000000   O    (48e)
   1.59380000000000    0.90654000000000    1.38618000000000   O    (48e)
   0.02036000000000    0.20762000000000   -0.88618000000000   O    (48e)
  -1.09380000000000    0.29238000000000   -0.68726000000000   O    (48e)
   0.47964000000000   -0.40654000000000    1.18726000000000   O    (48e)
   0.90654000000000    1.38618000000000    1.59380000000000   O    (48e)
   0.20762000000000   -0.88618000000000    0.02036000000000   O    (48e)
   0.29238000000000   -0.68726000000000   -1.09380000000000   O    (48e)
  -0.40654000000000    1.18726000000000    0.47964000000000   O    (48e)
   1.38618000000000    1.59380000000000    0.90654000000000   O    (48e)
  -0.88618000000000    0.02036000000000    0.20762000000000   O    (48e)
  -0.68726000000000   -1.09380000000000    0.29238000000000   O    (48e)
   1.18726000000000    0.47964000000000   -0.40654000000000   O    (48e)
```

$Al(PO_3)_3$ ($G5_2$): AB9C3_cI208_220_c_3e_e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM
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_chemical_formula_sum 'Al O9 P3'

loop_
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 'L. Pauling'
 'J. Sherman'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 96
_journal_year 1937
_journal_page_first 481
_journal_page_last 487
_publ_Section_title
;
 The Crystal Structure of Aluminum Metaphosphate, Al(PO$_{3}$)$_{3}$
;

# Found in The crystal structure of a monoclinic form of aluminium
      ↪ metaphosphate, Al(PO$_{3}$)$_{3}$, 1976

_aflow_title 'Al(PO$_{3}$)$_{3}$ ($G5_{2}$) Structure'
_aflow_proto 'AB9C3_cI208_220_c_3e_e'
_aflow_params 'a,x_{1},x_{2},y_{2},z_{2},x_{3},y_{3},z_{3},x_{4},y_{4},
      ↪ z_{4},x_{5},y_{5},z_{5}'
_aflow_params_values '13.63,0.117,0.09,0.11,0.8,0.095,0.141,0.245,0.137,
      ↪ 0.096,-0.014,0.34,0.063,0.124'
_aflow_Strukturbericht '$G5_{2}$'
_aflow_Pearson 'cI208'

_symmetry_space_group_name_H-M "I -4 3 d"
_symmetry_Int_Tables_number 220

_cell_length_a     13.63000
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_cell_length_c     13.63000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
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2 x,-y,-z+1/2
3 -x+1/2,y,-z
4 -x,-y+1/2,z
5 y,z,x
6 y,-z,-x+1/2
7 -y+1/2,z,-x
8 -y,-z+1/2,x
9 z,x,y
10 z,-x,-y+1/2
11 -z+1/2,x,-y
12 -z,-x+1/2,y
13 y+1/4,x+1/4,z+1/4
14 y+1/4,-x+3/4,-z+1/4
15 -y+1/4,x+1/4,-z+3/4
16 -y+3/4,-x+1/4,z+1/4
17 x+1/4,z+1/4,y+1/4
18 x+1/4,-z+3/4,-y+1/4
19 -x+1/4,z+1/4,-y+3/4
20 -x+3/4,-z+1/4,y+1/4
21 z+1/4,y+1/4,x+1/4
22 z+1/4,-y+3/4,-x+1/4
23 -z+1/4,y+1/4,-x+3/4
24 -z+3/4,-y+1/4,x+1/4
25 x+1/2,y+1/2,z+1/2
26 x+1/2,-y+1/2,-z
27 -x,y+1/2,-z+1/2
28 -x+1/2,-y,z+1/2
29 y+1/2,z+1/2,x+1/2
30 y+1/2,-z+1/2,-x
31 -y,z+1/2,-x+1/2
32 -y+1/2,-z,x+1/2
33 z+1/2,x+1/2,y+1/2
34 z+1/2,-x+1/2,-y
35 -z,x+1/2,-y+1/2
36 -z+1/2,-x,y+1/2
37 y+3/4,x+3/4,z+3/4
38 y+3/4,-x+1/4,-z+3/4
39 -y+3/4,x+3/4,-z+1/4
40 -y+1/4,-x+3/4,z+3/4
41 x+3/4,z+3/4,y+3/4
42 x+3/4,-z+1/4,-y+3/4
43 -x+3/4,z+3/4,-y+1/4
44 -x+1/4,-z+3/4,y+3/4
45 z+3/4,y+3/4,x+3/4
46 z+3/4,-y+1/4,-x+3/4
47 -z+3/4,y+3/4,-x+1/4
48 -z+1/4,-y+3/4,x+3/4

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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Al1 Al 16 c 0.11700 0.11700 0.11700 1.00000
O1 O 48 e 0.09000 0.11000 0.80000 1.00000
O2 O 48 e 0.09500 0.14100 0.24500 1.00000
O3 O 48 e 0.13700 0.09600 -0.01400 1.00000
P1 P 48 e 0.34000 0.06300 0.12400 1.00000
```

Al(PO$_3$)$_3$ ($G5_2$): AB9C3_cI208_220_c_3e_e - POSCAR

```
AB9C3_cI208_220_c_3e_e & a,x1,x2,y2,z2,x3,y3,z3,x4,y4,z4,x5,y5,z5 --
      ↪ params=13.63,0.117,0.09,0.11,0.8,0.095,0.141,0.245,0.137,0.096
      ↪ ,-0.014,0.34,0.063,0.124 & I-43d T_{d}^{6} #220 (ce^4) & cI208
      ↪ & $G5_{2}$ & AlO9P3 & AlO9P3 & L. Pauling and J. Sherman,
      ↪ Zeitschrift f{"u}r Kristallographie - Crystalline Materials 96,
      ↪ 481-487 (1937)
   1.00000000000000
  -6.81500000000000    6.81500000000000    6.81500000000000
   6.81500000000000   -6.81500000000000    6.81500000000000
   6.81500000000000    6.81500000000000   -6.81500000000000
    Al     O     P
     8    72    24
Direct
   0.23400000000000    0.23400000000000    0.23400000000000   Al  (16c)
   0.50000000000000    0.00000000000000    0.26600000000000   Al  (16c)
   0.00000000000000    0.26600000000000    0.50000000000000   Al  (16c)
   0.26600000000000    0.50000000000000    0.00000000000000   Al  (16c)
   0.73400000000000    0.73400000000000    0.73400000000000   Al  (16c)
   0.50000000000000    0.00000000000000   -0.23400000000000   Al  (16c)
  -0.23400000000000    0.50000000000000    0.00000000000000   Al  (16c)
   0.00000000000000   -0.23400000000000    0.50000000000000   Al  (16c)
   0.91000000000000    0.89000000000000    0.20000000000000   O  (48e)
   1.19000000000000    0.71000000000000    0.30000000000000   O  (48e)
  -0.69000000000000   -0.39000000000000    0.52000000000000   O  (48e)
  -0.41000000000000   -0.21000000000000   -0.02000000000000   O  (48e)
   0.20000000000000    0.91000000000000    0.89000000000000   O  (48e)
   0.30000000000000    1.19000000000000    0.71000000000000   O  (48e)
   0.52000000000000   -0.69000000000000   -0.39000000000000   O  (48e)
  -0.02000000000000   -0.41000000000000   -0.21000000000000   O  (48e)
   0.89000000000000    0.20000000000000    0.91000000000000   O  (48e)
   0.71000000000000    0.30000000000000    1.19000000000000   O  (48e)
  -0.39000000000000    0.52000000000000   -0.69000000000000   O  (48e)
  -0.21000000000000   -0.02000000000000   -0.41000000000000   O  (48e)
   1.39000000000000    1.41000000000000    0.70000000000000   O  (48e)
   1.21000000000000    0.69000000000000   -0.20000000000000   O  (48e)
  -0.89000000000000   -0.19000000000000    0.02000000000000   O  (48e)
  -0.71000000000000   -0.91000000000000    0.48000000000000   O  (48e)
   1.41000000000000    0.70000000000000    1.39000000000000   O  (48e)
   0.69000000000000   -0.20000000000000    1.21000000000000   O  (48e)
  -0.19000000000000    0.02000000000000   -0.89000000000000   O  (48e)
  -0.91000000000000    0.48000000000000   -0.71000000000000   O  (48e)
   0.70000000000000    1.39000000000000    1.41000000000000   O  (48e)
  -0.20000000000000    1.21000000000000    0.69000000000000   O  (48e)
   0.02000000000000   -0.89000000000000   -0.19000000000000   O  (48e)
   0.48000000000000   -0.71000000000000   -0.91000000000000   O  (48e)
   0.38600000000000    0.34000000000000    0.23600000000000   O  (48e)
   0.60400000000000    0.15000000000000    0.26400000000000   O  (48e)
  -0.10400000000000    0.16000000000000    0.54600000000000   O  (48e)
   0.11400000000000    0.35000000000000   -0.04600000000000   O  (48e)
   0.23600000000000    0.38600000000000    0.34000000000000   O  (48e)
   0.26400000000000    0.60400000000000    0.15000000000000   O  (48e)
   0.54600000000000   -0.10400000000000    0.16000000000000   O  (48e)
  -0.04600000000000    0.11400000000000    0.35000000000000   O  (48e)
   0.34000000000000    0.23600000000000    0.38600000000000   O  (48e)
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   0.16000000000000    0.54600000000000   -0.10400000000000   O  (48e)
   0.35000000000000   -0.04600000000000    0.11400000000000   O  (48e)
   0.84000000000000    0.88600000000000    0.73600000000000   O  (48e)
   0.65000000000000    0.10400000000000   -0.23600000000000   O  (48e)
  -0.34000000000000    0.39600000000000    0.04600000000000   O  (48e)
  -0.15000000000000   -0.38600000000000    0.45400000000000   O  (48e)
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   0.10400000000000   -0.23600000000000    0.65000000000000   O  (48e)
   0.39600000000000    0.04600000000000   -0.34000000000000   O  (48e)
  -0.38600000000000    0.45400000000000   -0.15000000000000   O  (48e)
   0.73600000000000    0.84000000000000    0.88600000000000   O  (48e)
  -0.23600000000000    0.65000000000000    0.10400000000000   O  (48e)
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   0.11000000000000    0.37700000000000    0.45900000000000   O  (48e)
   0.41800000000000    0.65100000000000    0.04100000000000   O  (48e)
   0.23300000000000    0.08200000000000    0.12300000000000   O  (48e)
   0.26700000000000    0.39000000000000   -0.15100000000000   O  (48e)
   0.45900000000000    0.11000000000000    0.37700000000000   O  (48e)
   0.04100000000000    0.41800000000000    0.65100000000000   O  (48e)
   0.12300000000000    0.23300000000000    0.08200000000000   O  (48e)
  -0.15100000000000    0.26700000000000    0.39000000000000   O  (48e)
   0.37700000000000    0.45900000000000    0.11000000000000   O  (48e)
   0.65100000000000    0.04100000000000    0.41800000000000   O  (48e)
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  -0.12300000000000    0.61000000000000   -0.04100000000000   O  (48e)
   0.15100000000000   -0.08200000000000    0.54100000000000   O  (48e)
   0.58200000000000    0.73300000000000    0.62300000000000   O  (48e)
  -0.11000000000000   -0.23300000000000    0.34900000000000   O  (48e)
   0.61000000000000   -0.04100000000000   -0.12300000000000   O  (48e)
  -0.08200000000000    0.54100000000000    0.15100000000000   O  (48e)
   0.73300000000000    0.62300000000000    0.58200000000000   O  (48e)
  -0.23300000000000    0.34900000000000   -0.11000000000000   O  (48e)
  -0.04100000000000   -0.12300000000000    0.61000000000000   O  (48e)
   0.54100000000000    0.15100000000000   -0.08200000000000   O  (48e)
   0.18700000000000    0.46400000000000    0.40300000000000   P  (48e)
   0.56100000000000   -0.21600000000000    0.09700000000000   P  (48e)
  -0.06100000000000    0.03600000000000    0.22300000000000   P  (48e)
   0.31300000000000    0.71600000000000    0.27700000000000   P  (48e)
   0.40300000000000    0.18700000000000    0.46400000000000   P  (48e)
   0.09700000000000    0.56100000000000   -0.21600000000000   P  (48e)
```

```
   0.22300000000000  -0.06100000000000   0.03600000000000   P  (48e)
   0.27700000000000   0.31300000000000   0.71600000000000   P  (48e)
   0.46400000000000   0.40300000000000   0.18700000000000   P  (48e)
  -0.21600000000000   0.09700000000000   0.56100000000000   P  (48e)
   0.03600000000000   0.22300000000000  -0.06100000000000   P  (48e)
   0.71600000000000   0.27700000000000   0.31300000000000   P  (48e)
   0.96400000000000   0.68700000000000   0.90300000000000   P  (48e)
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  -0.46400000000000   0.43900000000000  -0.27700000000000   P  (48e)
   0.21600000000000  -0.18700000000000   0.77700000000000   P  (48e)
   0.68700000000000   0.90300000000000   0.96400000000000   P  (48e)
   0.06100000000000  -0.40300000000000   0.28400000000000   P  (48e)
   0.43900000000000  -0.27700000000000  -0.46400000000000   P  (48e)
  -0.18700000000000   0.77700000000000   0.21600000000000   P  (48e)
   0.90300000000000   0.96400000000000   0.68700000000000   P  (48e)
  -0.40300000000000   0.28400000000000   0.06100000000000   P  (48e)
  -0.27700000000000  -0.46400000000000   0.43900000000000   P  (48e)
   0.77700000000000   0.21600000000000  -0.18700000000000   P  (48e)
```

AlN (cI24): AB_cI24_220_a_b - CIF

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_audit_creation_method FINDSYM

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_chemical_formula_sum 'Al N'

loop_
_publ_author_name
 'C. Liu'
 'M. Chen'
 'J. Li'
 'L. Liu'
 'P. Li'
 'M. Ma'
 'C. Shao'
 'J. He'
 'T. Liang'
_journal_name_full_name
;
 Journal of Physics and Chemistry of Solids
;
_journal_volume 130
_journal_year 2019
_journal_page_first 58
_journal_page_last 66
_publ_Section_title
;
 A first-principles study of novel cubic AlN phases
;

_aflow_title 'AlN (cI24) Structure'
_aflow_proto 'AB_cI24_220_a_b'
_aflow_params 'a'
_aflow_params_values '6.2951'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'cI24'

_symmetry_space_group_name_H-M "I -4 3 d"
_symmetry_Int_Tables_number 220

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_cell_length_c    6.29510
_cell_angle_alpha 90.00000
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_cell_angle_gamma 90.00000

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_space_group_symop_id
_space_group_symop_operation_xyz
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3 -x+1/2,y,-z
4 -x,-y+1/2,z
5 y,z,x
6 y,-z,-x+1/2
7 -y+1/2,z,-x
8 -y,-z+1/2,x
9 z,x,y
10 z,-x,-y+1/2
11 -z+1/2,x,-y
12 -z,-x+1/2,y
13 y+1/4,x+1/4,z+1/4
14 y+1/4,-x+3/4,-z+1/4
15 -y+1/4,x+1/4,-z+3/4
16 -y+3/4,-x+1/4,z+1/4
17 x+1/4,z+1/4,y+1/4
18 x+1/4,-z+3/4,-y+1/4
19 -x+1/4,z+1/4,-y+3/4
20 -x+3/4,-z+1/4,y+1/4
21 z+1/4,y+1/4,x+1/4
22 z+1/4,-y+3/4,-x+1/4
23 -z+1/4,y+1/4,-x+3/4
24 -z+3/4,-y+1/4,x+1/4
25 x+1/2,y+1/2,z+1/2
26 x+1/2,-y+1/2,-z
27 -x,y+1/2,-z+1/2
28 -x+1/2,-y,z+1/2
29 y+1/2,z+1/2,x+1/2
30 y+1/2,-z+1/2,-x
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33 z+1/2,x+1/2,y+1/2
34 z+1/2,-x+1/2,-y
35 -z,x+1/2,-y+1/2
36 -z+1/2,-x,y+1/2
37 y+3/4,x+3/4,z+3/4
38 y+3/4,-x+1/4,-z+3/4
39 -y+3/4,x+3/4,-z+1/4
40 -y+1/4,-x+3/4,z+3/4
41 x+3/4,z+3/4,y+3/4
42 x+3/4,-z+1/4,-y+3/4
43 -x+3/4,z+3/4,-y+1/4
44 -x+1/4,-z+3/4,y+3/4
45 z+3/4,y+3/4,x+3/4
46 z+3/4,-y+1/4,-x+3/4
47 -z+3/4,y+3/4,-x+1/4
48 -z+1/4,-y+3/4,x+3/4

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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
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N1 N 12 b 0.87500 0.00000 0.25000 1.00000
```

AlN (cI24): AB_cI24_220_a_b - POSCAR

```
AB_cI24_220_a_b & a --params=6.2951 & I-43d T_{d}^{6} #220 (ab) & cI24 &
    ↪ None & AlN & AlN & C. Liu et al., J. Phys. Chem. Solids 130,
    ↪ 58-66 (2019)
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  -3.14755000000000   3.14755000000000   3.14755000000000
   3.14755000000000  -3.14755000000000   3.14755000000000
   3.14755000000000   3.14755000000000  -3.14755000000000
    Al    N
     6     6
Direct
   0.25000000000000   0.62500000000000   0.37500000000000   Al  (12a)
   0.75000000000000   0.87500000000000   0.12500000000000   Al  (12a)
   0.37500000000000   0.25000000000000   0.62500000000000   Al  (12a)
   0.12500000000000   0.75000000000000   0.87500000000000   Al  (12a)
   0.62500000000000   0.37500000000000   0.25000000000000   Al  (12a)
   0.87500000000000   0.12500000000000   0.75000000000000   Al  (12a)
   0.25000000000000   0.12500000000000   0.87500000000000   N  (12b)
   0.75000000000000   0.37500000000000   0.62500000000000   N  (12b)
   0.87500000000000   0.25000000000000   0.12500000000000   N  (12b)
   0.62500000000000   0.75000000000000   0.37500000000000   N  (12b)
   0.12500000000000   0.87500000000000   0.25000000000000   N  (12b)
   0.37500000000000   0.62500000000000   0.75000000000000   N  (12b)
```

γ-Fe_4N ($L'1_0$): A4B_cP5_221_bc_a - CIF

```
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_audit_creation_method FINDSYM

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_chemical_formula_sum 'Fe4 N'

loop_
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 'H. Jacobs'
 'R. Rechenbach'
 'U. Zachwieja'
_journal_name_full_name
;
 Journal of Alloys and Compounds
;
_journal_volume 227
_journal_year 1995
_journal_page_first 10
_journal_page_last 17
_publ_Section_title
;
 Structure determination of $\gamma\'$-Fe$_{4}$N and $\epsilon$-Fe$_{3}
    ↪ $N
;

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_aflow_params 'a'
_aflow_params_values '3.79'
_aflow_Strukturbericht '$L'1_{0}$'
_aflow_Pearson 'cP5'

_symmetry_space_group_name_H-M "P 4/m -3 2/m"
_symmetry_Int_Tables_number 221

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_cell_length_b    3.79000
_cell_length_c    3.79000
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
```

```
5 y,z,x
6 y,-z,-x
7 -y,z,-x
8 -y,-z,x
9 z,x,y
10 z,-x,-y
11 -z,x,-y
12 -z,-x,y
13 -y,-x,-z
14 -y,x,z
15 y,-x,z
16 y,x,-z
17 -x,-z,-y
18 -x,z,y
19 x,-z,y
20 x,z,-y
21 -z,-y,-x
22 -z,y,x
23 z,-y,x
24 z,y,-x
25 -x,-y,-z
26 -x,y,z
27 x,-y,z
28 x,y,-z
29 -y,-z,-x
30 -y,z,x
31 y,-z,x
32 y,z,-x
33 -z,-x,-y
34 -z,x,y
35 z,-x,y
36 z,x,-y
37 y,x,z
38 y,-x,-z
39 -y,x,-z
40 -y,-x,z
41 x,z,y
42 x,-z,-y
43 -x,z,-y
44 -x,-z,y
45 z,y,x
46 z,-y,-x
47 -z,y,-x
48 -z,-y,x

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_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
N1 N 1 a 0.00000 0.00000 0.00000 1.00000
Fe1 Fe 1 b 0.50000 0.50000 0.50000 1.00000
Fe2 Fe 3 c 0.00000 0.50000 0.50000 1.00000
```

γ-Fe_4N ($L'1_0$): A4B_cP5_221_bc_a - POSCAR

```
A4B_cP5_221_bc_a & a --params=3.79 & Pm-3m O_{h}^{1} #221 (abc) & cP5 &
    ↪ $L'1_{0}$ & Fe4N & Fe4N & H. Jacobs and R. Rechenbach and U.
    ↪ Zachwieja, J. Alloys Compd. 227, 10-17 (1995)
   1.00000000000000
   3.79000000000000    0.00000000000000    0.00000000000000
   0.00000000000000    3.79000000000000    0.00000000000000
   0.00000000000000    0.00000000000000    3.79000000000000
    Fe     N
     4     1
Direct
   0.50000000000000    0.50000000000000    0.50000000000000   Fe    (1b)
   0.00000000000000    0.50000000000000    0.50000000000000   Fe    (3c)
   0.50000000000000    0.00000000000000    0.50000000000000   Fe    (3c)
   0.50000000000000    0.50000000000000    0.00000000000000   Fe    (3c)
   0.00000000000000    0.00000000000000    0.00000000000000    N    (1a)
```

Predicted High-Pressure $YCaH_{12}$: AB12C_cP14_221_a_h_b - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'CaH12Y'
_chemical_formula_sum 'Ca H12 Y'

loop_
_publ_author_name
 'H. Xie'
 'D. Duan'
 'Z. Shao'
 'H. Song'
 'Y. Wang'
 'X. Xiao'
 'D. Li'
 'F. Tian'
 'B. Liu'
 'T. Cui'
_journal_name_full_name
;
 Journal of Physics: Condensed Matter
;
_journal_volume 31
_journal_year 2019
_journal_page_first 245404
_journal_page_last 245404
_publ_Section_title
;
 High-temperature superconductivity in ternary clathrate YCaH$_{12}$
     ↪ under high pressures
;

_aflow_title 'Predicted High-Pressure YCaH$_{12}$ Structure'
_aflow_proto 'AB12C_cP14_221_a_h_b'
_aflow_params 'a,x_{3}'
_aflow_params_values '3.4708,0.24126'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'cP14'

_symmetry_space_group_name_H-M "P 4/m -3 2/m"
_symmetry_Int_Tables_number 221

_cell_length_a      3.47080
_cell_length_b      3.47080
_cell_length_c      3.47080
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 y,z,x
6 y,-z,-x
7 -y,z,-x
8 -y,-z,x
9 z,x,y
10 z,-x,-y
11 -z,x,-y
12 -z,-x,y
13 -y,-x,-z
14 -y,x,z
15 y,-x,z
16 y,x,-z
17 -x,-z,-y
18 -x,z,y
19 x,-z,y
20 x,z,-y
21 -z,-y,-x
22 -z,y,x
23 z,-y,x
24 z,y,-x
25 -x,-y,-z
26 -x,y,z
27 x,-y,z
28 x,y,-z
29 -y,-z,-x
30 -y,z,x
31 y,-z,x
32 y,z,-x
33 -z,-x,-y
34 -z,x,y
35 z,-x,y
36 z,x,-y
37 y,x,z
38 y,-x,-z
39 -y,x,-z
40 -y,-x,z
41 x,z,y
42 x,-z,-y
43 -x,z,-y
44 -x,-z,y
45 z,y,x
46 z,-y,-x
47 -z,y,-x
48 -z,-y,x

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Ca1 Ca 1 a 0.00000 0.00000 0.00000 1.00000
Y1 Y 1 b 0.50000 0.50000 0.50000 1.00000
H1 H 12 h 0.24126 0.50000 0.00000 1.00000
```

Predicted High-Pressure $YCaH_{12}$: AB12C_cP14_221_a_h_b - POSCAR

```
AB12C_cP14_221_a_h_b & a,x3 --params=3.4708,0.24126 & Pm-3m O_{h}^{1} #
    ↪ 221 (abh) & cP14 & None & CaH12Y & CaH12Y & H. Xie et al., J.
    ↪ Phys.: Condens. Matter 31, 245404(2019)
   1.00000000000000
   3.47080000000000    0.00000000000000    0.00000000000000
   0.00000000000000    3.47080000000000    0.00000000000000
   0.00000000000000    0.00000000000000    3.47080000000000
    Ca     H     Y
     1    12     1
Direct
   0.00000000000000    0.00000000000000    0.00000000000000   Ca    (1a)
   0.24126000000000    0.50000000000000    0.00000000000000    H   (12h)
  -0.24126000000000    0.50000000000000    0.00000000000000    H   (12h)
   0.00000000000000    0.24126000000000    0.50000000000000    H   (12h)
   0.00000000000000   -0.24126000000000    0.50000000000000    H   (12h)
```

```
   0.50000000000000   0.00000000000000   0.24126000000000   H   (12h)
   0.50000000000000   0.00000000000000  -0.24126000000000   H   (12h)
   0.50000000000000   0.24126000000000   0.00000000000000   H   (12h)
   0.50000000000000  -0.24126000000000   0.00000000000000   H   (12h)
   0.24126000000000   0.00000000000000   0.50000000000000   H   (12h)
  -0.24126000000000   0.00000000000000   0.50000000000000   H   (12h)
   0.00000000000000   0.50000000000000  -0.24126000000000   H   (12h)
   0.00000000000000   0.50000000000000   0.24126000000000   H   (12h)
   0.50000000000000   0.50000000000000   0.50000000000000   Y   (1b)
```

NH_4NO_3 I ($G0_8$): AB_cP2_221_a_b - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'NH4NO3'
_chemical_formula_sum '(NH4) (NO3)'

loop_
_publ_author_name
 'F. C. Kracek'
 'S. B. Hendricks'
 'E. Posnjak'
_journal_name_full_name
;
 Nature
;
_journal_volume 128
_journal_year 1931
_journal_page_first 410
_journal_page_last 411
_publ_Section_title
;
 Group Rotation in Solid Ammonium and Calcium Nitrates
;

# Found in Strukturbericht Band II 1928-1932, 1937

_aflow_title 'NH$_{4}$NO$_{3}$ I ($G0_{8}$) Structure'
_aflow_proto 'AB_cP2_221_a_b'
_aflow_params 'a'
_aflow_params_values '4.4'
_aflow_Strukturbericht '$G0_{8}$'
_aflow_Pearson 'cP2'

_symmetry_space_group_name_H-M "P 4/m -3 2/m"
_symmetry_Int_Tables_number 221

_cell_length_a     4.40000
_cell_length_b     4.40000
_cell_length_c     4.40000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 y,z,x
6 y,-z,-x
7 -y,z,-x
8 -y,-z,x
9 z,x,y
10 z,-x,-y
11 -z,x,-y
12 -z,-x,y
13 -y,-x,-z
14 -y,x,z
15 y,-x,z
16 y,x,-z
17 -x,-z,-y
18 -x,z,y
19 x,-z,y
20 x,z,-y
21 -z,-y,-x
22 -z,y,x
23 z,-y,x
24 z,y,-x
25 -x,-y,-z
26 -x,y,z
27 x,-y,z
28 x,y,-z
29 -y,-z,-x
30 -y,z,x
31 y,-z,x
32 y,z,-x
33 -z,-x,-y
34 -z,x,y
35 z,-x,y
36 z,x,-y
37 y,x,z
38 y,-x,-z
39 -y,x,-z
40 -y,-x,z
41 x,z,y
42 x,-z,-y
43 -x,z,-y
44 -x,-z,y
45 z,y,x
46 z,-y,-x
47 -z,y,-x
48 -z,-y,x

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
NH41 NH4 1 a 0.00000 0.00000 0.00000 1.00000
NO31 NO3 1 b 0.50000 0.50000 0.50000 1.00000
```

NH_4NO_3 I ($G0_8$): AB_cP2_221_a_b - POSCAR

```
AB_cP2_221_a_b & a --params=4.4 & Pm-3m O_{h}^{1} #221 (ab) & cP2 & $G0_
    ↪ {8}$ & NH4NO3 & NH4NO3 & F. C. Kracek and S. B. Hendricks and
    ↪ E. Posnjak, Nature 128, 410-411 (1931)
   1.00000000000000
   4.40000000000000   0.00000000000000   0.00000000000000
   0.00000000000000   4.40000000000000   0.00000000000000
   0.00000000000000   0.00000000000000   4.40000000000000
   NH4    NO3
     1      1
Direct
   0.00000000000000   0.00000000000000   0.00000000000000  NH4    (1a)
   0.50000000000000   0.50000000000000   0.50000000000000  NO3    (1b)
```

Dodecatungstophosphoric Acid Hexahydrate $[H_3PW_{12}O_{40}\cdot 6H_2O]$: A27B52CD12_cP184_224_dl_eh3k_a_k - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Dodecatungstophosphoric acid hexahydrate'
_chemical_formula_sum 'H27 O52 P W12'

loop_
_publ_author_name
 'G. M. Brown'
 'M.-R. {Noe-Spirlet}'
 'W. R. Busing'
 'H. A. Levy'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 33
_journal_year 1977
_journal_page_first 1038
_journal_page_last 1046
_publ_Section_title
;
 Dodecatungstophosphoric acid hexahydrate, (H$_{5}$O$_{2}$$^{+}$)$_{3}$(
    ↪ PW$_{12}$O$_{40}$$^{3-}$). The true structure of Keggin\'s '
    ↪ pentahydrate\' from single-crystal X-ray and neutron
    ↪ diffraction data
;

_aflow_title 'Dodecatungstophosphoric Acid Hexahydrate [H$_{3}$PW$_{12}
    ↪ $O$_{40}$$\cdot$6H$_{2}$O] Structure'
_aflow_proto 'A27B52CD12_cP184_224_dl_eh3k_a_k'
_aflow_params 'a,x_{3},x_{4},x_{5},z_{5},x_{6},z_{6},x_{7},z_{7},x_{8},
    ↪ z_{8},x_{9},y_{9},z_{9}'
_aflow_params_values '12.506,0.32273,0.34752,0.65683,-0.00684,0.37155,
    ↪ 0.52447,0.55441,0.23276,0.45688,0.25822,0.24525,0.61259,0.68288
    ↪ '
_aflow_Strukturbericht 'None'
_aflow_Pearson 'cP184'

_symmetry_space_group_name_H-M "P 42/n -3 2/m (origin choice 2)"
_symmetry_Int_Tables_number 224

_cell_length_a     12.50600
_cell_length_b     12.50600
_cell_length_c     12.50600
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y+1/2,-z+1/2
3 -x+1/2,y,-z+1/2
4 -x+1/2,-y+1/2,z
5 y,z,x
6 y,-z+1/2,-x+1/2
7 -y+1/2,z,-x+1/2
8 -y+1/2,-z+1/2,x
9 z,x,y
10 z,-x+1/2,-y+1/2
11 -z+1/2,x,-y+1/2
12 -z+1/2,-x+1/2,y
13 -y,-x,-z
14 -y,x+1/2,z+1/2
15 y+1/2,-x,z+1/2
16 y+1/2,x+1/2,-z
17 -x,-z,-y
18 -x,z+1/2,y+1/2
19 x+1/2,-z,y+1/2
```

```
20 x+1/2,z+1/2,-y
21 -z,-y,-x
22 -z,y+1/2,x+1/2
23 z+1/2,-y,x+1/2
24 z+1/2,y+1/2,-x
25 -x,-y,-z
26 -x,y+1/2,z+1/2
27 x+1/2,-y,z+1/2
28 x+1/2,y+1/2,-z
29 -y,-z,-x
30 -y,z+1/2,x+1/2
31 y+1/2,-z,x+1/2
32 y+1/2,z+1/2,-x
33 -z,-x,-y
34 -z,x+1/2,y+1/2
35 z+1/2,-x,y+1/2
36 z+1/2,x+1/2,-y
37 y,x,z
38 y,-x+1/2,-z+1/2
39 -y+1/2,x,-z+1/2
40 -y+1/2,-x+1/2,z
41 x,z,y
42 x,-z+1/2,-y+1/2
43 -x+1/2,z,-y+1/2
44 -x+1/2,-z+1/2,y
45 z,y,x
46 z,-y+1/2,-x+1/2
47 -z+1/2,y,-x+1/2
48 -z+1/2,-y+1/2,x

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
P1 P 2 a 0.25000 0.25000 0.25000 1.00000
H1 H 6 d 0.25000 0.75000 0.75000 1.00000
O1 O 8 e 0.32273 0.32273 0.32273 1.00000
O2 O 24 h 0.34752 0.25000 0.75000 0.50000
O3 O 24 k 0.65683 0.65683 -0.00684 1.00000
O4 O 24 k 0.37155 0.37155 0.52447 1.00000
O5 O 24 k 0.55441 0.55441 0.23276 1.00000
W1 W 24 k 0.45688 0.45688 0.25822 1.00000
H2 H 48 l 0.24525 0.61259 0.68288 0.50000
```

Dodecatungstophosphoric Acid Hexahydrate $[H_3PW_{12}O_{40}{\cdot}6H_2O]$: A27B52CD12_cP184_224_dl_eh3k_a_k - POSCAR

```
A27B52CD12_cP184_224_dl_eh3k_a_k & a,x3,x4,x5,z5,x6,z6,x7,z7,x8,z8,x9,y9
    ↪ ,z9 --params=12.506,0.32273,0.34752,0.65683,-0.00684,0.37155,
    ↪ 0.52447,0.55441,0.23276,0.45688,0.25822,0.24525,0.61259,0.68288
    ↪ & Pn-3m O_{h}^{4} #224 (adehk^41) & cP184 & None & H15O46PW12
    ↪ & Dodecatungstophosphoric acid hexahydrate & G. M. Brown et al.
    ↪ , Acta Crystallogr. Sect. B Struct. Sci. 33, 1038-1046 (1977)
   1.00000000000000
  12.50600000000000   0.00000000000000   0.00000000000000
   0.00000000000000  12.50600000000000   0.00000000000000
   0.00000000000000   0.00000000000000  12.50600000000000
     H     O     P     W
    54   104     2    24
Direct
   0.25000000000000   0.75000000000000   0.75000000000000   H   (6d)
   0.75000000000000   0.25000000000000   0.75000000000000   H   (6d)
   0.75000000000000   0.75000000000000   0.25000000000000   H   (6d)
   0.25000000000000   0.75000000000000   0.25000000000000   H   (6d)
   0.75000000000000   0.25000000000000   0.25000000000000   H   (6d)
   0.25000000000000   0.25000000000000   0.75000000000000   H   (6d)
   0.24525000000000   0.61259000000000   0.68288000000000   H  (48l)
   0.25475000000000  -0.11259000000000   0.68288000000000   H  (48l)
   0.25475000000000   0.61259000000000  -0.18288000000000   H  (48l)
   0.24525000000000  -0.11259000000000  -0.18288000000000   H  (48l)
   0.68288000000000   0.24525000000000   0.61259000000000   H  (48l)
   0.68288000000000   0.25475000000000  -0.11259000000000   H  (48l)
  -0.18288000000000   0.25475000000000   0.61259000000000   H  (48l)
  -0.18288000000000   0.24525000000000  -0.11259000000000   H  (48l)
   0.61259000000000   0.68288000000000   0.24525000000000   H  (48l)
  -0.11259000000000   0.68288000000000   0.25475000000000   H  (48l)
   0.61259000000000  -0.18288000000000   0.25475000000000   H  (48l)
  -0.11259000000000  -0.18288000000000   0.24525000000000   H  (48l)
   1.11259000000000   0.74525000000000  -0.68288000000000   H  (48l)
  -0.61259000000000  -0.24525000000000  -0.68288000000000   H  (48l)
   1.11259000000000  -0.24525000000000   1.18288000000000   H  (48l)
  -0.61259000000000   0.74525000000000   1.18288000000000   H  (48l)
   0.74525000000000   1.18288000000000  -0.61259000000000   H  (48l)
  -0.24525000000000   1.18288000000000   1.11259000000000   H  (48l)
  -0.24525000000000  -0.68288000000000  -0.61259000000000   H  (48l)
   0.74525000000000  -0.68288000000000   1.11259000000000   H  (48l)
   1.18288000000000   1.11259000000000  -0.24525000000000   H  (48l)
   1.18288000000000  -0.61259000000000   0.74525000000000   H  (48l)
  -0.68288000000000   1.11259000000000   0.74525000000000   H  (48l)
  -0.68288000000000  -0.61259000000000  -0.24525000000000   H  (48l)
  -0.24525000000000  -0.61259000000000  -0.68288000000000   H  (48l)
   0.74525000000000   1.11259000000000  -0.68288000000000   H  (48l)
   0.74525000000000  -0.61259000000000   1.18288000000000   H  (48l)
  -0.24525000000000   1.11259000000000   1.18288000000000   H  (48l)
  -0.68288000000000  -0.24525000000000  -0.61259000000000   H  (48l)
  -0.68288000000000   0.74525000000000   1.11259000000000   H  (48l)
   1.18288000000000   0.74525000000000  -0.61259000000000   H  (48l)
   1.18288000000000  -0.24525000000000   1.11259000000000   H  (48l)
  -0.61259000000000  -0.68288000000000  -0.24525000000000   H  (48l)
   1.11259000000000  -0.68288000000000   0.74525000000000   H  (48l)
  -0.61259000000000   1.18288000000000   0.74525000000000   H  (48l)
   1.11259000000000   1.18288000000000  -0.24525000000000   H  (48l)
  -0.11259000000000   0.25475000000000   0.68288000000000   H  (48l)
   0.61259000000000   0.24525000000000   0.68288000000000   H  (48l)
  -0.11259000000000   0.24525000000000  -0.18288000000000   H  (48l)
   0.61259000000000   0.25475000000000  -0.18288000000000   H  (48l)
   0.25475000000000  -0.18288000000000   0.61259000000000   H  (48l)
   0.24525000000000  -0.18288000000000  -0.11259000000000   H  (48l)
   0.24525000000000   0.68288000000000   0.61259000000000   H  (48l)
   0.25475000000000   0.68288000000000  -0.11259000000000   H  (48l)
  -0.18288000000000  -0.11259000000000   0.24525000000000   H  (48l)
  -0.18288000000000   0.61259000000000   0.25475000000000   H  (48l)
   0.68288000000000  -0.11259000000000   0.25475000000000   H  (48l)
   0.68288000000000   0.61259000000000   0.24525000000000   H  (48l)
   0.32273000000000   0.32273000000000   0.32273000000000   O   (8e)
   0.17727000000000   0.17727000000000   0.32273000000000   O   (8e)
   0.17727000000000   0.32273000000000   0.17727000000000   O   (8e)
   0.32273000000000   0.17727000000000   0.17727000000000   O   (8e)
   0.82273000000000   0.82273000000000  -0.32273000000000   O   (8e)
  -0.32273000000000  -0.32273000000000  -0.32273000000000   O   (8e)
   0.82273000000000  -0.32273000000000   0.82273000000000   O   (8e)
  -0.32273000000000   0.82273000000000   0.82273000000000   O   (8e)
   0.34752000000000   0.25000000000000   0.75000000000000   O  (24h)
   0.15248000000000   0.25000000000000   0.75000000000000   O  (24h)
   0.75000000000000   0.34752000000000   0.25000000000000   O  (24h)
   0.75000000000000   0.15248000000000   0.25000000000000   O  (24h)
   0.25000000000000   0.75000000000000   0.34752000000000   O  (24h)
   0.25000000000000   0.75000000000000   0.15248000000000   O  (24h)
   0.75000000000000   0.84752000000000   0.25000000000000   O  (24h)
   0.75000000000000  -0.34752000000000   0.25000000000000   O  (24h)
   0.84752000000000   0.25000000000000   0.75000000000000   O  (24h)
  -0.34752000000000   0.25000000000000   0.75000000000000   O  (24h)
   0.25000000000000   0.75000000000000  -0.34752000000000   O  (24h)
   0.25000000000000   0.75000000000000   0.84752000000000   O  (24h)
  -0.34752000000000   0.75000000000000   0.25000000000000   O  (24h)
   0.84752000000000   0.75000000000000   0.25000000000000   O  (24h)
   0.25000000000000  -0.34752000000000   0.75000000000000   O  (24h)
   0.25000000000000   0.84752000000000   0.75000000000000   O  (24h)
   0.75000000000000   0.25000000000000  -0.34752000000000   O  (24h)
   0.75000000000000   0.25000000000000   0.84752000000000   O  (24h)
   0.25000000000000   0.15248000000000   0.75000000000000   O  (24h)
   0.25000000000000   0.34752000000000   0.75000000000000   O  (24h)
   0.15248000000000   0.75000000000000   0.25000000000000   O  (24h)
   0.34752000000000   0.75000000000000   0.25000000000000   O  (24h)
   0.75000000000000   0.25000000000000   0.34752000000000   O  (24h)
   0.75000000000000   0.25000000000000   0.15248000000000   O  (24h)
   0.65683000000000   0.65683000000000  -0.00684000000000   O  (24k)
  -0.15683000000000  -0.15683000000000  -0.00684000000000   O  (24k)
  -0.15683000000000   0.65683000000000   0.50684000000000   O  (24k)
   0.65683000000000  -0.15683000000000   0.50684000000000   O  (24k)
  -0.00684000000000   0.65683000000000   0.65683000000000   O  (24k)
  -0.00684000000000  -0.15683000000000  -0.15683000000000   O  (24k)
   0.50684000000000  -0.15683000000000   0.65683000000000   O  (24k)
   0.50684000000000   0.65683000000000  -0.15683000000000   O  (24k)
   0.65683000000000  -0.00684000000000   0.65683000000000   O  (24k)
  -0.15683000000000  -0.00684000000000  -0.15683000000000   O  (24k)
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  -0.15683000000000   0.50684000000000   0.65683000000000   O  (24k)
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  -0.65683000000000  -0.65683000000000   0.00684000000000   O  (24k)
   1.15683000000000  -0.65683000000000   0.49316000000000   O  (24k)
  -0.65683000000000   1.15683000000000   0.49316000000000   O  (24k)
   1.15683000000000   0.49316000000000  -0.65683000000000   O  (24k)
  -0.65683000000000   0.49316000000000   1.15683000000000   O  (24k)
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   1.15683000000000   0.00684000000000   1.15683000000000   O  (24k)
   0.49316000000000   1.15683000000000  -0.65683000000000   O  (24k)
   0.49316000000000  -0.65683000000000   1.15683000000000   O  (24k)
   0.00684000000000   1.15683000000000   1.15683000000000   O  (24k)
   0.00684000000000  -0.65683000000000  -0.65683000000000   O  (24k)
   0.37155000000000   0.37155000000000   0.52447000000000   O  (24k)
   0.12845000000000   0.12845000000000   0.52447000000000   O  (24k)
   0.12845000000000   0.37155000000000  -0.02447000000000   O  (24k)
   0.37155000000000   0.12845000000000  -0.02447000000000   O  (24k)
   0.52447000000000   0.37155000000000   0.37155000000000   O  (24k)
   0.52447000000000   0.12845000000000   0.12845000000000   O  (24k)
  -0.02447000000000   0.12845000000000   0.37155000000000   O  (24k)
  -0.02447000000000   0.37155000000000   0.12845000000000   O  (24k)
   0.37155000000000   0.52447000000000   0.37155000000000   O  (24k)
   0.12845000000000   0.52447000000000   0.12845000000000   O  (24k)
   0.37155000000000  -0.02447000000000   0.12845000000000   O  (24k)
   0.12845000000000  -0.02447000000000   0.37155000000000   O  (24k)
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  -0.37155000000000  -0.37155000000000  -0.52447000000000   O  (24k)
   0.87155000000000  -0.37155000000000   1.02447000000000   O  (24k)
  -0.37155000000000   0.87155000000000   1.02447000000000   O  (24k)
   0.87155000000000   1.02447000000000  -0.37155000000000   O  (24k)
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  -0.37155000000000  -0.52447000000000  -0.37155000000000   O  (24k)
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   1.02447000000000   0.87155000000000  -0.37155000000000   O  (24k)
   1.02447000000000  -0.37155000000000   0.87155000000000   O  (24k)
  -0.52447000000000   0.87155000000000   0.87155000000000   O  (24k)
  -0.52447000000000  -0.37155000000000  -0.37155000000000   O  (24k)
   0.55441000000000   0.55441000000000   0.23276000000000   O  (24k)
  -0.05441000000000  -0.05441000000000   0.23276000000000   O  (24k)
  -0.05441000000000   0.55441000000000   0.26724000000000   O  (24k)
   0.55441000000000  -0.05441000000000   0.26724000000000   O  (24k)
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   0.26724000000000   0.55441000000000  -0.05441000000000   O  (24k)
   0.55441000000000   0.23276000000000   0.55441000000000   O  (24k)
  -0.05441000000000   0.23276000000000  -0.05441000000000   O  (24k)
   0.55441000000000   0.26724000000000  -0.05441000000000   O  (24k)
```

−0.05441000000000	0.26724000000000	0.55441000000000	O	(24k)
1.05441000000000	1.05441000000000	−0.23276000000000	O	(24k)
−0.55441000000000	−0.55441000000000	−0.23276000000000	O	(24k)
1.05441000000000	−0.55441000000000	0.73276000000000	O	(24k)
−0.55441000000000	1.05441000000000	0.73276000000000	O	(24k)
1.05441000000000	0.73276000000000	−0.55441000000000	O	(24k)
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−0.55441000000000	−0.23276000000000	−0.55441000000000	O	(24k)
1.05441000000000	−0.23276000000000	1.05441000000000	O	(24k)
0.73276000000000	1.05441000000000	−0.55441000000000	O	(24k)
0.73276000000000	−0.55441000000000	1.05441000000000	O	(24k)
−0.23276000000000	1.05441000000000	1.05441000000000	O	(24k)
−0.23276000000000	−0.55441000000000	−0.55441000000000	O	(24k)
0.25000000000000	0.25000000000000	0.25000000000000	P	(2a)
0.75000000000000	0.75000000000000	0.75000000000000	P	(2a)
0.45688000000000	0.45688000000000	0.25822000000000	W	(24k)
0.04312000000000	0.04312000000000	0.25822000000000	W	(24k)
0.04312000000000	0.45688000000000	0.24178000000000	W	(24k)
0.45688000000000	0.04312000000000	0.24178000000000	W	(24k)
0.25822000000000	0.45688000000000	0.45688000000000	W	(24k)
0.25822000000000	0.04312000000000	0.04312000000000	W	(24k)
0.24178000000000	0.04312000000000	0.45688000000000	W	(24k)
0.24178000000000	0.45688000000000	0.04312000000000	W	(24k)
0.45688000000000	0.25822000000000	0.45688000000000	W	(24k)
0.04312000000000	0.25822000000000	0.04312000000000	W	(24k)
0.45688000000000	0.24178000000000	0.04312000000000	W	(24k)
0.04312000000000	0.24178000000000	0.45688000000000	W	(24k)
0.95688000000000	0.95688000000000	−0.25822000000000	W	(24k)
−0.45688000000000	−0.45688000000000	−0.25822000000000	W	(24k)
0.95688000000000	−0.45688000000000	0.75822000000000	W	(24k)
−0.45688000000000	0.95688000000000	0.75822000000000	W	(24k)
0.95688000000000	0.75822000000000	−0.45688000000000	W	(24k)
−0.45688000000000	0.75822000000000	0.95688000000000	W	(24k)
−0.45688000000000	−0.25822000000000	−0.45688000000000	W	(24k)
0.95688000000000	−0.25822000000000	0.95688000000000	W	(24k)
0.75822000000000	0.95688000000000	−0.45688000000000	W	(24k)
0.75822000000000	−0.45688000000000	0.95688000000000	W	(24k)
−0.25822000000000	0.95688000000000	0.95688000000000	W	(24k)
−0.25822000000000	−0.45688000000000	−0.45688000000000	W	(24k)

Mg_3P_2 ($D5_5$): A3B2_cP10_224_d_b - CIF

```
# CIF file
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_audit_creation_method FINDSYM

_chemical_name_mineral 'Mg3P2'
_chemical_formula_sum 'Mg3 P2'

loop_
_publ_author_name
 'L. Passerini'
_journal_name_full_name
;
 Gazzetta Chimica Italiana
;
_journal_volume 58
_journal_year 1928
_journal_page_first 655
_journal_page_last 664
_publ_Section_title
;
 Struttura cristallina di alcuni fosfuri di metalli bivalenti e
      ↪ trivalenti
;

# Found in Strukturbericht Band II 1928-1932, 1937

_aflow_title 'Mg$_{3}$P$_{2}$ ($D5_{5}$) Structure'
_aflow_proto 'A3B2_cP10_224_d_b'
_aflow_params 'a'
_aflow_params_values '5.91207'
_aflow_Strukturbericht '$D5_{5}$'
_aflow_Pearson 'cP10'

_symmetry_space_group_name_H-M "P 42/n -3 2/m (origin choice 2)"
_symmetry_Int_Tables_number 224

_cell_length_a    5.91207
_cell_length_b    5.91207
_cell_length_c    5.91207
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

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_space_group_symop_operation_xyz
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3 -x+1/2,y,-z+1/2
4 -x+1/2,-y+1/2,z
5 y,z,x
6 y,-z+1/2,-x+1/2
7 -y+1/2,z,-x+1/2
8 -y+1/2,-z+1/2,x
9 z,x,y
10 z,-x+1/2,-y+1/2
11 -z+1/2,x,-y+1/2
12 -z+1/2,-x+1/2,y
13 -y,-x,-z
14 -y,x+1/2,z+1/2
15 y+1/2,-x,z+1/2
16 y+1/2,x+1/2,-z
17 -x,-z,-y
18 -x,z+1/2,y+1/2
19 x+1/2,-z,y+1/2
20 x+1/2,z+1/2,-y
21 -z,-y,-x
22 -z,y+1/2,x+1/2
23 z+1/2,-y,x+1/2
24 z+1/2,y+1/2,-x
25 -x,-y,-z
26 -x,y+1/2,z+1/2
27 x+1/2,-y,z+1/2
28 x+1/2,y+1/2,-z
29 -y,-z,-x
30 -y,z+1/2,x+1/2
31 y+1/2,-z,x+1/2
32 y+1/2,z+1/2,-x
33 -z,-x,-y
34 -z,x+1/2,y+1/2
35 z+1/2,-x,y+1/2
36 z+1/2,x+1/2,-y
37 y,x,z
38 y,-x+1/2,-z+1/2
39 -y+1/2,x,-z+1/2
40 -y+1/2,-x+1/2,z
41 x,z,y
42 x,-z+1/2,-y+1/2
43 -x+1/2,z,-y+1/2
44 -x+1/2,-z+1/2,y
45 z,y,x
46 z,-y+1/2,-x+1/2
47 -z+1/2,y,-x+1/2
48 -z+1/2,-y+1/2,x

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_atom_site_occupancy
P1 P 4 b 0.00000 0.00000 0.00000 1.00000
Mg1 Mg 6 d 0.25000 0.75000 0.75000 1.00000
```

Mg_3P_2 ($D5_5$): A3B2_cP10_224_d_b - POSCAR

```
A3B2_cP10_224_d_b & a --params=5.91207 & Pn-3m O_{h}^{4} #224 (bd) &
      ↪ cP10 & $D5_{5}$ & Mg3P2 & Mg3P2 & L. Passerini, Gazz. Chim.
      ↪ Ital. 58, 655-664 (1928)
   1.00000000000000
   5.91207000000000    0.00000000000000    0.00000000000000
   0.00000000000000    5.91207000000000    0.00000000000000
   0.00000000000000    0.00000000000000    5.91207000000000
    Mg      P
     6      4
Direct
   0.25000000000000    0.75000000000000    0.75000000000000  Mg   (6d)
   0.75000000000000    0.25000000000000    0.75000000000000  Mg   (6d)
   0.75000000000000    0.75000000000000    0.25000000000000  Mg   (6d)
   0.25000000000000    0.75000000000000    0.25000000000000  Mg   (6d)
   0.75000000000000    0.25000000000000    0.25000000000000  Mg   (6d)
   0.25000000000000    0.25000000000000    0.75000000000000  Mg   (6d)
   0.00000000000000    0.00000000000000    0.00000000000000  P    (4b)
   0.50000000000000    0.50000000000000    0.00000000000000  P    (4b)
   0.50000000000000    0.00000000000000    0.50000000000000  P    (4b)
   0.00000000000000    0.50000000000000    0.50000000000000  P    (4b)
```

$H_3PW_{12}O_{40}\cdot 3H_2O$: A3B40CD12_cP112_224_d_e3k_a_k - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral '(H3O)3O40PW12'
_chemical_formula_sum '(H3O)3 O40 P W12'

loop_
_publ_author_name
 'L. Marosi'
 'E. E. Platero'
 'J. Cifre'
 'C. O. Are\\'{a}na'
_journal_name_full_name
;
 Journal of Materials Chemistry
;
_journal_volume 10
_journal_year 2000
_journal_page_first 1949
_journal_page_last 1955
_publ_Section_title
;
 Thermal dehydration of H$_{3+x}$PV$_{x}$M${12-x}$O$_{40}$$\cdot$y$H$_{
      ↪ 2}$O Keggin type heteropolyacids; formation, thermal stability
      ↪  and structure of the anhydrous acids H$_{3}$PM$_{12}$O$_{40}$
      ↪ , of the corresponding anhydrides PM$_{12}$O$_{38.5}$ and of a
      ↪  novel trihydrate H$_{3}$PW$_{12}$O$_{40}$$\cdot$3H$_{2}$O
;

# Found in H$_{3}$PW$_{12}$O$_{40}$$\cdot$3H$_{2}$O Crystal Structure,
      ↪ 2016 Found in H$_{3}$PW$_{12}$O$_{40}$$\cdot$3H$_{2}$O Crystal
      ↪ Structure, {in: Inorganic Solid Phases, SpringerMaterials (
      ↪ online database)},

_aflow_title 'H$_{3}$PW$_{12}$O$_{40}$$\cdot$3H$_{2}$O Structure'
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```
_aflow_proto 'A3B40CD12_cP112_224_d_e3k_a_k'
_aflow_params 'a,x_{3},x_{4},z_{4},x_{5},z_{5},x_{6},z_{6},x_{7},z_{7}'
_aflow_params_values '11.75,0.3226,0.6463,0.0133,0.3687,0.5462,0.5629,
    ↪ 0.2412,0.4636,0.2579'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'cP112'

_symmetry_space_group_name_H-M "P 42/n -3 2/m (origin choice 2)"
_symmetry_Int_Tables_number 224

_cell_length_a 11.75000
_cell_length_b 11.75000
_cell_length_c 11.75000
_cell_angle_alpha 90.00000
_cell_angle_beta 90.00000
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_space_group_symop_id
_space_group_symop_operation_xyz
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2 x,-y+1/2,-z+1/2
3 -x+1/2,y,-z+1/2
4 -x+1/2,-y+1/2,z
5 y,z,x
6 y,-z+1/2,-x+1/2
7 -y+1/2,z,-x+1/2
8 -y+1/2,-z+1/2,x
9 z,x,y
10 z,-x+1/2,-y+1/2
11 -z+1/2,x,-y+1/2
12 -z+1/2,-x+1/2,y
13 -y,-x,-z
14 -y,x+1/2,z+1/2
15 y+1/2,-x,z+1/2
16 y+1/2,x+1/2,-z
17 -x,-z,-y
18 -x,z+1/2,y+1/2
19 x+1/2,-z,y+1/2
20 x+1/2,z+1/2,-y
21 -z,-y,-x
22 -z,y+1/2,x+1/2
23 z+1/2,-y,x+1/2
24 z+1/2,y+1/2,-x
25 -x,-y,-z
26 -x,y+1/2,z+1/2
27 x+1/2,-y,z+1/2
28 x+1/2,y+1/2,-z
29 -y,-z,-x
30 -y,z+1/2,x+1/2
31 y+1/2,-z,x+1/2
32 y+1/2,z+1/2,-x
33 -z,-x,-y
34 -z,x+1/2,y+1/2
35 z+1/2,-x,y+1/2
36 z+1/2,x+1/2,-y
37 y,x,z
38 y,-x+1/2,-z+1/2
39 -y+1/2,x,-z+1/2
40 -y+1/2,-x+1/2,z
41 x,z,y
42 x,-z+1/2,-y+1/2
43 -x+1/2,z,-y+1/2
44 -x+1/2,-z+1/2,y
45 z,y,x
46 z,-y+1/2,-x+1/2
47 -z+1/2,y,-x+1/2
48 -z+1/2,-y+1/2,x

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_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
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H3O1 H3O 6 d 0.25000 0.75000 0.75000 1.00000
O1 O 8 e 0.32260 0.32260 0.32260 1.00000
O2 O 24 k 0.64630 0.64630 0.01330 1.00000
O3 O 24 k 0.36870 0.36870 0.54620 1.00000
O4 O 24 k 0.56290 0.56290 0.24120 1.00000
W1 W 24 k 0.46360 0.46360 0.25790 1.00000
```

$H_3PW_{12}O_{40}\cdot 3H_2O$: A3B40CD12_cP112_224_d_e3k_a_k - POSCAR

```
A3B40CD12_cP112_224_d_e3k_a_k & a,x3,x4,z4,x5,z5,x6,z6,x7,z7 --params=
    ↪ 11.75,0.3226,0.6463,0.0133,0.3687,0.5462,0.5629,0.2412,0.4636,
    ↪ 0.2579 & Pn-3m O_{h}^{4} #224 (adek^4) & cP112 & None & (H3O)
    ↪ 3O40PW12 & (H3O)3O40PW12 & L. Marosi et al., J. Mater. Chem. 10
    ↪ , 1949-1955 (2000)
  1.00000000000000
 11.75000000000000  0.00000000000000  0.00000000000000
  0.00000000000000 11.75000000000000  0.00000000000000
  0.00000000000000  0.00000000000000 11.75000000000000
  H3O   O   P   W
    6  80   2  24
Direct
  0.25000000000000  0.75000000000000  0.75000000000000  H3O  (6d)
  0.75000000000000  0.25000000000000  0.75000000000000  H3O  (6d)
  0.75000000000000  0.75000000000000  0.25000000000000  H3O  (6d)
  0.25000000000000  0.75000000000000  0.25000000000000  H3O  (6d)
  0.75000000000000  0.25000000000000  0.25000000000000  H3O  (6d)
  0.25000000000000  0.25000000000000  0.75000000000000  H3O  (6d)
  0.32260000000000  0.32260000000000  0.32260000000000  O  (8e)
  0.17740000000000  0.17740000000000  0.32260000000000  O  (8e)
  0.17740000000000  0.32260000000000  0.17740000000000  O  (8e)
  0.32260000000000  0.17740000000000  0.17740000000000  O  (8e)
  0.82260000000000  0.82260000000000 -0.32260000000000  O  (8e)
 -0.32260000000000 -0.32260000000000 -0.32260000000000  O  (8e)
  0.82260000000000 -0.32260000000000  0.82260000000000  O  (8e)
 -0.32260000000000  0.82260000000000  0.82260000000000  O  (8e)
  0.64630000000000  0.64630000000000  0.01330000000000  O (24k)
 -0.14630000000000 -0.14630000000000  0.01330000000000  O (24k)
 -0.14630000000000  0.64630000000000  0.48670000000000  O (24k)
  0.64630000000000 -0.14630000000000  0.48670000000000  O (24k)
  0.01330000000000  0.64630000000000  0.64630000000000  O (24k)
  0.01330000000000 -0.14630000000000 -0.14630000000000  O (24k)
  0.48670000000000 -0.14630000000000  0.64630000000000  O (24k)
  0.48670000000000  0.64630000000000 -0.14630000000000  O (24k)
  0.64630000000000  0.01330000000000  0.64630000000000  O (24k)
 -0.14630000000000  0.01330000000000 -0.14630000000000  O (24k)
  0.64630000000000  0.48670000000000 -0.14630000000000  O (24k)
 -0.14630000000000  0.48670000000000  0.64630000000000  O (24k)
  1.14630000000000  1.14630000000000 -0.01330000000000  O (24k)
 -0.64630000000000 -0.64630000000000 -0.01330000000000  O (24k)
  1.14630000000000 -0.64630000000000  0.51330000000000  O (24k)
 -0.64630000000000  1.14630000000000  0.51330000000000  O (24k)
  1.14630000000000  0.51330000000000 -0.64630000000000  O (24k)
 -0.64630000000000  0.51330000000000  1.14630000000000  O (24k)
 -0.64630000000000 -0.01330000000000 -0.64630000000000  O (24k)
  1.14630000000000 -0.01330000000000  1.14630000000000  O (24k)
  0.51330000000000  1.14630000000000 -0.64630000000000  O (24k)
  0.51330000000000 -0.64630000000000  1.14630000000000  O (24k)
 -0.01330000000000  1.14630000000000  1.14630000000000  O (24k)
 -0.01330000000000 -0.64630000000000 -0.64630000000000  O (24k)
  0.36870000000000  0.36870000000000  0.54620000000000  O (24k)
  0.13130000000000  0.13130000000000  0.54620000000000  O (24k)
  0.13130000000000  0.36870000000000 -0.04620000000000  O (24k)
  0.36870000000000  0.13130000000000 -0.04620000000000  O (24k)
  0.54620000000000  0.36870000000000  0.36870000000000  O (24k)
  0.54620000000000  0.13130000000000  0.13130000000000  O (24k)
 -0.04620000000000  0.13130000000000  0.36870000000000  O (24k)
 -0.04620000000000  0.36870000000000  0.13130000000000  O (24k)
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  0.13130000000000  0.54620000000000  0.13130000000000  O (24k)
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  0.13130000000000 -0.04620000000000  0.36870000000000  O (24k)
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 -0.36870000000000 -0.36870000000000 -0.54620000000000  O (24k)
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  0.86870000000000 -0.54620000000000  0.86870000000000  O (24k)
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  1.04620000000000 -0.36870000000000  0.86870000000000  O (24k)
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 -0.06290000000000  0.56290000000000  0.25880000000000  O (24k)
  0.56290000000000 -0.06290000000000  0.25880000000000  O (24k)
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  0.56290000000000  0.24120000000000  0.56290000000000  O (24k)
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 -0.56290000000000 -0.56290000000000 -0.24120000000000  O (24k)
  1.06290000000000 -0.56290000000000  0.74120000000000  O (24k)
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  1.06290000000000  0.74120000000000 -0.56290000000000  O (24k)
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  0.74120000000000  1.06290000000000 -0.56290000000000  O (24k)
  0.74120000000000 -0.56290000000000  1.06290000000000  O (24k)
 -0.24120000000000  1.06290000000000  1.06290000000000  O (24k)
 -0.24120000000000 -0.56290000000000 -0.56290000000000  O (24k)
  0.25000000000000  0.25000000000000  0.25000000000000  P  (2a)
  0.75000000000000  0.75000000000000  0.75000000000000  P  (2a)
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  0.03640000000000  0.03640000000000  0.25790000000000  W (24k)
  0.03640000000000  0.46360000000000  0.24210000000000  W (24k)
  0.46360000000000  0.03640000000000  0.24210000000000  W (24k)
  0.25790000000000  0.46360000000000  0.46360000000000  W (24k)
  0.25790000000000  0.03640000000000  0.03640000000000  W (24k)
  0.24210000000000  0.03640000000000  0.46360000000000  W (24k)
  0.24210000000000  0.46360000000000  0.03640000000000  W (24k)
  0.46360000000000  0.25790000000000  0.46360000000000  W (24k)
  0.03640000000000  0.25790000000000  0.03640000000000  W (24k)
  0.46360000000000  0.24210000000000  0.03640000000000  W (24k)
  0.03640000000000  0.24210000000000  0.46360000000000  W (24k)
  0.96360000000000  0.96360000000000 -0.25790000000000  W (24k)
 -0.46360000000000 -0.46360000000000 -0.25790000000000  W (24k)
  0.96360000000000 -0.46360000000000  0.75790000000000  W (24k)
 -0.46360000000000  0.96360000000000  0.75790000000000  W (24k)
  0.96360000000000  0.75790000000000 -0.46360000000000  W (24k)
 -0.46360000000000  0.75790000000000  0.96360000000000  W (24k)
 -0.46360000000000 -0.25790000000000 -0.46360000000000  W (24k)
  0.96360000000000 -0.25790000000000  0.96360000000000  W (24k)
  0.75790000000000  0.96360000000000 -0.46360000000000  W (24k)
  0.75790000000000 -0.46360000000000  0.96360000000000  W (24k)
```

```
  −0.25790000000000    0.96360000000000    0.96360000000000    W  (24k)
  −0.25790000000000   −0.46360000000000   −0.46360000000000    W  (24k)
```

12-phosphotungstic acid [$H_3PW_{12}O_{40}{\cdot}5H_2O$ ($H4_{16}$)]: A5B40CD12_cP116_224_cd_e3k_a_k - CIF

```
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_audit_creation_method FINDSYM

_chemical_name_mineral '12-phosphotungstic acid'
_chemical_formula_sum '(H2O)5 O40 P W12'

loop_
_publ_author_name
 'J. F. Keggin'
_journal_name_full_name
;
 Proceedings of the Royal Society London A
;
_journal_volume 144
_journal_year 1934
_journal_page_first 75
_journal_page_last 100
_publ_Section_title
;
 The structure and formula of 12-phosphotungstic acid
;

# Found in Dodecatungstophosphoric acid hexahydrate, (H$_{5}$O$_{2}^{+}$
      ↪ )$_{3}$(PW$_{12}$O$_{40}^{3-}$). The true structure of Keggin's
      ↪  'pentahydrate' from single-crystal X-ray and neutron
      ↪  diffraction data, 1977

_aflow_title '12-phosphotungstic acid [H$_{3}$PW$_{12}$O$_{40}$$\
      ↪ cdot$5H$_{2}$O ($H4_{16}$)] Structure'
_aflow_proto 'A5B40CD12_cP116_224_cd_e3k_a_k'
_aflow_params 'a,x_{4},x_{5},z_{5},x_{6},z_{6},x_{7},z_{7},x_{8},z_{8}'
_aflow_params_values '12.141,0.83154,0.67011,0.51608,0.87272,0.04157,
      ↪ 0.43783,0.74176,-0.0445,0.75577'
_aflow_Strukturbericht '$H4_{16}$'
_aflow_Pearson 'cP116'

_symmetry_space_group_name_H-M "P 42/n -3 2/m (origin choice 2)"
_symmetry_Int_Tables_number 224

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_cell_length_b     12.14100
_cell_length_c     12.14100
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

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_space_group_symop_id
_space_group_symop_operation_xyz
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2 x,-y+1/2,-z+1/2
3 -x+1/2,y,-z+1/2
4 -x+1/2,-y+1/2,z
5 y,z,x
6 y,-z+1/2,-x+1/2
7 -y+1/2,z,-x+1/2
8 -y+1/2,-z+1/2,x
9 z,x,y
10 z,-x+1/2,-y+1/2
11 -z+1/2,x,-y+1/2
12 -z+1/2,-x+1/2,y
13 -y,-x,-z
14 -y,x+1/2,z+1/2
15 y+1/2,-x,z+1/2
16 y+1/2,x+1/2,-z
17 -x,-z,-y
18 -x,z+1/2,y+1/2
19 x+1/2,-z,y+1/2
20 x+1/2,z+1/2,-y
21 -z,-y,-x
22 -z,y+1/2,x+1/2
23 z+1/2,-y,x+1/2
24 z+1/2,y+1/2,-x
25 -x,-y,-z
26 -x,y+1/2,z+1/2
27 x+1/2,-y,z+1/2
28 x+1/2,y+1/2,-z
29 -y,-z,-x
30 -y,z+1/2,x+1/2
31 y+1/2,-z,x+1/2
32 y+1/2,z+1/2,-x
33 -z,-x,-y
34 -z,x+1/2,y+1/2
35 z+1/2,-x,y+1/2
36 z+1/2,x+1/2,-y
37 y,x,z
38 y,-x+1/2,-z+1/2
39 -y+1/2,x,-z+1/2
40 -y+1/2,-x+1/2,z
41 x,z,y
42 x,-z+1/2,-y+1/2
43 -x+1/2,z,-y+1/2
44 -x+1/2,-z+1/2,y
45 z,y,x
46 z,-y+1/2,-x+1/2
47 -z+1/2,y,-x+1/2
48 -z+1/2,-y+1/2,x

loop_
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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
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H2O1 H2O 4 c 0.50000 0.50000 0.50000 1.00000
H2O2 H2O 6 d 0.25000 0.75000 0.75000 1.00000
O1 O 8 e 0.83154 0.83154 0.83154 1.00000
O2 O 24 k 0.67011 0.67011 0.51608 1.00000
O3 O 24 k 0.87272 0.87272 0.04157 1.00000
O4 O 24 k 0.43783 0.43783 0.74176 1.00000
W1 W 24 k -0.04450 -0.04450 0.75577 1.00000
```

12-phosphotungstic acid [$H_3PW_{12}O_{40}{\cdot}5H_2O$ ($H4_{16}$)]: A5B40CD12_cP116_224_cd_e3k_a_k - POSCAR

```
A5B40CD12_cP116_224_cd_e3k_a_k & a,x4,x5,z5,x6,z6,x7,z7,x8,z8 --params=
      ↪ 12.141,0.83154,0.67011,0.51608,0.87272,0.04157,0.43783,0.74176
      ↪ ,-0.0445,0.75577 & Pn-3m O_{h}^{4} #224 (acdek^4) & cP116 &
      ↪ $H4_{16}$ & [5H2.6O]O40PW12 & 12-phosphotungstic acid & J. F.
      ↪ Keggin, Proc. Roy. Soc. Lond. A 144, 75-100 (1934)
   1.00000000000000
  12.14100000000000    0.00000000000000    0.00000000000000
   0.00000000000000   12.14100000000000    0.00000000000000
   0.00000000000000    0.00000000000000   12.14100000000000
   H2O     O     P     W
    10    80     2    24
Direct
   0.50000000000000    0.50000000000000    0.50000000000000  H2O   (4c)
   0.00000000000000    0.00000000000000    0.50000000000000  H2O   (4c)
   0.00000000000000    0.50000000000000    0.00000000000000  H2O   (4c)
   0.50000000000000    0.00000000000000    0.00000000000000  H2O   (4c)
   0.25000000000000    0.75000000000000    0.75000000000000  H2O   (6d)
   0.75000000000000    0.25000000000000    0.75000000000000  H2O   (6d)
   0.75000000000000    0.75000000000000    0.25000000000000  H2O   (6d)
   0.25000000000000    0.75000000000000    0.25000000000000  H2O   (6d)
   0.75000000000000    0.25000000000000    0.25000000000000  H2O   (6d)
   0.25000000000000    0.25000000000000    0.75000000000000  H2O   (6d)
   0.83154000000000    0.83154000000000    0.83154000000000    O   (8e)
  -0.33154000000000   -0.33154000000000    0.83154000000000    O   (8e)
  -0.33154000000000    0.83154000000000   -0.33154000000000    O   (8e)
   0.83154000000000   -0.33154000000000   -0.33154000000000    O   (8e)
   1.33154000000000    1.33154000000000   -0.83154000000000    O   (8e)
  -0.83154000000000   -0.83154000000000   -0.83154000000000    O   (8e)
   1.33154000000000   -0.83154000000000    1.33154000000000    O   (8e)
  -0.83154000000000    1.33154000000000    1.33154000000000    O   (8e)
   0.67011000000000    0.67011000000000    0.51608000000000    O  (24k)
  -0.17011000000000   -0.17011000000000    0.51608000000000    O  (24k)
  -0.17011000000000    0.67011000000000   -0.01608000000000    O  (24k)
   0.67011000000000   -0.17011000000000   -0.01608000000000    O  (24k)
   0.51608000000000    0.67011000000000    0.67011000000000    O  (24k)
   0.51608000000000   -0.17011000000000   -0.17011000000000    O  (24k)
  -0.01608000000000   -0.17011000000000    0.67011000000000    O  (24k)
  -0.01608000000000    0.67011000000000   -0.17011000000000    O  (24k)
   0.67011000000000    0.51608000000000    0.67011000000000    O  (24k)
  -0.17011000000000    0.51608000000000   -0.17011000000000    O  (24k)
   0.67011000000000   -0.01608000000000   -0.17011000000000    O  (24k)
  -0.17011000000000   -0.01608000000000    0.67011000000000    O  (24k)
   1.17011000000000    1.17011000000000   -0.51608000000000    O  (24k)
  -0.67011000000000   -0.67011000000000   -0.51608000000000    O  (24k)
   1.17011000000000   -0.67011000000000    1.01608000000000    O  (24k)
  -0.67011000000000    1.17011000000000    1.01608000000000    O  (24k)
   1.17011000000000    1.01608000000000   -0.67011000000000    O  (24k)
  -0.67011000000000    1.01608000000000    1.17011000000000    O  (24k)
  -0.67011000000000   -0.51608000000000   -0.67011000000000    O  (24k)
   1.17011000000000   -0.51608000000000    1.17011000000000    O  (24k)
   1.01608000000000    1.17011000000000   -0.67011000000000    O  (24k)
   1.01608000000000   -0.67011000000000    1.17011000000000    O  (24k)
  -0.51608000000000    1.17011000000000    1.17011000000000    O  (24k)
  -0.51608000000000   -0.67011000000000   -0.67011000000000    O  (24k)
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  -0.37272000000000   -0.37272000000000    0.04157000000000    O  (24k)
  -0.37272000000000    0.87272000000000    0.45843000000000    O  (24k)
   0.87272000000000   -0.37272000000000    0.45843000000000    O  (24k)
   0.04157000000000    0.87272000000000    0.87272000000000    O  (24k)
   0.04157000000000   -0.37272000000000   -0.37272000000000    O  (24k)
   0.45843000000000   -0.37272000000000    0.87272000000000    O  (24k)
   0.45843000000000    0.87272000000000   -0.37272000000000    O  (24k)
   0.87272000000000    0.04157000000000    0.87272000000000    O  (24k)
  -0.37272000000000    0.04157000000000   -0.37272000000000    O  (24k)
   0.87272000000000    0.45843000000000   -0.37272000000000    O  (24k)
  -0.37272000000000    0.45843000000000    0.87272000000000    O  (24k)
   1.37272000000000    1.37272000000000   -0.04157000000000    O  (24k)
  -0.87272000000000   -0.87272000000000   -0.04157000000000    O  (24k)
   1.37272000000000   -0.87272000000000    0.54157000000000    O  (24k)
  -0.87272000000000    1.37272000000000    0.54157000000000    O  (24k)
   1.37272000000000    0.54157000000000   -0.87272000000000    O  (24k)
  -0.87272000000000    0.54157000000000    1.37272000000000    O  (24k)
  -0.87272000000000   -0.04157000000000   -0.87272000000000    O  (24k)
   1.37272000000000   -0.04157000000000    1.37272000000000    O  (24k)
   0.54157000000000    1.37272000000000   -0.87272000000000    O  (24k)
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   0.06217000000000    0.43783000000000   -0.24176000000000    O  (24k)
   0.43783000000000    0.06217000000000   -0.24176000000000    O  (24k)
   0.74176000000000    0.43783000000000    0.43783000000000    O  (24k)
   0.74176000000000    0.06217000000000    0.06217000000000    O  (24k)
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  -0.24176000000000    0.43783000000000    0.06217000000000    O  (24k)
```

```
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  0.06217000000000   0.74176000000000   0.06217000000000   O  (24k)
  0.43783000000000  -0.24176000000000   0.06217000000000   O  (24k)
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 -0.43783000000000  -0.74176000000000  -0.43783000000000   O  (24k)
  0.93783000000000  -0.74176000000000   0.93783000000000   O  (24k)
  1.24176000000000   0.93783000000000  -0.43783000000000   O  (24k)
  1.24176000000000  -0.43783000000000   0.93783000000000   O  (24k)
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 -0.74176000000000  -0.43783000000000  -0.43783000000000   O  (24k)
  0.25000000000000   0.25000000000000   0.25000000000000   P   (2a)
  0.75000000000000   0.75000000000000   0.75000000000000   P   (2a)
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  0.54450000000000   0.54450000000000   0.75577000000000   W  (24k)
  0.54450000000000  -0.04450000000000  -0.25577000000000   W  (24k)
 -0.04450000000000   0.54450000000000  -0.25577000000000   W  (24k)
  0.75577000000000  -0.04450000000000  -0.04450000000000   W  (24k)
  0.75577000000000   0.54450000000000   0.54450000000000   W  (24k)
 -0.25577000000000   0.54450000000000  -0.04450000000000   W  (24k)
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  0.54450000000000   0.75577000000000   0.54450000000000   W  (24k)
 -0.04450000000000  -0.25577000000000   0.54450000000000   W  (24k)
  0.54450000000000  -0.25577000000000  -0.04450000000000   W  (24k)
  0.45550000000000   0.45550000000000  -0.75577000000000   W  (24k)
  0.04450000000000   0.04450000000000  -0.75577000000000   W  (24k)
  0.45550000000000   0.04450000000000   1.25577000000000   W  (24k)
  0.04450000000000   0.45550000000000   1.25577000000000   W  (24k)
  0.45550000000000   1.25577000000000   0.04450000000000   W  (24k)
  0.04450000000000   1.25577000000000   0.45550000000000   W  (24k)
  0.04450000000000  -0.75577000000000   0.04450000000000   W  (24k)
  0.45550000000000  -0.75577000000000   0.45550000000000   W  (24k)
  1.25577000000000   0.45550000000000   0.04450000000000   W  (24k)
  1.25577000000000   0.04450000000000   0.45550000000000   W  (24k)
 -0.75577000000000   0.45550000000000   0.45550000000000   W  (24k)
 -0.75577000000000   0.04450000000000   0.04450000000000   W  (24k)
```

LaH_{10} High-T_c Superconductor: A10B_cF44_225_cf_b - CIF

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_chemical_formula_sum 'H10 La'

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_aflow_proto 'A10B_cF44_225_cf_b'
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_aflow_params_values '4.78,0.12'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'cF44'

_symmetry_space_group_name_H-M "F 4/m -3 2/m"
_symmetry_Int_Tables_number 225

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_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

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_space_group_symop_operation_xyz
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2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 y,z,x
6 y,-z,-x
7 -y,z,-x
8 -y,-z,x
9 z,x,y
10 z,-x,-y
11 -z,x,-y
12 -z,-x,y
13 -y,-x,-z
14 -y,x,z
15 y,-x,z
16 y,x,-z
17 -x,-z,-y
18 -x,z,y
19 x,-z,y
20 x,z,-y
21 -z,-y,-x
22 -z,y,x
23 z,-y,x
24 z,y,-x
25 -x,-y,-z
26 -x,y,z
27 x,-y,z
28 x,y,-z
29 -y,-z,-x
30 -y,z,x
31 y,-z,x
32 y,z,-x
33 -z,-x,-y
34 -z,x,y
35 z,-x,y
36 z,x,-y
37 y,x,z
38 y,-x,-z
39 -y,x,-z
40 -y,-x,z
41 x,z,y
42 x,-z,-y
43 -x,z,-y
44 -x,-z,y
45 z,y,x
46 z,-y,-x
47 -z,y,-x
48 -z,-y,x
49 x,y+1/2,z+1/2
50 x,-y+1/2,-z+1/2
51 -x,y+1/2,-z+1/2
52 -x,-y+1/2,z+1/2
53 y,z+1/2,x+1/2
54 y,-z+1/2,-x+1/2
55 -y,z+1/2,-x+1/2
56 -y,-z+1/2,x+1/2
57 z,x+1/2,y+1/2
58 z,-x+1/2,-y+1/2
59 -z,x+1/2,-y+1/2
60 -z,-x+1/2,y+1/2
61 -y,-x+1/2,-z+1/2
62 -y,x+1/2,z+1/2
63 y,-x+1/2,z+1/2
64 y,x+1/2,-z+1/2
65 -x,-z+1/2,-y+1/2
66 -x,z+1/2,y+1/2
67 x,-z+1/2,y+1/2
68 x,z+1/2,-y+1/2
69 -z,-y+1/2,-x+1/2
70 -z,y+1/2,x+1/2
71 z,-y+1/2,x+1/2
72 z,y+1/2,-x+1/2
73 -x,-y+1/2,-z+1/2
74 -x,y+1/2,z+1/2
75 x,-y+1/2,z+1/2
76 x,y+1/2,-z+1/2
77 -y,-z+1/2,-x+1/2
78 -y,z+1/2,x+1/2
79 y,-z+1/2,x+1/2
80 y,z+1/2,-x+1/2
81 -z,-x+1/2,-y+1/2
82 -z,x+1/2,y+1/2
83 z,-x+1/2,y+1/2
84 z,x+1/2,-y+1/2
85 y,x+1/2,z+1/2
86 y,-x+1/2,-z+1/2
87 -y,x+1/2,-z+1/2
88 -y,-x+1/2,z+1/2
89 x,z+1/2,y+1/2
90 x,-z+1/2,-y+1/2
91 -x,z+1/2,-y+1/2
92 -x,-z+1/2,y+1/2
93 z,y+1/2,x+1/2
94 z,-y+1/2,-x+1/2
95 -z,y+1/2,-x+1/2
96 -z,-y+1/2,x+1/2
97 x+1/2,y,z+1/2
98 x+1/2,-y,-z+1/2
99 -x+1/2,y,-z+1/2
100 -x+1/2,-y,z+1/2
101 y+1/2,z,x+1/2
102 y+1/2,-z,-x+1/2
103 -y+1/2,z,-x+1/2
104 -y+1/2,-z,x+1/2
105 z+1/2,x,y+1/2
106 z+1/2,-x,-y+1/2
107 -z+1/2,x,-y+1/2
108 -z+1/2,-x,y+1/2
109 -y+1/2,-x,-z+1/2
110 -y+1/2,x,z+1/2
111 y+1/2,-x,z+1/2
112 y+1/2,x,-z+1/2
113 -x+1/2,-z,-y+1/2
114 -x+1/2,z,y+1/2
115 x+1/2,-z,y+1/2
116 x+1/2,z,-y+1/2
117 -z+1/2,-y,-x+1/2
118 -z+1/2,y,x+1/2
119 z+1/2,-y,x+1/2
120 z+1/2,y,-x+1/2
121 -x+1/2,-y,-z+1/2
122 -x+1/2,y,z+1/2
123 x+1/2,-y,z+1/2
124 x+1/2,y,-z+1/2
125 -y+1/2,-z,-x+1/2
126 -y+1/2,z,x+1/2
127 y+1/2,-z,x+1/2
128 y+1/2,z,-x+1/2
129 -z+1/2,-x,-y+1/2
130 -z+1/2,x,y+1/2
131 z+1/2,-x,y+1/2
132 z+1/2,x,-y+1/2
133 y+1/2,x,z+1/2
134 y+1/2,-x,-z+1/2
135 -y+1/2,x,-z+1/2
136 -y+1/2,-x,z+1/2
137 x+1/2,z,y+1/2
138 x+1/2,-z,-y+1/2
```

```
139 -x+1/2,z,-y+1/2
140 -x+1/2,-z,y+1/2
141 z+1/2,y,x+1/2
142 z+1/2,-y,-x+1/2
143 -z+1/2,y,-x+1/2
144 -z+1/2,-y,x+1/2
145 x+1/2,y+1/2,z
146 x+1/2,-y+1/2,-z
147 -x+1/2,y+1/2,-z
148 -x+1/2,-y+1/2,z
149 y+1/2,z+1/2,x
150 y+1/2,-z+1/2,-x
151 -y+1/2,z+1/2,-x
152 -y+1/2,-z+1/2,x
153 z+1/2,x+1/2,y
154 z+1/2,-x+1/2,-y
155 -z+1/2,x+1/2,-y
156 -z+1/2,-x+1/2,y
157 -y+1/2,-x+1/2,-z
158 -y+1/2,x+1/2,z
159 y+1/2,-x+1/2,z
160 y+1/2,x+1/2,-z
161 -x+1/2,-z+1/2,-y
162 -x+1/2,z+1/2,y
163 x+1/2,-z+1/2,y
164 x+1/2,z+1/2,-y
165 -z+1/2,-y+1/2,-x
166 -z+1/2,y+1/2,x
167 z+1/2,-y+1/2,x
168 z+1/2,y+1/2,-x
169 -x+1/2,-y+1/2,-z
170 -x+1/2,y+1/2,z
171 x+1/2,-y+1/2,z
172 x+1/2,y+1/2,-z
173 -y+1/2,-z+1/2,-x
174 -y+1/2,z+1/2,x
175 y+1/2,-z+1/2,x
176 y+1/2,z+1/2,-x
177 -z+1/2,-x+1/2,-y
178 -z+1/2,x+1/2,y
179 z+1/2,-x+1/2,y
180 z+1/2,x+1/2,-y
181 y+1/2,x+1/2,z
182 y+1/2,-x+1/2,-z
183 -y+1/2,x+1/2,-z
184 -y+1/2,-x+1/2,z
185 x+1/2,z+1/2,y
186 x+1/2,-z+1/2,-y
187 -x+1/2,z+1/2,-y
188 -x+1/2,-z+1/2,y
189 z+1/2,y+1/2,x
190 z+1/2,-y+1/2,-x
191 -z+1/2,y+1/2,-x
192 -z+1/2,-y+1/2,x

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
La1 La 4 b 0.50000 0.50000 0.50000 1.00000
H1 H 8 c 0.25000 0.25000 0.25000 1.00000
H2 H 32 f 0.12000 0.12000 0.12000 1.00000
```

LaH_{10} High-T_c Superconductor: A10B_cF44_225_cf_b - POSCAR

```
A10B_cF44_225_cf_b & a,x3 --params=4.78,0.12 & Fm-3m O_{h}^{5} #225 (bcf
    ↪ ) & cF44 & None & H10La & H10La &
   1.00000000000000
     0.00000000000000     2.39000000000000     2.39000000000000
     2.39000000000000     0.00000000000000     2.39000000000000
     2.39000000000000     2.39000000000000     0.00000000000000
     H    La
    10     1
Direct
     0.25000000000000     0.25000000000000     0.25000000000000   H   (8c)
     0.75000000000000     0.75000000000000     0.75000000000000   H   (8c)
     0.12000000000000     0.12000000000000     0.12000000000000   H  (32f)
     0.12000000000000     0.12000000000000    -0.36000000000000   H  (32f)
     0.12000000000000    -0.36000000000000     0.12000000000000   H  (32f)
    -0.36000000000000     0.12000000000000     0.12000000000000   H  (32f)
    -0.12000000000000    -0.12000000000000     0.36000000000000   H  (32f)
    -0.12000000000000    -0.12000000000000    -0.12000000000000   H  (32f)
    -0.12000000000000     0.36000000000000    -0.12000000000000   H  (32f)
     0.36000000000000    -0.12000000000000    -0.12000000000000   H  (32f)
     0.50000000000000     0.50000000000000     0.50000000000000  La  (4b)
```

Double Perovskite (Ba_2MnWO_6): A2BC6D_cF40_225_c_a_e_b - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Double perovskite'
_chemical_formula_sum 'Ba2 Mn O6 W'

loop_
_publ_author_name
 'A. K. Azad'
 'S. A. Ivanov'
 'S.-G. Eriksson'
 'J. Eriksen'
 'H. Rundl\"{o}f'
 'R. Mathieu'
 'P. Svedlindh'
_journal_name_full_name
;
 Materials Research Bulletin
;
_journal_volume 36
_journal_year 2001
_journal_page_first 2215
_journal_page_last 2228
_publ_Section_title
;
 Synthesis, crystal structure, and magnetic characterization of the
    ↪ double perovskite Ba$_{2}$MnWO$_{6}$
;

_aflow_title 'Double Perovskite (Ba$_{2}$MnWO$_{6}$) Structure'
_aflow_proto 'A2BC6D_cF40_225_c_a_e_b'
_aflow_params 'a,x_{4}'
_aflow_params_values '8.19849,0.2654'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'cF40'

_symmetry_space_group_name_H-M "F 4/m -3 2/m"
_symmetry_Int_Tables_number 225

_cell_length_a    8.19849
_cell_length_b    8.19849
_cell_length_c    8.19849
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 y,z,x
6 y,-z,-x
7 -y,z,-x
8 -y,-z,x
9 z,x,y
10 z,-x,-y
11 -z,x,-y
12 -z,-x,y
13 -y,-x,-z
14 -y,x,z
15 y,-x,z
16 y,x,-z
17 -x,-z,-y
18 -x,z,y
19 x,-z,y
20 x,z,-y
21 -z,-y,-x
22 -z,y,x
23 z,-y,x
24 z,y,-x
25 -x,-y,-z
26 -x,y,z
27 x,-y,z
28 x,y,-z
29 -y,-z,-x
30 -y,z,x
31 y,-z,x
32 y,z,-x
33 -z,-x,-y
34 -z,x,y
35 z,-x,y
36 z,x,-y
37 y,x,z
38 y,-x,-z
39 -y,x,-z
40 -y,-x,z
41 x,z,y
42 x,-z,-y
43 -x,z,-y
44 -x,-z,y
45 z,y,x
46 z,-y,-x
47 -z,y,-x
48 -z,-y,x
49 x,y+1/2,z+1/2
50 x,-y+1/2,-z+1/2
51 -x,y+1/2,-z+1/2
52 -x,-y+1/2,z+1/2
53 y,z+1/2,x+1/2
54 y,-z+1/2,-x+1/2
55 -y,z+1/2,-x+1/2
56 -y,-z+1/2,x+1/2
57 z,x+1/2,y+1/2
58 z,-x+1/2,-y+1/2
59 -z,x+1/2,-y+1/2
60 -z,-x+1/2,y+1/2
61 -y,-x+1/2,-z+1/2
62 -y,x+1/2,z+1/2
63 y,-x+1/2,z+1/2
64 y,x+1/2,-z+1/2
65 -x,-z+1/2,-y+1/2
66 -x,z+1/2,y+1/2
67 x,-z+1/2,y+1/2
```

```
68 x,z+1/2,-y+1/2
69 -z,-y+1/2,-x+1/2
70 -z,y+1/2,x+1/2
71 z,-y+1/2,x+1/2
72 z,y+1/2,-x+1/2
73 -x,-y+1/2,-z+1/2
74 -x,y+1/2,z+1/2
75 x,-y+1/2,z+1/2
76 x,y+1/2,-z+1/2
77 -y,-z+1/2,-x+1/2
78 -y,z+1/2,x+1/2
79 y,-z+1/2,x+1/2
80 y,z+1/2,-x+1/2
81 -z,-x+1/2,-y+1/2
82 -z,x+1/2,y+1/2
83 z,-x+1/2,y+1/2
84 z,x+1/2,-y+1/2
85 y,x+1/2,z+1/2
86 y,-x+1/2,-z+1/2
87 -y,x+1/2,-z+1/2
88 -y,-x+1/2,z+1/2
89 x,z+1/2,y+1/2
90 x,-z+1/2,-y+1/2
91 -x,z+1/2,-y+1/2
92 -x,-z+1/2,y+1/2
93 z,y+1/2,x+1/2
94 z,-y+1/2,-x+1/2
95 -z,y+1/2,-x+1/2
96 -z,-y+1/2,x+1/2
97 x+1/2,y,z+1/2
98 x+1/2,-y,-z+1/2
99 -x+1/2,y,-z+1/2
100 -x+1/2,-y,z+1/2
101 y+1/2,z,x+1/2
102 y+1/2,-z,-x+1/2
103 -y+1/2,z,-x+1/2
104 -y+1/2,-z,x+1/2
105 z+1/2,x,y+1/2
106 z+1/2,-x,-y+1/2
107 -z+1/2,x,-y+1/2
108 -z+1/2,-x,y+1/2
109 -y+1/2,-x,-z+1/2
110 -y+1/2,x,z+1/2
111 y+1/2,-x,z+1/2
112 y+1/2,x,-z+1/2
113 -x+1/2,-z,-y+1/2
114 -x+1/2,z,y+1/2
115 x+1/2,-z,y+1/2
116 x+1/2,z,-y+1/2
117 -z+1/2,-y,-x+1/2
118 -z+1/2,y,x+1/2
119 z+1/2,-y,x+1/2
120 z+1/2,y,-x+1/2
121 -x+1/2,-y,-z+1/2
122 -x+1/2,y,z+1/2
123 x+1/2,-y,z+1/2
124 x+1/2,y,-z+1/2
125 -y+1/2,-z,-x+1/2
126 -y+1/2,z,x+1/2
127 y+1/2,-z,x+1/2
128 y+1/2,z,-x+1/2
129 -z+1/2,-x,-y+1/2
130 -z+1/2,x,y+1/2
131 z+1/2,-x,y+1/2
132 z+1/2,x,-y+1/2
133 y+1/2,x,z+1/2
134 y+1/2,-x,-z+1/2
135 -y+1/2,x,-z+1/2
136 -y+1/2,-x,z+1/2
137 x+1/2,z,y+1/2
138 x+1/2,-z,-y+1/2
139 -x+1/2,z,-y+1/2
140 -x+1/2,-z,y+1/2
141 z+1/2,y,x+1/2
142 z+1/2,-y,-x+1/2
143 -z+1/2,y,-x+1/2
144 -z+1/2,-y,x+1/2
145 x+1/2,y+1/2,z
146 x+1/2,-y+1/2,-z
147 -x+1/2,y+1/2,-z
148 -x+1/2,-y+1/2,z
149 y+1/2,z+1/2,x
150 y+1/2,-z+1/2,-x
151 -y+1/2,z+1/2,-x
152 -y+1/2,-z+1/2,x
153 z+1/2,x+1/2,y
154 z+1/2,-x+1/2,-y
155 -z+1/2,x+1/2,-y
156 -z+1/2,-x+1/2,y
157 -y+1/2,-x+1/2,-z
158 -y+1/2,x+1/2,z
159 y+1/2,-x+1/2,z
160 y+1/2,x+1/2,-z
161 -x+1/2,-z+1/2,-y
162 -x+1/2,z+1/2,y
163 x+1/2,-z+1/2,y
164 x+1/2,z+1/2,-y
165 -z+1/2,-y+1/2,-x
166 -z+1/2,y+1/2,x
167 z+1/2,-y+1/2,x
168 z+1/2,y+1/2,-x
169 -x+1/2,-y+1/2,-z
170 -x+1/2,y+1/2,z
171 x+1/2,-y+1/2,z
172 x+1/2,y+1/2,-z
173 -y+1/2,-z+1/2,-x
174 -y+1/2,z+1/2,x
175 y+1/2,-z+1/2,x
176 y+1/2,z+1/2,-x
177 -z+1/2,-x+1/2,-y
178 -z+1/2,x+1/2,y
179 z+1/2,-x+1/2,y
180 z+1/2,x+1/2,-y
181 y+1/2,x+1/2,z
182 y+1/2,-x+1/2,-z
183 -y+1/2,x+1/2,-z
184 -y+1/2,-x+1/2,z
185 x+1/2,z+1/2,y
186 x+1/2,-z+1/2,-y
187 -x+1/2,z+1/2,-y
188 -x+1/2,-z+1/2,y
189 z+1/2,y+1/2,x
190 z+1/2,-y+1/2,-x
191 -z+1/2,y+1/2,-x
192 -z+1/2,-y+1/2,x

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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Mn1 Mn 4 a 0.00000 0.00000 0.00000 1.00000
W1 W 4 b 0.50000 0.50000 0.50000 1.00000
Ba1 Ba 8 c 0.25000 0.25000 0.25000 1.00000
O1 O 24 e 0.26540 0.00000 0.00000 1.00000
```

Double Perovskite (Ba_2MnWO_6): A2BC6D_cF40_225_c_a_e_b - POSCAR

```
A2BC6D_cF40_225_c_a_e_b & a,x4 --params=8.19849,0.2654 & Fm-3m O_{h}^{5}
    ↪ #225 (abce) & cF40 & None & Ba2MnO6W & Double perovskite & A.
    ↪ K. Azad et al., Mater. Res. Bull. 36, 2215-2228 (2001)
   1.00000000000000
    0.00000000000000    4.09924500000000    4.09924500000000
    4.09924500000000    0.00000000000000    4.09924500000000
    4.09924500000000    4.09924500000000    0.00000000000000
     Ba     Mn      O      W
      2      1      6      1
Direct
    0.25000000000000    0.25000000000000    0.25000000000000    Ba    (8c)
    0.75000000000000    0.75000000000000    0.75000000000000    Ba    (8c)
    0.00000000000000    0.00000000000000    0.00000000000000    Mn    (4a)
   -0.26540000000000    0.26540000000000    0.26540000000000    O    (24e)
    0.26540000000000   -0.26540000000000   -0.26540000000000    O    (24e)
    0.26540000000000   -0.26540000000000    0.26540000000000    O    (24e)
   -0.26540000000000    0.26540000000000   -0.26540000000000    O    (24e)
    0.26540000000000    0.26540000000000   -0.26540000000000    O    (24e)
   -0.26540000000000   -0.26540000000000    0.26540000000000    O    (24e)
    0.50000000000000    0.50000000000000    0.50000000000000    W    (4b)
```

$Cu_3[Fe(CN)_6]_2{\cdot}xH_2O$ ($J2_5$, $x \approx 3$): A6B9CD2E6_cF96_225_e_bf_a_c_e - CIF

```
# CIF file
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_audit_creation_method FINDSYM

_chemical_name_mineral 'Prussian blue analog'
_chemical_formula_sum 'C6 Cu9 Fe (H2O)2 N6'

loop_
_publ_author_name
 'A. K. {van Bever}'
_journal_name_full_name
;
 Recueil des Travaux Chimiques des Pays-Bas
;
_journal_volume 57
_journal_year 1938
_journal_page_first 1259
_journal_page_last 1268
_publ_Section_title
;
 The Crystal Structure of Some Ferricyanides with Bivalent Kations
;

_aflow_title 'Cu${_}{3}$[Fe(CN)$_{6}$]$_{2}$\cdot x$H$_{2}$O ($J2_{5}$)
    ↪ Structure'
_aflow_proto 'A6B9CD2E6_cF96_225_e_bf_a_c_e'
_aflow_params 'a,x_{4},x_{5},x_{6}'
_aflow_params_values '10.12,0.20257,0.31126,0.16667'
_aflow_Strukturbericht '$J2_{5}$'
_aflow_Pearson 'cF96'

_symmetry_space_group_name_H-M "F 4/m -3 2/m"
_symmetry_Int_Tables_number 225

_cell_length_a     10.12000
_cell_length_b     10.12000
_cell_length_c     10.12000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
```

```
2 x,−y,−z
3 −x,y,−z
4 −x,−y,z
5 y,z,x
6 y,−z,−x
7 −y,z,−x
8 −y,−z,x
9 z,x,y
10 z,−x,−y
11 −z,x,−y
12 −z,−x,y
13 −y,−x,−z
14 −y,x,z
15 y,−x,z
16 y,x,−z
17 −x,−z,−y
18 −x,z,y
19 x,−z,y
20 x,z,−y
21 −z,−y,−x
22 −z,y,x
23 z,−y,x
24 z,y,−x
25 −x,−y,−z
26 −x,y,z
27 x,−y,z
28 x,y,−z
29 −y,−z,−x
30 −y,z,x
31 y,−z,x
32 y,z,−x
33 −z,−x,−y
34 −z,x,y
35 z,−x,y
36 z,x,−y
37 y,x,z
38 y,−x,−z
39 −y,x,−z
40 −y,−x,z
41 x,z,y
42 x,−z,−y
43 −x,z,−y
44 −x,−z,y
45 z,y,x
46 z,−y,−x
47 −z,y,−x
48 −z,−y,x
49 x,y+1/2,z+1/2
50 x,−y+1/2,−z+1/2
51 −x,y+1/2,−z+1/2
52 −x,−y+1/2,z+1/2
53 y,z+1/2,x+1/2
54 y,−z+1/2,−x+1/2
55 −y,z+1/2,−x+1/2
56 −y,−z+1/2,x+1/2
57 z,x+1/2,y+1/2
58 z,−x+1/2,−y+1/2
59 −z,x+1/2,−y+1/2
60 −z,−x+1/2,y+1/2
61 −y,−x+1/2,−z+1/2
62 −y,x+1/2,z+1/2
63 y,−x+1/2,z+1/2
64 y,x+1/2,−z+1/2
65 −x,−z+1/2,−y+1/2
66 −x,z+1/2,y+1/2
67 x,−z+1/2,y+1/2
68 x,z+1/2,−y+1/2
69 −z,−y+1/2,−x+1/2
70 −z,y+1/2,x+1/2
71 z,−y+1/2,x+1/2
72 z,y+1/2,−x+1/2
73 −x,−y+1/2,−z+1/2
74 −x,y+1/2,z+1/2
75 x,−y+1/2,z+1/2
76 x,y+1/2,−z+1/2
77 −y,−z+1/2,−x+1/2
78 −y,z+1/2,x+1/2
79 y,−z+1/2,x+1/2
80 y,z+1/2,−x+1/2
81 −z,−x+1/2,−y+1/2
82 −z,x+1/2,y+1/2
83 z,−x+1/2,y+1/2
84 z,x+1/2,−y+1/2
85 y,x+1/2,z+1/2
86 y,−x+1/2,−z+1/2
87 −y,x+1/2,−z+1/2
88 −y,−x+1/2,z+1/2
89 x,z+1/2,y+1/2
90 x,−z+1/2,−y+1/2
91 −x,z+1/2,−y+1/2
92 −x,−z+1/2,y+1/2
93 z,y+1/2,x+1/2
94 z,−y+1/2,−x+1/2
95 −z,y+1/2,−x+1/2
96 −z,−y+1/2,x+1/2
97 x+1/2,y,z+1/2
98 x+1/2,−y,−z+1/2
99 −x+1/2,y,−z+1/2
100 −x+1/2,−y,z+1/2
101 y+1/2,z,x+1/2
102 y+1/2,−z,−x+1/2
103 −y+1/2,z,−x+1/2
104 −y+1/2,−z,x+1/2
105 z+1/2,x,y+1/2
106 z+1/2,−x,−y+1/2
107 −z+1/2,x,−y+1/2
108 −z+1/2,−x,y+1/2
109 −y+1/2,−x,−z+1/2
110 −y+1/2,x,z+1/2
111 y+1/2,−x,z+1/2
112 y+1/2,x,−z+1/2
113 −x+1/2,−z,−y+1/2
114 −x+1/2,z,y+1/2
115 x+1/2,−z,y+1/2
116 x+1/2,z,−y+1/2
117 −z+1/2,−y,−x+1/2
118 −z+1/2,y,x+1/2
119 z+1/2,−y,x+1/2
120 z+1/2,y,−x+1/2
121 −x+1/2,−y,−z+1/2
122 −x+1/2,y,z+1/2
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124 x+1/2,y,−z+1/2
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126 −y+1/2,z,x+1/2
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128 y+1/2,z,−x+1/2
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131 z+1/2,−x,y+1/2
132 z+1/2,x,−y+1/2
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134 y+1/2,−x,−z+1/2
135 −y+1/2,x,−z+1/2
136 −y+1/2,−x,z+1/2
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138 x+1/2,−z,−y+1/2
139 −x+1/2,z,−y+1/2
140 −x+1/2,−z,y+1/2
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142 z+1/2,−y,−x+1/2
143 −z+1/2,y,−x+1/2
144 −z+1/2,−y,x+1/2
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146 x+1/2,−y+1/2,−z
147 −x+1/2,y+1/2,−z
148 −x+1/2,−y+1/2,z
149 y+1/2,z+1/2,x
150 y+1/2,−z+1/2,−x
151 −y+1/2,z+1/2,−x
152 −y+1/2,−z+1/2,x
153 z+1/2,x+1/2,y
154 z+1/2,−x+1/2,−y
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156 −z+1/2,−x+1/2,y
157 −y+1/2,−x+1/2,−z
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160 y+1/2,x+1/2,−z
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162 −x+1/2,z+1/2,y
163 x+1/2,−z+1/2,y
164 x+1/2,z+1/2,−y
165 −z+1/2,−y+1/2,−x
166 −z+1/2,y+1/2,x
167 z+1/2,−y+1/2,x
168 z+1/2,y+1/2,−x
169 −x+1/2,−y+1/2,−z
170 −x+1/2,y+1/2,z
171 x+1/2,−y+1/2,z
172 x+1/2,y+1/2,−z
173 −y+1/2,−z+1/2,−x
174 −y+1/2,z+1/2,x
175 y+1/2,−z+1/2,x
176 y+1/2,z+1/2,−x
177 −z+1/2,−x+1/2,−y
178 −z+1/2,x+1/2,y
179 z+1/2,−x+1/2,y
180 z+1/2,x+1/2,−y
181 y+1/2,x+1/2,z
182 y+1/2,−x+1/2,−z
183 −y+1/2,x+1/2,−z
184 −y+1/2,−x+1/2,z
185 x+1/2,z+1/2,y
186 x+1/2,−z+1/2,−y
187 −x+1/2,z+1/2,−y
188 −x+1/2,−z+1/2,y
189 z+1/2,y+1/2,x
190 z+1/2,−y+1/2,−x
191 −z+1/2,y+1/2,−x
192 −z+1/2,−y+1/2,x

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_atom_site_fract_x
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_atom_site_fract_z
_atom_site_occupancy
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Cu1 Cu 4 b 0.50000 0.50000 0.50000 1.00000
H2O1 H2O 8 c 0.25000 0.25000 0.25000 0.75000
C1 C 24 e 0.20257 0.00000 0.00000 1.00000
N1 N 24 e 0.31126 0.00000 0.00000 1.00000
Cu2 Cu 32 f 0.16667 0.16667 0.16667 0.06250
```

$Cu_3[Fe(CN)_6]_2{\cdot}xH_2O$ ($J2_5$, $x \approx 3$): A6B9CD2E6_cF96_225_e_bf_a_c_e - POSCAR

```
A6B9CD2E6_cF96_225_e_bf_a_c_e & a,x4,x5,x6 --params=10.12,0.20257,
    ↪ 0.31126,0.16667 & Fm-3m O_{h}^{5} #225 (abce^2f) & cF96 & $J2_{
    ↪ 5}$ & C12Cu3Fe2N12\cdot xH2O & Prussian blue analog & A. K. {
    ↪ van Bever}, Rec. Trav. Chim. Pays-Bas 57, 1259-1268 (1938)
   1.00000000000000
   0.00000000000000    5.06000000000000    5.06000000000000
   5.06000000000000    0.00000000000000    5.06000000000000
   5.06000000000000    5.06000000000000    0.00000000000000
     C     Cu     Fe    H2O      N
     6      9      1      2      6
Direct
  -0.20257000000000    0.20257000000000    0.20257000000000     C   (24e)
   0.20257000000000   -0.20257000000000   -0.20257000000000     C   (24e)
   0.20257000000000   -0.20257000000000    0.20257000000000     C   (24e)
  -0.20257000000000    0.20257000000000   -0.20257000000000     C   (24e)
   0.20257000000000    0.20257000000000   -0.20257000000000     C   (24e)
  -0.20257000000000   -0.20257000000000    0.20257000000000     C   (24e)
   0.50000000000000    0.50000000000000    0.50000000000000    Cu    (4b)
   0.16667000000000    0.16667000000000    0.16667000000000    Cu   (32f)
   0.16667000000000    0.16667000000000   -0.50001000000000    Cu   (32f)
   0.16667000000000   -0.50001000000000    0.16667000000000    Cu   (32f)
  -0.50001000000000    0.16667000000000    0.16667000000000    Cu   (32f)
  -0.16667000000000   -0.16667000000000    0.50001000000000    Cu   (32f)
  -0.16667000000000   -0.16667000000000   -0.16667000000000    Cu   (32f)
  -0.16667000000000    0.50001000000000   -0.16667000000000    Cu   (32f)
   0.50001000000000   -0.16667000000000   -0.16667000000000    Cu   (32f)
   0.00000000000000    0.00000000000000    0.00000000000000    Fe    (4a)
   0.25000000000000    0.25000000000000    0.25000000000000   H2O    (8c)
   0.75000000000000    0.75000000000000    0.75000000000000   H2O    (8c)
  -0.31126000000000    0.31126000000000    0.31126000000000     N   (24e)
   0.31126000000000   -0.31126000000000   -0.31126000000000     N   (24e)
   0.31126000000000   -0.31126000000000    0.31126000000000     N   (24e)
  -0.31126000000000    0.31126000000000   -0.31126000000000     N   (24e)
   0.31126000000000    0.31126000000000   -0.31126000000000     N   (24e)
  -0.31126000000000   -0.31126000000000    0.31126000000000     N   (24e)
```

Co_9S_8 ($D8_9$): A9B8_cF68_225_af_ce - CIF

```
# CIF file
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_audit_creation_method FINDSYM

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_chemical_formula_sum 'Co9 S8'

loop_
_publ_author_name
 'S. Geller'
_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 15
_journal_year 1962
_journal_page_first 1195
_journal_page_last 1198
_publ_Section_title
;
 Refinement of the crystal structure of Co$_{9}$S$_{8}$
;

_aflow_title 'Co$_{9}$S$_{8}$ ($D8_{9}$) Structure'
_aflow_proto 'A9B8_cF68_225_af_ce'
_aflow_params 'a,x_{3},x_{4}'
_aflow_params_values '9.928,0.7591,0.374'
_aflow_Strukturbericht '$D8_{9}$'
_aflow_Pearson 'cF68'

_symmetry_space_group_name_H-M "F 4/m -3 2/m"
_symmetry_Int_Tables_number 225

_cell_length_a    9.92800
_cell_length_b    9.92800
_cell_length_c    9.92800
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

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_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 y,z,x
6 y,-z,-x
7 -y,z,-x
8 -y,-z,x
9 z,x,y
10 z,-x,-y
11 -z,x,-y
12 -z,-x,y
13 -y,-x,-z
14 -y,x,z
15 y,-x,z
16 y,x,-z
17 -x,-z,-y
18 -x,z,y
19 x,-z,y
20 x,z,-y
21 -z,-y,-x
22 -z,y,x
23 z,-y,x
24 z,y,-x
25 -x,-y,-z
26 -x,y,z
27 x,-y,z
28 x,y,-z
29 -y,-z,-x
30 -y,z,x
31 y,-z,x
32 y,z,-x
33 -z,-x,-y
34 -z,x,y
35 z,-x,y
36 z,x,-y
37 y,x,z
38 y,-x,-z
39 -y,x,-z
40 -y,-x,z
41 x,z,y
42 x,-z,-y
43 -x,z,-y
44 -x,-z,y
45 z,y,x
46 z,-y,-x
47 -z,y,-x
48 -z,-y,x
49 x,y+1/2,z+1/2
50 x,-y+1/2,-z+1/2
51 -x,y+1/2,-z+1/2
52 -x,-y+1/2,z+1/2
53 y,z+1/2,x+1/2
54 y,-z+1/2,-x+1/2
55 -y,z+1/2,-x+1/2
56 -y,-z+1/2,x+1/2
57 z,x+1/2,y+1/2
58 z,-x+1/2,-y+1/2
59 -z,x+1/2,-y+1/2
60 -z,-x+1/2,y+1/2
61 -y,-x+1/2,-z+1/2
62 -y,x+1/2,z+1/2
63 y,-x+1/2,z+1/2
64 y,x+1/2,-z+1/2
65 -x,-z+1/2,-y+1/2
66 -x,z+1/2,y+1/2
67 x,-z+1/2,y+1/2
68 x,z+1/2,-y+1/2
69 -z,-y+1/2,-x+1/2
70 -z,y+1/2,x+1/2
71 z,-y+1/2,x+1/2
72 z,y+1/2,-x+1/2
73 -x,-y+1/2,-z+1/2
74 -x,y+1/2,z+1/2
75 x,-y+1/2,z+1/2
76 x,y+1/2,-z+1/2
77 -y,-z+1/2,-x+1/2
78 -y,z+1/2,x+1/2
79 y,-z+1/2,x+1/2
80 y,z+1/2,-x+1/2
81 -z,-x+1/2,-y+1/2
82 -z,x+1/2,y+1/2
83 z,-x+1/2,y+1/2
84 z,x+1/2,-y+1/2
85 y,x+1/2,z+1/2
86 y,-x+1/2,-z+1/2
87 -y,x+1/2,-z+1/2
88 -y,-x+1/2,z+1/2
89 x,z+1/2,y+1/2
90 x,-z+1/2,-y+1/2
91 -x,z+1/2,-y+1/2
92 -x,-z+1/2,y+1/2
93 z,y+1/2,x+1/2
94 z,-y+1/2,-x+1/2
95 -z,y+1/2,-x+1/2
96 -z,-y+1/2,x+1/2
97 x+1/2,y,z+1/2
98 x+1/2,-y,-z+1/2
99 -x+1/2,y,-z+1/2
100 -x+1/2,-y,z+1/2
101 y+1/2,z,x+1/2
102 y+1/2,-z,-x+1/2
103 -y+1/2,z,-x+1/2
104 -y+1/2,-z,x+1/2
105 z+1/2,x,y+1/2
106 z+1/2,-x,-y+1/2
107 -z+1/2,x,-y+1/2
108 -z+1/2,-x,y+1/2
109 -y+1/2,-x,-z+1/2
110 -y+1/2,x,z+1/2
111 y+1/2,-x,z+1/2
112 y+1/2,x,-z+1/2
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114 -x+1/2,z,y+1/2
115 x+1/2,-z,y+1/2
116 x+1/2,z,-y+1/2
117 -z+1/2,-y,-x+1/2
118 -z+1/2,y,x+1/2
119 z+1/2,-y,x+1/2
120 z+1/2,y,-x+1/2
121 -x+1/2,-y,-z+1/2
122 -x+1/2,y,z+1/2
123 x+1/2,-y,z+1/2
124 x+1/2,y,-z+1/2
125 -y+1/2,-z,-x+1/2
126 -y+1/2,z,x+1/2
127 y+1/2,-z,x+1/2
128 y+1/2,z,-x+1/2
129 -z+1/2,-x,-y+1/2
```

```
130 -z+1/2,x,y+1/2
131 z+1/2,-x,y+1/2
132 z+1/2,x,-y+1/2
133 y+1/2,x,z+1/2
134 y+1/2,-x,-z+1/2
135 -y+1/2,x,-z+1/2
136 -y+1/2,-x,z+1/2
137 x+1/2,z,y+1/2
138 x+1/2,-z,-y+1/2
139 -x+1/2,z,-y+1/2
140 -x+1/2,-z,y+1/2
141 z+1/2,y,x+1/2
142 z+1/2,-y,-x+1/2
143 -z+1/2,y,-x+1/2
144 -z+1/2,-y,x+1/2
145 x+1/2,y+1/2,z
146 x+1/2,-y+1/2,-z
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148 -x+1/2,-y+1/2,z
149 y+1/2,z+1/2,x
150 y+1/2,-z+1/2,-x
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152 -y+1/2,-z+1/2,x
153 z+1/2,x+1/2,y
154 z+1/2,-x+1/2,-y
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161 -x+1/2,-z+1/2,-y
162 -x+1/2,z+1/2,y
163 x+1/2,-z+1/2,y
164 x+1/2,z+1/2,-y
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166 -z+1/2,y+1/2,x
167 z+1/2,-y+1/2,x
168 z+1/2,y+1/2,-x
169 -x+1/2,-y+1/2,-z
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172 x+1/2,y+1/2,-z
173 -y+1/2,-z+1/2,-x
174 -y+1/2,z+1/2,x
175 y+1/2,-z+1/2,x
176 y+1/2,z+1/2,-x
177 -z+1/2,-x+1/2,-y
178 -z+1/2,x+1/2,y
179 z+1/2,-x+1/2,y
180 z+1/2,x+1/2,-y
181 y+1/2,x+1/2,z
182 y+1/2,-x+1/2,-z
183 -y+1/2,x+1/2,-z
184 -y+1/2,-x+1/2,z
185 x+1/2,z+1/2,y
186 x+1/2,-z+1/2,-y
187 -x+1/2,z+1/2,-y
188 -x+1/2,-z+1/2,y
189 z+1/2,y+1/2,x
190 z+1/2,-y+1/2,-x
191 -z+1/2,y+1/2,-x
192 -z+1/2,-y+1/2,x

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_atom_site_Wyckoff_label
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_atom_site_occupancy
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S1 S 8 c 0.25000 0.25000 0.25000 1.00000
S2 S 24 e 0.75910 0.00000 0.00000 1.00000
Co2 Co 32 f 0.37400 0.37400 0.37400 1.00000
```

Co_9S_8 ($D8_9$): A9B8_cF68_225_af_ce - POSCAR

```
A9B8_cF68_225_af_ce & a,x3,x4 --params=9.928,0.7591,0.374 & Fm-3m O_{h
    ↪ }^{5} #225 (acef) & cF68 & $D8_{9}$ & Co9S8 & Co9S8 & S. Geller
    ↪ , Acta Cryst. 15, 1195-1198 (1962)
   1.00000000000000
     0.00000000000000     4.96400000000000     4.96400000000000
     4.96400000000000     0.00000000000000     4.96400000000000
     4.96400000000000     4.96400000000000     0.00000000000000
      Co       S
       9       8
Direct
     0.00000000000000     0.00000000000000     0.00000000000000     Co     (4a)
     0.37400000000000     0.37400000000000     0.37400000000000     Co     (32f)
     0.37400000000000     0.37400000000000    -1.12200000000000     Co     (32f)
     0.37400000000000    -1.12200000000000     0.37400000000000     Co     (32f)
    -1.12200000000000     0.37400000000000     0.37400000000000     Co     (32f)
    -0.37400000000000    -0.37400000000000     1.12200000000000     Co     (32f)
    -0.37400000000000    -0.37400000000000    -0.37400000000000     Co     (32f)
    -0.37400000000000     1.12200000000000    -0.37400000000000     Co     (32f)
     1.12200000000000    -0.37400000000000    -0.37400000000000     Co     (32f)
     0.25000000000000     0.25000000000000     0.25000000000000     S      (8c)
     0.75000000000000     0.75000000000000     0.75000000000000     S      (8c)
    -0.75910000000000     0.75910000000000     0.75910000000000     S      (24e)
     0.75910000000000    -0.75910000000000    -0.75910000000000     S      (24e)
     0.75910000000000    -0.75910000000000     0.75910000000000     S      (24e)
    -0.75910000000000     0.75910000000000    -0.75910000000000     S      (24e)
     0.75910000000000     0.75910000000000    -0.75910000000000     S      (24e)
    -0.75910000000000    -0.75910000000000     0.75910000000000     S      (24e)
```

$(NH_4)_3AlF_6$ ($J2_1$): AB30C16D3_cF200_225_a_ej_2f_bc - CIF

```
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_publ_author_name
 'A. A. Udovenko'
 'N. M. Laptash'
 'I. G. Maslennikova'
_journal_name_full_name
;
 Journal of Fluorine Chemistry
;
_journal_volume 124
_journal_year 2003
_journal_page_first 5
_journal_page_last 15
_publ_Section_title
;
 Orientation disorder in ammonium elpasolites: Crystal structures of (
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;

_aflow_title '(NH$_{4}$)$_{3}$AlF$_{6}$ ($J2_{1}$) Structure'
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```

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189 z+1/2,y+1/2,x
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N2 N 8 c 0.25000 0.25000 0.25000 1.00000
F1 F 24 e 0.19690 0.00000 0.00000 0.33333
H1 H 32 f 0.30700 0.30700 0.30700 1.00000
H2 H 32 f 0.55700 0.55700 0.55700 0.50000
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$(NH_4)_3AlF_6$ ($J2_1$): AB30C16D3_cF200_225_a_ej_2f_bc - POSCAR

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    ↪ 2j) & cF200 & $J2_{1}$ & AlF6H12N3 & AlF6H12N3 & A. A. Udovenko
    ↪  and N. M. Laptash and I. G. Maslennikova, J. Fluor. Chem. 124,
    ↪  5-15 (2003)
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   4.47005500000000   4.47005500000000   0.00000000000000
    Al    F    H    N
     1   30   16    3
Direct
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   0.19690000000000  -0.19690000000000  -0.19690000000000   F   (24e)
   0.19690000000000  -0.19690000000000   0.19690000000000   F   (24e)
  -0.19690000000000   0.19690000000000  -0.19690000000000   F   (24e)
   0.19690000000000   0.19690000000000  -0.19690000000000   F   (24e)
  -0.19690000000000  -0.19690000000000   0.19690000000000   F   (24e)
   0.24590000000000   0.13870000000000  -0.13870000000000   F   (96j)
   0.13870000000000   0.24590000000000  -0.24590000000000   F   (96j)
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  -0.24590000000000   0.13870000000000   0.24590000000000   F   (96j)
   0.24590000000000  -0.13870000000000   0.24590000000000   F   (96j)
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  -0.13870000000000   0.13870000000000  -0.24590000000000   F   (96j)
  -0.24590000000000  -0.13870000000000   0.24590000000000   F   (96j)
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   0.13870000000000  -0.13870000000000  -0.24590000000000   F   (96j)
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   0.30700000000000   0.30700000000000  -0.92100000000000   H   (32f)
   0.30700000000000  -0.92100000000000   0.30700000000000   H   (32f)
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   0.92100000000000  -0.30700000000000  -0.30700000000000   H   (32f)
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  -1.67100000000000   0.55700000000000   0.55700000000000   H   (32f)
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   1.67100000000000  -0.55700000000000  -0.55700000000000   H   (32f)
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$L1_a$ (disputed $CuPt_3$): AB7_cF32_225_b_ad - CIF

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_journal_name_full_name
;
 Acta Cristallographica
;
_journal_volume 4
_journal_year 1951
_journal_page_first 377
_journal_page_last 378
_publ_Section_title
;
A cubic structure for the phase Pt$_{3}$Cu
;

# Found in Revisiting the CuPt$_{3}$ prototype and the $L1_{3}$
      ↪ structure, 2014

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_symmetry_Int_Tables_number 225

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2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
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6 y,-z,-x
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8 -y,-z,x
9 z,x,y
10 z,-x,-y
11 -z,x,-y
12 -z,-x,y
13 -y,-x,-z
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15 y,-x,z
16 y,x,-z
17 -x,-z,-y
18 -x,z,y
19 x,-z,y
20 x,z,-y
21 -z,-y,-x
22 -z,y,x
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24 z,y,-x
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27 x,-y,z
28 x,y,-z
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30 -y,z,x
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35 z,-x,y
36 z,x,-y
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38 y,-x,-z
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44 -x,-z,y
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53 y,z+1/2,x+1/2
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76 x,y+1/2,-z+1/2
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87 -y,x+1/2,-z+1/2
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150 y+1/2,-z+1/2,-x
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165 -z+1/2,-y+1/2,-x
166 -z+1/2,y+1/2,x
167 z+1/2,-y+1/2,x
168 z+1/2,y+1/2,-x
169 -x+1/2,-y+1/2,-z
170 -x+1/2,y+1/2,z
171 x+1/2,-y+1/2,z
172 x+1/2,y+1/2,-z
173 -y+1/2,-z+1/2,-x
174 -y+1/2,z+1/2,x
175 y+1/2,-z+1/2,x
176 y+1/2,z+1/2,-x
177 -z+1/2,-x+1/2,-y
178 -z+1/2,x+1/2,y
179 z+1/2,-x+1/2,y
180 z+1/2,x+1/2,-y
181 y+1/2,x+1/2,z
182 y+1/2,-x+1/2,-z
183 -y+1/2,x+1/2,-z
184 -y+1/2,-x+1/2,z
185 x+1/2,z+1/2,y
186 x+1/2,-z+1/2,-y
187 -x+1/2,z+1/2,-y
188 -x+1/2,-z+1/2,y
189 z+1/2,y+1/2,x
190 z+1/2,-y+1/2,-x
191 -z+1/2,y+1/2,-x
192 -z+1/2,-y+1/2,x

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Pt1 Pt 4 a 0.00000 0.00000 0.00000 1.00000
Cu1 Cu 4 b 0.50000 0.50000 0.50000 1.00000
Pt2 Pt 24 d 0.00000 0.25000 0.25000 1.00000
```

$L1_a$ (disputed $CuPt_3$): AB7_cF32_225_b_ad - POSCAR

```
AB7_cF32_225_b_ad & a --params=5.6 & Fm-3m O_{h}^{5} #225 (abd) & cF32 &
    ↪ $L1_{a}$ & CuPt3 & CuPt3 & Y.-C. Tang, Acta Cryst. 4, 377-378
    ↪ (1951)
   1.00000000000000
   0.00000000000000   2.80000000000000   2.80000000000000
   2.80000000000000   0.00000000000000   2.80000000000000
   2.80000000000000   2.80000000000000   0.00000000000000
    Cu    Pt
     1     7
Direct
   0.50000000000000   0.50000000000000   0.50000000000000   Cu   (4b)
   0.00000000000000   0.00000000000000   0.00000000000000   Pt   (4a)
   0.50000000000000   0.00000000000000   0.00000000000000   Pt   (24d)
   0.00000000000000   0.50000000000000   0.50000000000000   Pt   (24d)
   0.00000000000000   0.50000000000000   0.00000000000000   Pt   (24d)
   0.50000000000000   0.00000000000000   0.50000000000000   Pt   (24d)
   0.00000000000000   0.00000000000000   0.50000000000000   Pt   (24d)
   0.50000000000000   0.50000000000000   0.00000000000000   Pt   (24d)
```

Sulphohalite [$Na_6ClF(SO_4)_2$, $H5_8$]: ABC6D8E2_cF72_225_b_a_e_f_c - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Sulphohalite'
_chemical_formula_sum 'Cl F Na6 O8 S2'

loop_
_publ_author_name
 'A. Pabst'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 89
_journal_year 1934
_journal_page_first 514
_journal_page_last 517
_publ_Section_title
;
 The Crystal Structure of Sulphohalite
;

# Found in Strukturbericht Band III 1933-1935, 1937

_aflow_title 'Sulphohalite [Na$_{6}$ClF(SO$_{4}$)$_{2}$, $H5_{8}$]
    ↪ Structure'
_aflow_proto 'ABC6D8E2_cF72_225_b_a_e_f_c'
_aflow_params 'a,x_{4},x_{5}'
_aflow_params_values '10.08,0.226,0.164'
_aflow_Strukturbericht '$H5_{8}$'
_aflow_Pearson 'cF72'

_symmetry_space_group_name_H-M "F 4/m -3 2/m"
_symmetry_Int_Tables_number 225

_cell_length_a     10.08000
_cell_length_b     10.08000
_cell_length_c     10.08000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 y,z,x
6 y,-z,-x
7 -y,z,-x
8 -y,-z,x
9 z,x,y
10 z,-x,-y
11 -z,x,-y
12 -z,-x,y
13 -y,-x,-z
14 -y,x,z
15 y,-x,z
16 y,x,-z
17 -x,-z,-y
18 -x,z,y
19 x,-z,y
20 x,z,-y
21 -z,-y,-x
22 -z,y,x
23 z,-y,x
24 z,y,-x
25 -x,-y,-z
26 -x,y,z
27 x,-y,z
28 x,y,-z
29 -y,-z,-x
30 -y,z,x
31 y,-z,x
32 y,z,-x
33 -z,-x,-y
34 -z,x,y
35 z,-x,y
36 z,x,-y
37 y,x,z
38 y,-x,-z
39 -y,x,-z
40 -y,-x,z
41 x,z,y
42 x,-z,-y
43 -x,z,-y
44 -x,-z,y
45 z,y,x
46 z,-y,-x
47 -z,y,-x
48 -z,-y,x
49 x,y+1/2,z+1/2
50 x,-y+1/2,-z+1/2
51 -x,y+1/2,-z+1/2
52 -x,-y+1/2,z+1/2
53 y,z+1/2,x+1/2
54 y,-z+1/2,-x+1/2
55 -y,z+1/2,-x+1/2
56 -y,-z+1/2,x+1/2
57 z,x+1/2,y+1/2
58 z,-x+1/2,-y+1/2
59 -z,x+1/2,-y+1/2
60 -z,-x+1/2,y+1/2
61 -y,-x+1/2,-z+1/2
62 -y,x+1/2,z+1/2
63 y,-x+1/2,z+1/2
64 y,x+1/2,-z+1/2
65 -x,-z+1/2,-y+1/2
66 -x,z+1/2,y+1/2
67 x,-z+1/2,y+1/2
68 x,z+1/2,-y+1/2
69 -z,-y+1/2,-x+1/2
70 -z,y+1/2,x+1/2
71 z,-y+1/2,x+1/2
72 z,y+1/2,-x+1/2
73 -x,-y+1/2,-z+1/2
74 -x,y+1/2,z+1/2
75 x,-y+1/2,z+1/2
76 x,y+1/2,-z+1/2
77 -y,-z+1/2,-x+1/2
78 -y,z+1/2,x+1/2
79 y,-z+1/2,x+1/2
80 y,z+1/2,-x+1/2
81 -z,-x+1/2,-y+1/2
82 -z,x+1/2,y+1/2
83 z,-x+1/2,y+1/2
84 z,x+1/2,-y+1/2
```

```
85 y,x+1/2,z+1/2
86 y,-x+1/2,-z+1/2
87 -y,x+1/2,-z+1/2
88 -y,-x+1/2,z+1/2
89 x,z+1/2,y+1/2
90 x,-z+1/2,-y+1/2
91 -x,z+1/2,-y+1/2
92 -x,-z+1/2,y+1/2
93 z,y+1/2,x+1/2
94 z,-y+1/2,-x+1/2
95 -z,y+1/2,-x+1/2
96 -z,-y+1/2,x+1/2
97 x+1/2,y,z+1/2
98 x+1/2,-y,-z+1/2
99 -x+1/2,y,-z+1/2
100 -x+1/2,-y,z+1/2
101 y+1/2,z,x+1/2
102 y+1/2,-z,-x+1/2
103 -y+1/2,z,-x+1/2
104 -y+1/2,-z,x+1/2
105 z+1/2,x,y+1/2
106 z+1/2,-x,-y+1/2
107 -z+1/2,x,-y+1/2
108 -z+1/2,-x,y+1/2
109 -y+1/2,-x,-z+1/2
110 -y+1/2,x,z+1/2
111 y+1/2,-x,z+1/2
112 y+1/2,x,-z+1/2
113 -x+1/2,-z,-y+1/2
114 -x+1/2,z,y+1/2
115 x+1/2,-z,y+1/2
116 x+1/2,z,-y+1/2
117 -z+1/2,-y,-x+1/2
118 -z+1/2,y,x+1/2
119 z+1/2,-y,x+1/2
120 z+1/2,y,-x+1/2
121 -x+1/2,-y,-z+1/2
122 -x+1/2,y,z+1/2
123 x+1/2,-y,z+1/2
124 x+1/2,y,-z+1/2
125 -y+1/2,-z,-x+1/2
126 -y+1/2,z,x+1/2
127 y+1/2,-z,x+1/2
128 y+1/2,z,-x+1/2
129 -z+1/2,-x,-y+1/2
130 -z+1/2,x,y+1/2
131 z+1/2,-x,y+1/2
132 z+1/2,x,-y+1/2
133 y+1/2,x,z+1/2
134 y+1/2,-x,-z+1/2
135 -y+1/2,x,-z+1/2
136 -y+1/2,-x,z+1/2
137 x+1/2,z,y+1/2
138 x+1/2,-z,-y+1/2
139 -x+1/2,z,-y+1/2
140 -x+1/2,-z,y+1/2
141 z+1/2,y,x+1/2
142 z+1/2,-y,-x+1/2
143 -z+1/2,y,-x+1/2
144 -z+1/2,-y,x+1/2
145 x+1/2,y+1/2,z
146 x+1/2,-y+1/2,-z
147 -x+1/2,y+1/2,-z
148 -x+1/2,-y+1/2,z
149 y+1/2,z+1/2,x
150 y+1/2,-z+1/2,-x
151 -y+1/2,z+1/2,-x
152 -y+1/2,-z+1/2,x
153 z+1/2,x+1/2,y
154 z+1/2,-x+1/2,-y
155 -z+1/2,x+1/2,-y
156 -z+1/2,-x+1/2,y
157 -y+1/2,-x+1/2,-z
158 -y+1/2,x+1/2,z
159 y+1/2,-x+1/2,z
160 y+1/2,x+1/2,-z
161 -x+1/2,-z+1/2,-y
162 -x+1/2,z+1/2,y
163 x+1/2,-z+1/2,y
164 x+1/2,z+1/2,-y
165 -z+1/2,-y+1/2,-x
166 -z+1/2,y+1/2,x
167 z+1/2,-y+1/2,x
168 z+1/2,y+1/2,-x
169 -x+1/2,-y+1/2,-z
170 -x+1/2,y+1/2,z
171 x+1/2,-y+1/2,z
172 x+1/2,y+1/2,-z
173 -y+1/2,-z+1/2,-x
174 -y+1/2,z+1/2,x
175 y+1/2,-z+1/2,x
176 y+1/2,z+1/2,-x
177 -z+1/2,-x+1/2,-y
178 -z+1/2,x+1/2,y
179 z+1/2,-x+1/2,y
180 z+1/2,x+1/2,-y
181 y+1/2,x+1/2,z
182 y+1/2,-x+1/2,-z
183 -y+1/2,x+1/2,-z
184 -y+1/2,-x+1/2,z
185 x+1/2,z+1/2,y
186 x+1/2,-z+1/2,-y
187 -x+1/2,z+1/2,-y
188 -x+1/2,-z+1/2,y
189 z+1/2,y+1/2,x
190 z+1/2,-y+1/2,-x
191 -z+1/2,y+1/2,-x
192 -z+1/2,-y+1/2,x

loop_
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_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
F1 F 4 a 0.00000 0.00000 0.00000 1.00000
Cl1 Cl 4 b 0.50000 0.50000 0.50000 1.00000
S1 S 8 c 0.25000 0.25000 0.25000 1.00000
Na1 Na 24 e 0.22600 0.00000 0.00000 1.00000
O1 O 32 f 0.16400 0.16400 0.16400 1.00000
```

Sulphohalite [$Na_6ClF(SO_4)_2$, $H5_8$]: ABC6D8E2_cF72_225_b_a_e_f_c - POSCAR

```
ABC6D8E2_cF72_225_b_a_e_f_c & a,x4,x5 --params=10.08,0.226,0.164 & Fm-3m
    ↪ O_{h}^{5} #225 (abcef) & cF72 & $H5_{8}$ & ClFNa6O8S2 &
    ↪ Sulphohalite & A. Pabst, Zeitschrift f{"u}r Kristallographie -
    ↪ Crystalline Materials 89, 514-517 (1934)
   1.00000000000000
   0.00000000000000   5.04000000000000   5.04000000000000
   5.04000000000000   0.00000000000000   5.04000000000000
   5.04000000000000   5.04000000000000   0.00000000000000
    Cl    F    Na    O    S
     1    1     6    8    2
Direct
   0.50000000000000   0.50000000000000   0.50000000000000   Cl   (4b)
   0.00000000000000   0.00000000000000   0.00000000000000   F    (4a)
  -0.22600000000000   0.22600000000000   0.22600000000000   Na   (24e)
   0.22600000000000  -0.22600000000000  -0.22600000000000   Na   (24e)
   0.22600000000000  -0.22600000000000   0.22600000000000   Na   (24e)
  -0.22600000000000   0.22600000000000  -0.22600000000000   Na   (24e)
   0.22600000000000   0.22600000000000  -0.22600000000000   Na   (24e)
  -0.22600000000000  -0.22600000000000   0.22600000000000   Na   (24e)
   0.16400000000000   0.16400000000000   0.16400000000000   O    (32f)
   0.16400000000000   0.16400000000000  -0.49200000000000   O    (32f)
   0.16400000000000  -0.49200000000000   0.16400000000000   O    (32f)
  -0.49200000000000   0.16400000000000   0.16400000000000   O    (32f)
  -0.16400000000000  -0.16400000000000   0.49200000000000   O    (32f)
  -0.16400000000000  -0.16400000000000  -0.16400000000000   O    (32f)
  -0.16400000000000   0.49200000000000  -0.16400000000000   O    (32f)
   0.49200000000000  -0.16400000000000  -0.16400000000000   O    (32f)
   0.25000000000000   0.25000000000000   0.25000000000000   S    (8c)
   0.75000000000000   0.75000000000000   0.75000000000000   S    (8c)
```

γ-Ga_2O_3: A11B4_cF120_227_acdf_e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Ga2O3'
_chemical_formula_sum 'Ga1 O4'

loop_
_publ_author_name
 'H. Y. Playford'
 'A. C. Hannon'
 'E. R. Barney'
 'R. I. Walton'
_journal_name_full_name
;
 Chemistry - A European Journal
;
_journal_volume 19
_journal_year 2013
_journal_page_first 2803
_journal_page_last 2813
_publ_Section_title
;
 Structures of Uncharacterised Polymorphs of Gallium Oxide from Total
    ↪ Neutron Diffraction
;

_aflow_title '$\gamma$-Ga$_{2}$O$_{3}$ Structure'
_aflow_proto 'A11B4_cF120_227_acdf_e'
_aflow_params 'a,x_{4},x_{5}'
_aflow_params_values '8.2376,0.2552,0.368'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'cF120'

_symmetry_space_group_name_H-M "F 41/d -3 2/m (origin choice 2)"
_symmetry_Int_Tables_number 227

_cell_length_a     8.23760
_cell_length_b     8.23760
_cell_length_c     8.23760
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y+1/4,-z+1/4
3 -x+1/4,y,-z+1/4
4 -x+1/4,-y+1/4,z
5 y,z,x
```

```
6 y,-z+1/4,-x+1/4
7 -y+1/4,z,-x+1/4
8 -y+1/4,-z+1/4,x
9 z,x,y
10 z,-x+1/4,-y+1/4
11 -z+1/4,x,-y+1/4
12 -z+1/4,-x+1/4,y
13 -y,-x,-z
14 -y,x+1/4,z+1/4
15 y+1/4,-x,z+1/4
16 y+1/4,x+1/4,-z
17 -x,-z,-y
18 -x,z+1/4,y+1/4
19 x+1/4,-z,y+1/4
20 x+1/4,z+1/4,-y
21 -z,-y,-x
22 -z,y+1/4,x+1/4
23 z+1/4,-y,x+1/4
24 z+1/4,y+1/4,-x
25 -x,-y,-z
26 -x,y+1/4,z+1/4
27 x+1/4,-y,z+1/4
28 x+1/4,y+1/4,-z
29 -y,-z,-x
30 -y,z+1/4,x+1/4
31 y+1/4,-z,x+1/4
32 y+1/4,z+1/4,-x
33 -z,-x,-y
34 -z,x+1/4,y+1/4
35 z+1/4,-x,y+1/4
36 z+1/4,x+1/4,-y
37 y,x,z
38 y,-x+1/4,-z+1/4
39 -y+1/4,x,-z+1/4
40 -y+1/4,-x+1/4,z
41 x,z,y
42 x,-z+1/4,-y+1/4
43 -x+1/4,z,-y+1/4
44 -x+1/4,-z+1/4,y
45 z,y,x
46 z,-y+1/4,-x+1/4
47 -z+1/4,y,-x+1/4
48 -z+1/4,-y+1/4,x
49 x,y+1/2,z+1/2
50 x,-y+3/4,-z+3/4
51 -x+1/4,y+1/2,-z+3/4
52 -x+1/4,-y+3/4,z+1/2
53 y,z+1/2,x+1/2
54 y,-z+3/4,-x+3/4
55 -y+1/4,z+1/2,-x+3/4
56 -y+1/4,-z+3/4,x+1/2
57 z,x+1/2,y+1/2
58 z,-x+3/4,-y+3/4
59 -z+1/4,x+1/2,-y+3/4
60 -z+1/4,-x+3/4,y+1/2
61 -y,-x+1/2,-z+1/2
62 -y,x+3/4,z+3/4
63 y+1/4,-x+1/2,z+3/4
64 y+1/4,x+3/4,-z+1/2
65 -x,-z+1/2,-y+1/2
66 -x,z+3/4,y+3/4
67 x+1/4,-z+1/2,y+3/4
68 x+1/4,z+3/4,-y+1/2
69 -z,-y+1/2,-x+1/2
70 -z,y+3/4,x+3/4
71 z+1/4,-y+1/2,x+3/4
72 z+1/4,y+3/4,-x+1/2
73 -x,-y+1/2,-z+1/2
74 -x,y+3/4,z+3/4
75 x+1/4,-y+1/2,z+3/4
76 x+1/4,y+3/4,-z+1/2
77 -y,-z+1/2,-x+1/2
78 -y,z+3/4,x+3/4
79 y+1/4,-z+1/2,x+3/4
80 y+1/4,z+3/4,-x+1/2
81 -z,-x+1/2,-y+1/2
82 -z,x+3/4,y+3/4
83 z+1/4,-x+1/2,y+3/4
84 z+1/4,x+3/4,-y+1/2
85 y,x+1/2,z+1/2
86 y,-x+3/4,-z+3/4
87 -y+1/4,x+1/2,-z+3/4
88 -y+1/4,-x+3/4,z+1/2
89 x,z+1/2,y+1/2
90 x,-z+3/4,-y+3/4
91 -x+1/4,z+1/2,-y+3/4
92 -x+1/4,-z+3/4,y+1/2
93 z,y+1/2,x+1/2
94 z,-y+3/4,-x+3/4
95 -z+1/4,y+1/2,-x+3/4
96 -z+1/4,-y+3/4,x+1/2
97 x+1/2,y,z+1/2
98 x+1/2,-y+1/4,-z+3/4
99 -x+3/4,y,-z+3/4
100 -x+3/4,-y+1/4,z+1/2
101 y+1/2,z,x+1/2
102 y+1/2,-z+1/4,-x+3/4
103 -y+3/4,z,-x+3/4
104 -y+3/4,-z+1/4,x+1/2
105 z+1/2,x,y+1/2
106 z+1/2,-x+1/4,-y+3/4
107 -z+3/4,x,-y+3/4
108 -z+3/4,-x+1/4,y+1/2
109 -y+1/2,-x,-z+1/2
110 -y+1/2,x+1/4,z+3/4
111 y+3/4,-x,z+3/4
112 y+3/4,x+1/4,-z+1/2
113 -x+1/2,-z,-y+1/2
114 -x+1/2,z+1/4,y+3/4
115 x+3/4,-z,y+3/4
116 x+3/4,z+1/4,-y+1/2
117 -z+1/2,-y,-x+1/2
118 -z+1/2,y+1/4,x+3/4
119 z+3/4,-y,x+3/4
120 z+3/4,y+1/4,-x+1/2
121 -x+1/2,-y,-z+1/2
122 -x+1/2,y+1/4,z+3/4
123 x+3/4,-y,z+3/4
124 x+3/4,y+1/4,-z+1/2
125 -y+1/2,-z,-x+1/2
126 -y+1/2,z+1/4,x+3/4
127 y+3/4,-z,x+3/4
128 y+3/4,z+1/4,-x+1/2
129 -z+1/2,-x,-y+1/2
130 -z+1/2,x+1/4,y+3/4
131 z+3/4,-x,y+3/4
132 z+3/4,x+1/4,-y+1/2
133 y+1/2,x,z+1/2
134 y+1/2,-x+1/4,-z+3/4
135 -y+3/4,x,-z+3/4
136 -y+3/4,-x+1/4,z+1/2
137 x+1/2,z,y+1/2
138 x+1/2,-z+1/4,-y+3/4
139 -x+3/4,z,-y+3/4
140 -x+3/4,-z+1/4,y+1/2
141 z+1/2,y,x+1/2
142 z+1/2,-y+1/4,-x+3/4
143 -z+3/4,y,-x+3/4
144 -z+3/4,-y+1/4,x+1/2
145 x+1/2,y+1/2,z
146 x+1/2,-y+3/4,-z+1/4
147 -x+3/4,y+1/2,-z+1/4
148 -x+3/4,-y+3/4,z
149 y+1/2,z+1/2,x
150 y+1/2,-z+3/4,-x+1/4
151 -y+3/4,z+1/2,-x+1/4
152 -y+3/4,-z+3/4,x
153 z+1/2,x+1/2,y
154 z+1/2,-x+3/4,-y+1/4
155 -z+3/4,x+1/2,-y+1/4
156 -z+3/4,-x+3/4,y
157 -y+1/2,-x+1/2,-z
158 -y+1/2,x+3/4,z+1/4
159 y+3/4,-x+1/2,z+1/4
160 y+3/4,x+3/4,-z
161 -x+1/2,-z+1/2,-y
162 -x+1/2,z+3/4,y+1/4
163 x+3/4,-z+1/2,y+1/4
164 x+3/4,z+3/4,-y
165 -z+1/2,-y+1/2,-x
166 -z+1/2,y+3/4,x+1/4
167 z+3/4,-y+1/2,x+1/4
168 z+3/4,y+3/4,-x
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170 -x+1/2,y+3/4,z+1/4
171 x+3/4,-y+1/2,z+1/4
172 x+3/4,y+3/4,-z
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174 -y+1/2,z+3/4,x+1/4
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176 y+3/4,z+3/4,-x
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178 -z+1/2,x+3/4,y+1/4
179 z+3/4,-x+1/2,y+1/4
180 z+3/4,x+3/4,-y
181 y+1/2,x+1/2,z
182 y+1/2,-x+3/4,-z+1/4
183 -y+3/4,x+1/2,-z+1/4
184 -y+3/4,-x+3/4,z
185 x+1/2,z+1/2,y
186 x+1/2,-z+3/4,-y+1/4
187 -x+3/4,z+1/2,-y+1/4
188 -x+3/4,-z+3/4,y
189 z+1/2,y+1/2,x
190 z+1/2,-y+3/4,-x+1/4
191 -z+3/4,y+1/2,-x+1/4
192 -z+3/4,-y+3/4,x

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Ga3 Ga 16 d 0.50000 0.50000 0.50000 1.00000
O1 O 32 e 0.25520 0.25520 0.25520 1.00000
Ga4 Ga 48 f 0.36800 0.12500 0.12500 1.00000
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γ-Ga_2O_3: A11B4_cF120_227_acdf_e - POSCAR

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    ↪ Playford et al., Chem. Eur. J. 19, 2803-2813 (2013)
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    4.11880000000000    4.11880000000000    0.00000000000000
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     22      8
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    0.87500000000000    0.87500000000000    0.87500000000000   Ga    (8a)
    0.00000000000000    0.00000000000000    0.00000000000000   Ga   (16c)
    0.00000000000000    0.00000000000000    0.50000000000000   Ga   (16c)
    0.00000000000000    0.50000000000000    0.00000000000000   Ga   (16c)
    0.50000000000000    0.00000000000000    0.00000000000000   Ga   (16c)
    0.50000000000000    0.50000000000000    0.50000000000000   Ga   (16d)
    0.50000000000000    0.50000000000000    0.00000000000000   Ga   (16d)
    0.50000000000000    0.00000000000000    0.50000000000000   Ga   (16d)
    0.00000000000000    0.50000000000000    0.50000000000000   Ga   (16d)
   -0.11800000000000    0.36800000000000    0.36800000000000   Ga   (48f)
    0.36800000000000   -0.11800000000000   -0.11800000000000   Ga   (48f)
    0.36800000000000   -0.11800000000000    0.36800000000000   Ga   (48f)
   -0.11800000000000    0.36800000000000   -0.11800000000000   Ga   (48f)
    0.36800000000000    0.36800000000000   -0.11800000000000   Ga   (48f)
   -0.11800000000000   -0.11800000000000    0.36800000000000   Ga   (48f)
    1.11800000000000   -0.36800000000000    1.11800000000000   Ga   (48f)
   -0.36800000000000    1.11800000000000   -0.36800000000000   Ga   (48f)
   -0.36800000000000    1.11800000000000    1.11800000000000   Ga   (48f)
    1.11800000000000   -0.36800000000000   -0.36800000000000   Ga   (48f)
   -0.36800000000000   -0.36800000000000    1.11800000000000   Ga   (48f)
    1.11800000000000    1.11800000000000   -0.36800000000000   Ga   (48f)
    0.25520000000000    0.25520000000000    0.25520000000000    O   (32e)
    0.25520000000000    0.25520000000000   -0.26560000000000    O   (32e)
    0.25520000000000   -0.26560000000000    0.25520000000000    O   (32e)
   -0.26560000000000    0.25520000000000    0.25520000000000    O   (32e)
   -0.25520000000000   -0.25520000000000    1.26560000000000    O   (32e)
   -0.25520000000000   -0.25520000000000   -0.25520000000000    O   (32e)
   -0.25520000000000    1.26560000000000   -0.25520000000000    O   (32e)
    1.26560000000000   -0.25520000000000   -0.25520000000000    O   (32e)
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Predicted Li_2MgH_{16} High-Temperature Superconductor (250 GPa): A16B2C_cF152_227_eg_d_a - CIF

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_chemical_formula_sum 'H16 Li2 Mg'

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 'J. Lv'
 'Y. Xie'
 'H. Liu'
 'Y. Ma'
_journal_name_full_name
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 Physical Review Letters
;
_journal_volume 123
_journal_year 2019
_journal_page_first 097001
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 Route to a Superconducting Phase above Room Temperature in
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_aflow_Pearson 'cF152'

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_cell_length_b     6.71851
_cell_length_c     6.71851
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

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_space_group_symop_operation_xyz
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3 -x+1/4,y,-z+1/4
4 -x+1/4,-y+1/4,z
5 y,z,x
6 y,-z+1/4,-x+1/4
7 -y+1/4,z,-x+1/4
8 -y+1/4,-z+1/4,x
9 z,x,y
10 z,-x+1/4,-y+1/4
11 -z+1/4,x,-y+1/4
12 -z+1/4,-x+1/4,y
13 -y,-x,-z
14 -y,x+1/4,z+1/4
15 y+1/4,-x,z+1/4
16 y+1/4,x+1/4,-z
17 -x,-z,-y
18 -x,z+1/4,y+1/4
19 x+1/4,-z,y+1/4
20 x+1/4,z+1/4,-y
21 -z,-y,-x
22 -z,y+1/4,x+1/4
23 z+1/4,-y,x+1/4
24 z+1/4,y+1/4,-x
25 -x,-y,-z
26 -x,y+1/4,z+1/4
27 x+1/4,-y,z+1/4
28 x+1/4,y+1/4,-z
29 -y,-z,-x
30 -y,z+1/4,x+1/4
31 y+1/4,-z,x+1/4
32 y+1/4,z+1/4,-x
33 -z,-x,-y
34 -z,x+1/4,y+1/4
35 z+1/4,-x,y+1/4
36 z+1/4,x+1/4,-y
37 y,x,z
38 y,-x+1/4,-z+1/4
39 -y+1/4,x,-z+1/4
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43 -x+1/4,z,-y+1/4
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45 z,y,x
46 z,-y+1/4,-x+1/4
47 -z+1/4,y,-x+1/4
48 -z+1/4,-y+1/4,x
49 x,y+1/2,z+1/2
50 x,-y+3/4,-z+3/4
51 -x+1/4,y+1/2,-z+3/4
52 -x+1/4,-y+3/4,z+1/2
53 y,z+1/2,x+1/2
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55 -y+1/4,z+1/2,-x+3/4
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68 x+1/4,z+3/4,-y+1/2
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70 -z,y+3/4,x+3/4
71 z+1/4,-y+1/2,x+3/4
72 z+1/4,y+3/4,-x+1/2
73 -x,-y+1/2,-z+1/2
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84 z+1/4,x+3/4,-y+1/2
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90 x,-z+3/4,-y+3/4
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93 z,y+1/2,x+1/2
94 z,-y+3/4,-x+3/4
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97 x+1/2,y,z+1/2
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99 -x+3/4,y,-z+3/4
100 -x+3/4,-y+1/4,z+1/2
101 y+1/2,z,x+1/2
102 y+1/2,-z+1/4,-x+3/4
103 -y+3/4,z,-x+3/4
104 -y+3/4,-z+1/4,x+1/2
105 z+1/2,x,y+1/2
106 z+1/2,-x+1/4,-y+3/4
107 -z+3/4,x,-y+3/4
108 -z+3/4,-x+1/4,y+1/2
109 -y+1/2,-x,-z+1/2
110 -y+1/2,x+1/4,z+3/4
111 y+3/4,-x,z+3/4
112 y+3/4,x+1/4,-z+1/2
113 -x+1/2,-z,-y+1/2
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115 x+3/4,-z,y+3/4
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118 -z+1/2,y+1/4,x+3/4
119 z+3/4,-y,x+3/4
120 z+3/4,y+1/4,-x+1/2
121 -x+1/2,-y,-z+1/2
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123 x+3/4,-y,z+3/4
```

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128 y+3/4,z+1/4,-x+1/2
129 -z+1/2,-x,-y+1/2
130 -z+1/2,x+1/4,y+3/4
131 z+3/4,-x,y+3/4
132 z+3/4,x+1/4,-y+1/2
133 y+1/2,x,z+1/2
134 y+1/2,-x+1/4,-z+3/4
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185 x+1/2,z+1/2,y
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189 z+1/2,y+1/2,x
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Li1 Li 16 d 0.50000 0.50000 0.50000 1.00000
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H2 H 96 g 0.56034 0.56034 0.87060 1.00000
```

Predicted Li_2MgH_{16} High-Temperature Superconductor (250 GPa): A16B2C_cF152_227_eg_d_a - POSCAR

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    ↪ & H16Li2Mg & Y. Sun et al., Phys. Rev. Lett. 123, 097001(2019)
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   0.000000000000000   3.359255000000000   3.359255000000000
   3.359255000000000   0.000000000000000   3.359255000000000
   3.359255000000000   3.359255000000000   0.000000000000000
     H     Li     Mg
    32      4      2
Direct
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   0.291500000000000   0.291500000000000  -0.374500000000000   H   (32e)
   0.291500000000000  -0.374500000000000   0.291500000000000   H   (32e)
  -0.374500000000000   0.291500000000000   0.291500000000000   H   (32e)
  -0.291500000000000  -0.291500000000000   1.374500000000000   H   (32e)
  -0.291500000000000  -0.291500000000000  -0.291500000000000   H   (32e)
  -0.291500000000000   1.374500000000000  -0.291500000000000   H   (32e)
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   0.870600000000000   0.870600000000000  -1.491280000000000   H   (96g)
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   0.870600000000000   0.250080000000000   0.870600000000000   H   (96g)
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  -1.491280000000000   0.870600000000000   0.250080000000000   H   (96g)
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  -0.870600000000000  -0.870600000000000  -0.250080000000000   H   (96g)
  -0.250080000000000   2.491280000000000  -0.870600000000000   H   (96g)
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  -0.250080000000000  -0.870600000000000   2.491280000000000   H   (96g)
   2.491280000000000  -0.870600000000000  -0.250080000000000   H   (96g)
  -0.870600000000000  -0.250080000000000  -0.870600000000000   H   (96g)
  -0.870600000000000   2.491280000000000  -0.870600000000000   H   (96g)
  -0.870600000000000  -0.250080000000000   2.491280000000000   H   (96g)
  -0.870600000000000   2.491280000000000  -0.250080000000000   H   (96g)
   2.491280000000000  -0.870600000000000  -0.870600000000000   H   (96g)
  -0.250080000000000  -0.870600000000000  -0.870600000000000   H   (96g)
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   0.500000000000000   0.500000000000000   0.000000000000000   Li  (16d)
   0.500000000000000   0.000000000000000   0.500000000000000   Li  (16d)
   0.000000000000000   0.500000000000000   0.500000000000000   Li  (16d)
   0.125000000000000   0.125000000000000   0.125000000000000   Mg   (8a)
   0.875000000000000   0.875000000000000   0.875000000000000   Mg   (8a)
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$Mg_3Cr_2Al_{18}$: A18B2C3_cF184_227_fg_d_ac - CIF

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     ↪ $Al$_{18}$
;

# Found in The Crystal Structure of the Intermetallic Compound ZrZn$_{22
    ↪ }$, 1961

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_aflow_Strukturbericht 'None'
_aflow_Pearson 'cF184'

_symmetry_space_group_name_H-M "F 41/d -3 2/m (origin choice 2)"
_symmetry_Int_Tables_number 227

_cell_length_a    14.53000
_cell_length_b    14.53000
_cell_length_c    14.53000
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y+1/4,-z+1/4
3 -x+1/4,y,-z+1/4
4 -x+1/4,-y+1/4,z
5 y,z,x
6 y,-z+1/4,-x+1/4
7 -y+1/4,z,-x+1/4
8 -y+1/4,-z+1/4,x
9 z,x,y
10 z,-x+1/4,-y+1/4
11 -z+1/4,x,-y+1/4
12 -z+1/4,-x+1/4,y
13 -y,-x,-z
14 -y,x+1/4,z+1/4
15 y+1/4,-x,z+1/4
16 y+1/4,x+1/4,-z
17 -x,-z,-y
18 -x,z+1/4,y+1/4
19 x+1/4,-z,y+1/4
20 x+1/4,z+1/4,-y
21 -z,-y,-x
22 -z,y+1/4,x+1/4
23 z+1/4,-y,x+1/4
24 z+1/4,y+1/4,-x
25 -x,-y,-z
26 -x,y+1/4,z+1/4
27 x+1/4,-y,z+1/4
28 x+1/4,y+1/4,-z
29 -y,-z,-x
30 -y,z+1/4,x+1/4
31 y+1/4,-z,x+1/4
32 y+1/4,z+1/4,-x
```

```
33 -z,-x,-y
34 -z,x+1/4,y+1/4
35 z+1/4,-x,y+1/4
36 z+1/4,x+1/4,-y
37 y,x,z
38 y,-x+1/4,-z+1/4
39 -y+1/4,x,-z+1/4
40 -y+1/4,-x+1/4,z
41 x,z,y
42 x,-z+1/4,-y+1/4
43 -x+1/4,z,-y+1/4
44 -x+1/4,-z+1/4,y
45 z,y,x
46 z,-y+1/4,-x+1/4
47 -z+1/4,y,-x+1/4
48 -z+1/4,-y+1/4,x
49 x,y+1/2,z+1/2
50 x,-y+3/4,-z+3/4
51 -x+1/4,y+1/2,-z+3/4
52 -x+1/4,-y+3/4,z+1/2
53 y,z+1/2,x+1/2
54 y,-z+3/4,-x+3/4
55 -y+1/4,z+1/2,-x+3/4
56 -y+1/4,-z+3/4,x+1/2
57 z,x+1/2,y+1/2
58 z,-x+3/4,-y+3/4
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60 -z+1/4,-x+3/4,y+1/2
61 -y,-x+1/2,-z+1/2
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64 y+1/4,x+3/4,-z+1/2
65 -x,-z+1/2,-y+1/2
66 -x,z+3/4,y+3/4
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68 x+1/4,z+3/4,-y+1/2
69 -z,-y+1/2,-x+1/2
70 -z,y+3/4,x+3/4
71 z+1/4,-y+1/2,x+3/4
72 z+1/4,y+3/4,-x+1/2
73 -x,-y+1/2,-z+1/2
74 -x,y+3/4,z+3/4
75 x+1/4,-y+1/2,z+3/4
76 x+1/4,y+3/4,-z+1/2
77 -y,-z+1/2,-x+1/2
78 -y,z+3/4,x+3/4
79 y+1/4,-z+1/2,x+3/4
80 y+1/4,z+3/4,-x+1/2
81 -z,-x+1/2,-y+1/2
82 -z,x+3/4,y+3/4
83 z+1/4,-x+1/2,y+3/4
84 z+1/4,x+3/4,-y+1/2
85 y,x+1/2,z+1/2
86 y,-x+3/4,-z+3/4
87 -y+1/4,x+1/2,-z+3/4
88 -y+1/4,-x+3/4,z+1/2
89 x,z+1/2,y+1/2
90 x,-z+3/4,-y+3/4
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92 -x+1/4,-z+3/4,y+1/2
93 z,y+1/2,x+1/2
94 z,-y+3/4,-x+3/4
95 -z+1/4,y+1/2,-x+3/4
96 -z+1/4,-y+3/4,x+1/2
97 x+1/2,y,z+1/2
98 x+1/2,-y+1/4,-z+3/4
99 -x+3/4,y,-z+3/4
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101 y+1/2,z,x+1/2
102 y+1/2,-z+1/4,-x+3/4
103 -y+3/4,z,-x+3/4
104 -y+3/4,-z+1/4,x+1/2
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107 -z+3/4,x,-y+3/4
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109 -y+1/2,-x,-z+1/2
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113 -x+1/2,-z,-y+1/2
114 -x+1/2,z+1/4,y+3/4
115 x+3/4,-z,y+3/4
116 x+3/4,z+1/4,-y+1/2
117 -z+1/2,-y,-x+1/2
118 -z+1/2,y+1/4,x+3/4
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120 z+3/4,y+1/4,-x+1/2
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128 y+3/4,z+1/4,-x+1/2
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130 -z+1/2,x+1/4,y+3/4
131 z+3/4,-x,y+3/4
132 z+3/4,x+1/4,-y+1/2
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138 x+1/2,-z+1/4,-y+3/4
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141 z+1/2,y,x+1/2
142 z+1/2,-y+1/4,-x+3/4
143 -z+3/4,y,-x+3/4
144 -z+3/4,-y+1/4,x+1/2
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146 x+1/2,-y+3/4,-z+1/4
147 -x+3/4,y+1/2,-z+1/4
148 -x+3/4,-y+3/4,z
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150 y+1/2,-z+3/4,-x+1/4
151 -y+3/4,z+1/2,-x+1/4
152 -y+3/4,-z+3/4,x
153 z+1/2,x+1/2,y
154 z+1/2,-x+3/4,-y+1/4
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157 -y+1/2,-x+1/2,-z
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160 y+3/4,x+3/4,-z
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162 -x+1/2,z+3/4,y+1/4
163 x+3/4,-z+1/2,y+1/4
164 x+3/4,z+3/4,-y
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166 -z+1/2,y+3/4,x+1/4
167 z+3/4,-y+1/2,x+1/4
168 z+3/4,y+3/4,-x
169 -x+1/2,-y+1/2,-z
170 -x+1/2,y+3/4,z+1/4
171 x+3/4,-y+1/2,z+1/4
172 x+3/4,y+3/4,-z
173 -y+1/2,-z+1/2,-x
174 -y+1/2,z+3/4,x+1/4
175 y+3/4,-z+1/2,x+1/4
176 y+3/4,z+3/4,-x
177 -z+1/2,-x+1/2,-y
178 -z+1/2,x+3/4,y+1/4
179 z+3/4,-x+1/2,y+1/4
180 z+3/4,x+3/4,-y
181 y+1/2,x+1/2,z
182 y+1/2,-x+3/4,-z+1/4
183 -y+3/4,x+1/2,-z+1/4
184 -y+3/4,-x+3/4,z
185 x+1/2,z+1/2,y
186 x+1/2,-z+3/4,-y+1/4
187 -x+3/4,z+1/2,-y+1/4
188 -x+3/4,-z+3/4,y
189 z+1/2,y+1/2,x
190 z+1/2,-y+3/4,-x+1/4
191 -z+3/4,y+1/2,-x+1/4
192 -z+3/4,-y+3/4,x

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_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
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Mg2 Mg 16 c 0.00000 0.00000 0.00000 1.00000
Cr1 Cr 16 d 0.50000 0.50000 0.50000 1.00000
Al1 Al 48 f 0.76570 0.12500 0.12500 1.00000
Al2 Al 96 g 0.55840 0.55840 0.32520 1.00000
```

$Mg_3Cr_2Al_{18}$: A18B2C3_cF184_227_fg_d_ac - POSCAR

```
A18B2C3_cF184_227_fg_d_ac & a,x4,x5,z5 --params=14.53,0.7657,0.5584,
    ↪ 0.3252 & Fd-3m O_{h}^{7} #227 (acdfg) & cF184 & None &
    ↪ Al18Cr2Mg3 & Al18Cr2Mg3 & S. Samson, (1958)
   1.00000000000000
   0.00000000000000   7.26500000000000   7.26500000000000
   7.26500000000000   0.00000000000000   7.26500000000000
   7.26500000000000   7.26500000000000   0.00000000000000
    Al   Cr   Mg
    36    4    6
Direct
  -0.51570000000000   0.76570000000000   0.76570000000000   Al   (48f)
   0.76570000000000  -0.51570000000000  -0.51570000000000   Al   (48f)
   0.76570000000000  -0.51570000000000   0.76570000000000   Al   (48f)
  -0.51570000000000   0.76570000000000  -0.51570000000000   Al   (48f)
   0.76570000000000   0.76570000000000  -0.51570000000000   Al   (48f)
  -0.51570000000000  -0.51570000000000   0.76570000000000   Al   (48f)
   1.51570000000000  -0.76570000000000   1.51570000000000   Al   (48f)
  -0.76570000000000   1.51570000000000  -0.76570000000000   Al   (48f)
  -0.76570000000000   1.51570000000000   1.51570000000000   Al   (48f)
   1.51570000000000  -0.76570000000000  -0.76570000000000   Al   (48f)
  -0.76570000000000  -0.76570000000000   1.51570000000000   Al   (48f)
   1.51570000000000   1.51570000000000  -0.76570000000000   Al   (48f)
   0.32520000000000   0.32520000000000   0.79160000000000   Al   (96g)
   0.32520000000000   0.32520000000000  -0.94200000000000   Al   (96g)
   0.79160000000000  -0.94200000000000   0.32520000000000   Al   (96g)
  -0.94200000000000   0.79160000000000   0.32520000000000   Al   (96g)
   0.79160000000000   0.32520000000000   0.32520000000000   Al   (96g)
  -0.94200000000000   0.32520000000000   0.32520000000000   Al   (96g)
   0.32520000000000   0.79160000000000  -0.94200000000000   Al   (96g)
   0.32520000000000  -0.94200000000000   0.79160000000000   Al   (96g)
   0.32520000000000   0.79160000000000   0.32520000000000   Al   (96g)
   0.32520000000000  -0.94200000000000   0.32520000000000   Al   (96g)
```

−0.94200000000000	0.32520000000000	0.79160000000000	Al	(96g)
0.79160000000000	0.32520000000000	−0.94200000000000	Al	(96g)
−0.32520000000000	−0.32520000000000	1.94200000000000	Al	(96g)
−0.32520000000000	−0.32520000000000	−0.79160000000000	Al	(96g)
−0.79160000000000	1.94200000000000	−0.32520000000000	Al	(96g)
1.94200000000000	−0.79160000000000	−0.32520000000000	Al	(96g)
−0.79160000000000	−0.32520000000000	1.94200000000000	Al	(96g)
1.94200000000000	−0.32520000000000	−0.79160000000000	Al	(96g)
−0.32520000000000	−0.79160000000000	−0.32520000000000	Al	(96g)
−0.32520000000000	1.94200000000000	−0.32520000000000	Al	(96g)
−0.32520000000000	−0.79160000000000	1.94200000000000	Al	(96g)
−0.32520000000000	1.94200000000000	−0.79160000000000	Al	(96g)
1.94200000000000	−0.32520000000000	−0.32520000000000	Al	(96g)
−0.79160000000000	−0.32520000000000	−0.32520000000000	Al	(96g)
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0.50000000000000	0.50000000000000	0.00000000000000	Cr	(16d)
0.50000000000000	0.00000000000000	0.50000000000000	Cr	(16d)
0.00000000000000	0.50000000000000	0.50000000000000	Cr	(16d)
0.12500000000000	0.12500000000000	0.12500000000000	Mg	(8a)
0.87500000000000	0.87500000000000	0.87500000000000	Mg	(8a)
0.00000000000000	0.00000000000000	0.00000000000000	Mg	(16c)
0.00000000000000	0.00000000000000	0.50000000000000	Mg	(16c)
0.00000000000000	0.50000000000000	0.00000000000000	Mg	(16c)
0.50000000000000	0.00000000000000	0.00000000000000	Mg	(16c)

$Zn_{22}Zr$: A22B_cF184_227_cdfg_a - CIF

```
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_audit_creation_method FINDSYM

_chemical_name_mineral 'Zn22Zr'
_chemical_formula_sum 'Zn22 Zr'

loop_
_publ_author_name
 'S. Samson'
_journal_year 1961
_publ_Section_title
;
 The Crystal Structure of the Intermetallic Compound ZrZn$_{22}$
;

# Found in Binary Alloy Phase Diagrams, {Hf-Re to Zn-Zr}, 1990

_aflow_title 'Zn$_{22}$Zr Structure'
_aflow_proto 'A22B_cF184_227_cdfg_a'
_aflow_params 'a,x_{4},x_{5},z_{5}'
_aflow_params_values '14.101,0.765,0.562,0.3195'
_aflow_Strukturbericht 'None'
_aflow_Pearson 'cF184'

_symmetry_space_group_name_H-M "F 41/d -3 2/m (origin choice 2)"
_symmetry_Int_Tables_number 227

_cell_length_a     14.10100
_cell_length_b     14.10100
_cell_length_c     14.10100
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y+1/4,-z+1/4
3 -x+1/4,y,-z+1/4
4 -x+1/4,-y+1/4,z
5 y,z,x
6 y,-z+1/4,-x+1/4
7 -y+1/4,z,-x+1/4
8 -y+1/4,-z+1/4,x
9 z,x,y
10 z,-x+1/4,-y+1/4
11 -z+1/4,x,-y+1/4
12 -z+1/4,-x+1/4,y
13 -y,-x,-z
14 -y,x+1/4,z+1/4
15 y+1/4,-x,z+1/4
16 y+1/4,x+1/4,-z
17 -x,-z,-y
18 -x,z+1/4,y+1/4
19 x+1/4,-z,y+1/4
20 x+1/4,z+1/4,-y
21 -z,-y,-x
22 -z,y+1/4,x+1/4
23 z+1/4,-y,x+1/4
24 z+1/4,y+1/4,-x
25 -x,-y,-z
26 -x,y+1/4,z+1/4
27 x+1/4,-y,z+1/4
28 x+1/4,y+1/4,-z
29 -y,-z,-x
30 -y,z+1/4,x+1/4
31 y+1/4,-z,x+1/4
32 y+1/4,z+1/4,-x
33 -z,-x,-y
34 -z,x+1/4,y+1/4
35 z+1/4,-x,y+1/4
36 z+1/4,x+1/4,-y
37 y,x,z
38 y,-x+1/4,-z+1/4
39 -y+1/4,x,-z+1/4
40 -y+1/4,-x+1/4,z
41 x,z,y
42 x,-z+1/4,-y+1/4
43 -x+1/4,z,-y+1/4
44 -x+1/4,-z+1/4,y
45 z,y,x
46 z,-y+1/4,-x+1/4
47 -z+1/4,y,-x+1/4
48 -z+1/4,-y+1/4,x
49 x,y+1/2,z+1/2
50 x,-y+3/4,-z+3/4
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53 y,z+1/2,x+1/2
54 y,-z+3/4,-x+3/4
55 -y+1/4,z+1/2,-x+3/4
56 -y+1/4,-z+3/4,x+1/2
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61 -y,-x+1/2,-z+1/2
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73 -x,-y+1/2,-z+1/2
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83 z+1/4,-x+1/2,y+3/4
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85 y,x+1/2,z+1/2
86 y,-x+3/4,-z+3/4
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89 x,z+1/2,y+1/2
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95 -z+1/4,y+1/2,-x+3/4
96 -z+1/4,-y+3/4,x+1/2
97 x+1/2,y,z+1/2
98 x+1/2,-y+1/4,-z+3/4
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102 y+1/2,-z+1/4,-x+3/4
103 -y+3/4,z,-x+3/4
104 -y+3/4,-z+1/4,x+1/2
105 z+1/2,x,y+1/2
106 z+1/2,-x+1/4,-y+3/4
107 -z+3/4,x,-y+3/4
108 -z+3/4,-x+1/4,y+1/2
109 -y+1/2,-x,-z+1/2
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111 y+3/4,-x,z+3/4
112 y+3/4,x+1/4,-z+1/2
113 -x+1/2,-z,-y+1/2
114 -x+1/2,z+1/4,y+3/4
115 x+3/4,-z,y+3/4
116 x+3/4,z+1/4,-y+1/2
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133 y+1/2,x,z+1/2
134 y+1/2,-x+1/4,-z+3/4
135 -y+3/4,x,-z+3/4
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138 x+1/2,-z+1/4,-y+3/4
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140 -x+3/4,-z+1/4,y+1/2
141 z+1/2,y,x+1/2
142 z+1/2,-y+1/4,-x+3/4
143 -z+3/4,y,-x+3/4
144 -z+3/4,-y+1/4,x+1/2
145 x+1/2,y+1/2,z
```

```
146 x+1/2,-y+3/4,-z+1/4
147 -x+3/4,y+1/2,-z+1/4
148 -x+3/4,-y+3/4,z
149 y+1/2,z+1/2,x
150 y+1/2,-z+3/4,-x+1/4
151 -y+3/4,z+1/2,-x+1/4
152 -y+3/4,-z+3/4,x
153 z+1/2,x+1/2,y
154 z+1/2,-x+3/4,-y+1/4
155 -z+3/4,x+1/2,-y+1/4
156 -z+3/4,-x+3/4,y
157 -y+1/2,-x+1/2,-z
158 -y+1/2,x+3/4,z+1/4
159 y+3/4,-x+1/2,z+1/4
160 y+3/4,x+3/4,-z
161 -x+1/2,-z+1/2,-y
162 -x+1/2,z+3/4,y+1/4
163 x+3/4,-z+1/2,y+1/4
164 x+3/4,z+3/4,-y
165 -z+1/2,-y+1/2,-x
166 -z+1/2,y+3/4,x+1/4
167 z+3/4,-y+1/2,x+1/4
168 z+3/4,y+3/4,-x
169 -x+1/2,-y+1/2,-z
170 -x+1/2,y+3/4,z+1/4
171 x+3/4,-y+1/2,z+1/4
172 x+3/4,y+3/4,-z
173 -y+1/2,-z+1/2,-x
174 -y+1/2,z+3/4,x+1/4
175 y+3/4,-z+1/2,x+1/4
176 y+3/4,z+3/4,-x
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178 -z+1/2,x+3/4,y+1/4
179 z+3/4,-x+1/2,y+1/4
180 z+3/4,x+3/4,-y
181 y+1/2,x+1/2,z
182 y+1/2,-x+3/4,-z+1/4
183 -y+3/4,x+1/2,-z+1/4
184 -y+3/4,-x+3/4,z
185 x+1/2,z+1/2,y
186 x+1/2,-z+3/4,-y+1/4
187 -x+3/4,z+1/2,-y+1/4
188 -x+3/4,-z+3/4,y
189 z+1/2,y+1/2,x
190 z+1/2,-y+3/4,-x+1/4
191 -z+3/4,y+1/2,-x+1/4
192 -z+3/4,-y+3/4,x

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Zn1 Zn 16 c 0.00000 0.00000 0.00000 1.00000
Zn2 Zn 16 d 0.50000 0.50000 0.50000 1.00000
Zn3 Zn 48 f 0.76500 0.12500 0.12500 1.00000
Zn4 Zn 96 g 0.56200 0.56200 0.31950 1.00000
```

$Zn_{22}Zr$: A22B_cF184_227_cdfg_a - POSCAR

```
A22B_cF184_227_cdfg_a & a,x4,x5,z5 --params=14.101,0.765,0.562,0.3195 &
    ↪ Fd-3m O_{h}^{7} #227 (acdfg) & cF184 & None & Zn22Zr & Zn22Zr &
    ↪ S. Samson, (1961)
   1.00000000000000
   0.00000000000000   7.05050000000000   7.05050000000000
   7.05050000000000   0.00000000000000   7.05050000000000
   7.05050000000000   7.05050000000000   0.00000000000000
    Zn    Zr
    44     2
Direct
   0.00000000000000   0.00000000000000   0.00000000000000   Zn   (16c)
   0.00000000000000   0.00000000000000   0.50000000000000   Zn   (16c)
   0.00000000000000   0.50000000000000   0.00000000000000   Zn   (16c)
   0.50000000000000   0.00000000000000   0.00000000000000   Zn   (16c)
   0.50000000000000   0.50000000000000   0.50000000000000   Zn   (16d)
   0.50000000000000   0.50000000000000   0.00000000000000   Zn   (16d)
   0.50000000000000   0.00000000000000   0.50000000000000   Zn   (16d)
   0.00000000000000   0.50000000000000   0.50000000000000   Zn   (16d)
  -0.51500000000000   0.76500000000000   0.76500000000000   Zn   (48f)
   0.76500000000000  -0.51500000000000  -0.51500000000000   Zn   (48f)
   0.76500000000000  -0.51500000000000   0.76500000000000   Zn   (48f)
  -0.51500000000000   0.76500000000000  -0.51500000000000   Zn   (48f)
   0.76500000000000   0.76500000000000  -0.51500000000000   Zn   (48f)
  -0.51500000000000  -0.51500000000000   0.76500000000000   Zn   (48f)
   1.51500000000000  -0.76500000000000   1.51500000000000   Zn   (48f)
  -0.76500000000000   1.51500000000000  -0.76500000000000   Zn   (48f)
  -0.76500000000000   1.51500000000000   1.51500000000000   Zn   (48f)
   1.51500000000000  -0.76500000000000  -0.76500000000000   Zn   (48f)
  -0.76500000000000  -0.76500000000000   1.51500000000000   Zn   (48f)
   1.51500000000000   1.51500000000000  -0.76500000000000   Zn   (48f)
   0.31950000000000   0.31950000000000   0.80450000000000   Zn   (96g)
   0.31950000000000   0.31950000000000  -0.94350000000000   Zn   (96g)
   0.80450000000000  -0.94350000000000   0.31950000000000   Zn   (96g)
  -0.94350000000000   0.80450000000000   0.31950000000000   Zn   (96g)
   0.80450000000000   0.31950000000000   0.31950000000000   Zn   (96g)
  -0.94350000000000   0.31950000000000   0.31950000000000   Zn   (96g)
   0.31950000000000   0.80450000000000  -0.94350000000000   Zn   (96g)
   0.31950000000000  -0.94350000000000   0.80450000000000   Zn   (96g)
   0.31950000000000   0.80450000000000   0.31950000000000   Zn   (96g)
   0.31950000000000  -0.94350000000000   0.31950000000000   Zn   (96g)
  -0.94350000000000   0.31950000000000   0.80450000000000   Zn   (96g)
   0.80450000000000   0.31950000000000  -0.94350000000000   Zn   (96g)
  -0.31950000000000  -0.31950000000000   1.94350000000000   Zn   (96g)
  -0.31950000000000   1.94350000000000  -0.31950000000000   Zn   (96g)
  -0.80450000000000  -0.31950000000000   1.94350000000000   Zn   (96g)
  -0.80450000000000   1.94350000000000  -0.31950000000000   Zn   (96g)
   1.94350000000000  -0.80450000000000  -0.31950000000000   Zn   (96g)
   1.94350000000000  -0.31950000000000  -0.80450000000000   Zn   (96g)
  -0.31950000000000  -0.80450000000000  -0.31950000000000   Zn   (96g)
  -0.31950000000000   1.94350000000000  -0.31950000000000   Zn   (96g)
  -0.31950000000000  -0.80450000000000   1.94350000000000   Zn   (96g)
  -0.31950000000000   1.94350000000000  -0.80450000000000   Zn   (96g)
   1.94350000000000  -0.31950000000000  -0.31950000000000   Zn   (96g)
  -0.80450000000000  -0.31950000000000  -0.31950000000000   Zn   (96g)
   0.12500000000000   0.12500000000000   0.12500000000000   Zr    (8a)
   0.87500000000000   0.87500000000000   0.87500000000000   Zr    (8a)
```

$H_3PW_{12}O_{40}{\cdot}29H_2O$ ($H4_{21}$): A29B40CD12_cF656_227_ae2fg_e3g_b_g - CIF

```
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_chemical_formula_sum '(H2O)29 O40 P W12'

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 'A. J. Bradley'
 'J. W. Illingworth'
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;
 Proceedings of the Royal Society London A
;
_journal_volume 157
_journal_year 1936
_journal_page_first 113
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;
 The Crystal Structure of H$_{3}$PW$_{12}$O$_{40}$$\cdot$29H$_{2}$O
;

# Found in Strukturbericht Band IV 1936, 1938

_aflow_title 'H$_{3}$PW$_{12}$O$_{40}$$\cdot$29H$_{2}$O ($H4_{21}$)
    ↪ Structure'
_aflow_proto 'A29B40CD12_cF656_227_ae2fg_e3g_b_g'
_aflow_params 'a,x_{3},x_{4},x_{5},x_{6},x_{7},z_{7},x_{8},z_{8},x_{9},
    ↪ z_{9},x_{10},z_{10},x_{11},z_{11}'
_aflow_params_values '23.28001,-0.05284,0.41753,-0.01933,0.85395,0.71864
    ↪ ,0.85825,0.83333,0.25301,-0.061,0.52706,0.5378,0.3707,-0.01783,
    ↪ 0.38144'
_aflow_Strukturbericht '$H4_{21}$'
_aflow_Pearson 'cF656'

_symmetry_space_group_name_H-M "F 41/d -3 2/m (origin choice 2)"
_symmetry_Int_Tables_number 227

_cell_length_a    23.28001
_cell_length_b    23.28001
_cell_length_c    23.28001
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

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_space_group_symop_id
_space_group_symop_operation_xyz
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2 x,-y+1/4,-z+1/4
3 -x+1/4,y,-z+1/4
4 -x+1/4,-y+1/4,z
5 y,z,x
6 y,-z+1/4,-x+1/4
7 -y+1/4,z,-x+1/4
8 -y+1/4,-z+1/4,x
9 z,x,y
10 z,-x+1/4,-y+1/4
11 -z+1/4,x,-y+1/4
12 -z+1/4,-x+1/4,y
13 -y,-x,-z
14 -y,x+1/4,z+1/4
15 y+1/4,-x,z+1/4
16 y+1/4,x+1/4,-z
17 -x,-z,-y
18 -x,z+1/4,y+1/4
19 x+1/4,-z,y+1/4
20 x+1/4,z+1/4,-y
21 -z,-y,-x
22 -z,y+1/4,x+1/4
23 z+1/4,-y,x+1/4
24 z+1/4,y+1/4,-x
25 -x,-y,-z
26 -x,y+1/4,z+1/4
27 x+1/4,-y,z+1/4
28 x+1/4,y+1/4,-z
29 -y,-z,-x
30 -y,z+1/4,x+1/4
31 y+1/4,-z,x+1/4
32 y+1/4,z+1/4,-x
33 -z,-x,-y
34 -z,x+1/4,y+1/4
35 z+1/4,-x,y+1/4
36 z+1/4,x+1/4,-y
```

```
37 y,x,z
38 y,-x+1/4,-z+1/4
39 -y+1/4,x,-z+1/4
40 -y+1/4,-x+1/4,z
41 x,z,y
42 x,-z+1/4,-y+1/4
43 -x+1/4,z,-y+1/4
44 -x+1/4,-z+1/4,y
45 z,y,x
46 z,-y+1/4,-x+1/4
47 -z+1/4,y,-x+1/4
48 -z+1/4,-y+1/4,x
49 x,y+1/2,z+1/2
50 x,-y+3/4,-z+3/4
51 -x+1/4,y+1/2,-z+3/4
52 -x+1/4,-y+3/4,z+1/2
53 y,z+1/2,x+1/2
54 y,-z+3/4,-x+3/4
55 -y+1/4,z+1/2,-x+3/4
56 -y+1/4,-z+3/4,x+1/2
57 z,x+1/2,y+1/2
58 z,-x+3/4,-y+3/4
59 -z+1/4,x+1/2,-y+3/4
60 -z+1/4,-x+3/4,y+1/2
61 -y,-x+1/2,-z+1/2
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66 -x,z+3/4,y+3/4
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69 -z,-y+1/2,-x+1/2
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72 z+1/4,y+3/4,-x+1/2
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74 -x,y+3/4,z+3/4
75 x+1/4,-y+1/2,z+3/4
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77 -y,-z+1/2,-x+1/2
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83 z+1/4,-x+1/2,y+3/4
84 z+1/4,x+3/4,-y+1/2
85 y,x+1/2,z+1/2
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88 -y+1/4,-x+3/4,z+1/2
89 x,z+1/2,y+1/2
90 x,-z+3/4,-y+3/4
91 -x+1/4,z+1/2,-y+3/4
92 -x+1/4,-z+3/4,y+1/2
93 z,y+1/2,x+1/2
94 z,-y+3/4,-x+3/4
95 -z+1/4,y+1/2,-x+3/4
96 -z+1/4,-y+3/4,x+1/2
97 x+1/2,y,z+1/2
98 x+1/2,-y+1/4,-z+3/4
99 -x+3/4,y,-z+3/4
100 -x+3/4,-y+1/4,z+1/2
101 y+1/2,z,x+1/2
102 y+1/2,-z+1/4,-x+3/4
103 -y+3/4,z,-x+3/4
104 -y+3/4,-z+1/4,x+1/2
105 z+1/2,x,y+1/2
106 z+1/2,-x+1/4,-y+3/4
107 -z+3/4,x,-y+3/4
108 -z+3/4,-x+1/4,y+1/2
109 -y+1/2,-x,-z+1/2
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130 -z+1/2,x+1/4,y+3/4
131 z+3/4,-x,y+3/4
132 z+3/4,x+1/4,-y+1/2
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141 z+1/2,y,x+1/2
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144 -z+3/4,-y+1/4,x+1/2
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167 z+3/4,-y+1/2,x+1/4
168 z+3/4,y+3/4,-x
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172 x+3/4,y+3/4,-z
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175 y+3/4,-z+1/2,x+1/4
176 y+3/4,z+3/4,-x
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182 y+1/2,-x+3/4,-z+1/4
183 -y+3/4,x+1/2,-z+1/4
184 -y+3/4,-x+3/4,z
185 x+1/2,z+1/2,y
186 x+1/2,-z+3/4,-y+1/4
187 -x+3/4,z+1/2,-y+1/4
188 -x+3/4,-z+3/4,y
189 z+1/2,y+1/2,x
190 z+1/2,-y+3/4,-x+1/4
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192 -z+3/4,-y+3/4,x

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_atom_site_Wyckoff_label
_atom_site_fract_x
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P1 P 8 b 0.37500 0.37500 0.37500 1.00000
H2O2 H2O 32 e -0.05284 -0.05284 -0.05284 1.00000
O1 O 32 e 0.41753 0.41753 0.41753 1.00000
H2O3 H2O 48 f -0.01933 0.12500 0.12500 1.00000
H2O4 H2O 48 f 0.85395 0.12500 0.12500 1.00000
H2O5 H2O 96 g 0.71864 0.71864 0.85825 1.00000
O2 O 96 g 0.83333 0.83333 0.25301 1.00000
O3 O 96 g -0.06100 -0.06100 0.52706 1.00000
O4 O 96 g 0.53780 0.53780 0.37070 1.00000
W1 W 96 g -0.01783 -0.01783 0.38144 1.00000
```

$H_3PW_{12}O_{40}{\cdot}29H_2O$ ($H4_{21}$): A29B40CD12_cF656_227_ae2fg_e3g_b_g - POSCAR

```
A29B40CD12_cF656_227_ae2fg_e3g_b_g & a,x3,x4,x5,x6,x7,z7,x8,z8,x9,z9,x10
    ↪ ,z10,x11,z11 --params=23.28001,-0.05284,0.41753,-0.01933,
    ↪ 0.85395,0.71864,0.85825,0.83333,0.25301,-0.061,0.52706,0.5378,
    ↪ 0.3707,-0.01783,0.38144 & Fd-3m O_{h}^{7} #227 (abe^2f^2g^5) &
    ↪ cF656 & $H4_{21}$ & (H2O)29O40PW12 & 29-phosphotungstic acid (
    ↪ 29-wpa) & A. J. Bradley and J. W. Illingworth, Proc. Roy. Soc.
    ↪ Lond. A 157, 113-131 (1936)
   1.00000000000000
   0.00000000000000  11.64000500000000  11.64000500000000
  11.64000500000000   0.00000000000000  11.64000500000000
  11.64000500000000  11.64000500000000   0.00000000000000
   H2O       O       P       W
    58      80       2      24
Direct
   0.12500000000000   0.12500000000000   0.12500000000000  H2O    (8a)
   0.87500000000000   0.87500000000000   0.87500000000000  H2O    (8a)
  -0.05284000000000  -0.05284000000000  -0.05284000000000  H2O   (32e)
  -0.05284000000000  -0.05284000000000   0.65852000000000  H2O   (32e)
  -0.05284000000000   0.65852000000000  -0.05284000000000  H2O   (32e)
   0.65852000000000  -0.05284000000000  -0.05284000000000  H2O   (32e)
   0.05284000000000   0.05284000000000   0.34148000000000  H2O   (32e)
   0.05284000000000   0.05284000000000   0.05284000000000  H2O   (32e)
   0.05284000000000   0.34148000000000   0.05284000000000  H2O   (32e)
   0.34148000000000   0.05284000000000   0.05284000000000  H2O   (32e)
   0.26933000000000  -0.01933000000000  -0.01933000000000  H2O   (48f)
  -0.01933000000000   0.26933000000000   0.26933000000000  H2O   (48f)
  -0.01933000000000   0.26933000000000  -0.01933000000000  H2O   (48f)
   0.26933000000000  -0.01933000000000   0.26933000000000  H2O   (48f)
  -0.01933000000000  -0.01933000000000   0.26933000000000  H2O   (48f)
   0.26933000000000   0.26933000000000  -0.01933000000000  H2O   (48f)
```

0.73067000000000	0.01933000000000	0.73067000000000	H2O	(48f)
0.01933000000000	0.73067000000000	0.01933000000000	H2O	(48f)
0.01933000000000	0.73067000000000	0.73067000000000	H2O	(48f)
0.73067000000000	0.01933000000000	0.01933000000000	H2O	(48f)
0.01933000000000	0.01933000000000	0.73067000000000	H2O	(48f)
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−0.60395000000000	0.85395000000000	0.85395000000000	H2O	(48f)
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0.85395000000000	−0.60395000000000	0.85395000000000	H2O	(48f)
−0.60395000000000	0.85395000000000	−0.60395000000000	H2O	(48f)
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−0.85395000000000	1.60395000000000	−0.85395000000000	H2O	(48f)
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0.85825000000000	0.85825000000000	−1.79553000000000	H2O	(96g)
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−0.85825000000000	−0.85825000000000	2.79553000000000	H2O	(96g)
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0.38144000000000	−0.41710000000000	0.38144000000000	W	(96g)
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0.41710000000000	0.84578000000000	−0.38144000000000	W	(96g)
0.84578000000000	0.41710000000000	−0.38144000000000	W	(96g)
0.41710000000000	−0.38144000000000	0.84578000000000	W	(96g)
0.84578000000000	−0.38144000000000	0.41710000000000	W	(96g)
−0.38144000000000	0.41710000000000	−0.38144000000000	W	(96g)
−0.38144000000000	0.84578000000000	−0.38144000000000	W	(96g)
−0.38144000000000	0.41710000000000	0.84578000000000	W	(96g)
−0.38144000000000	0.84578000000000	0.41710000000000	W	(96g)
0.84578000000000	−0.38144000000000	−0.38144000000000	W	(96g)
0.41710000000000	−0.38144000000000	−0.38144000000000	W	(96g)

$G7_3$ [Northupite, $Na_3MgCl(CO_3)_2$] (*obsolete*): A2BCD3E6_cF208_227_e_c_d_f_g - CIF

```
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_chemical_formula_sum 'C2 Cl Mg Na3 O6'

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 'T. Watanab\\'{e}'
_journal_name_full_name
;
 Comptes Rendus de l'Acad{\'e}mie des Sciences
;
_journal_volume 193
_journal_year 1931
_journal_page_first 1421
_journal_page_last 1423
_publ_Section_title
;
 Les structures des cristaux de northupite, de northupite brom\\'{e}e et
      ↪ de tychite
;

# Found in The American Mineralogist Crystal Structure Database, 2003

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      ↪ obsolete}}) Structure'
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_aflow_params_values '14.05,0.715,0.35,0.75,0.65'
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_aflow_Pearson 'cF208'

_symmetry_space_group_name_H-M "F 41/d -3 2/m (origin choice 2)"
_symmetry_Int_Tables_number 227

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_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
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2 x,-y+1/4,-z+1/4
3 -x+1/4,y,-z+1/4
4 -x+1/4,-y+1/4,z
5 y,z,x
6 y,-z+1/4,-x+1/4
7 -y+1/4,z,-x+1/4
8 -y+1/4,-z+1/4,x
9 z,x,y
10 z,-x+1/4,-y+1/4
11 -z+1/4,x,-y+1/4
```

```
12 -z+1/4,-x+1/4,y
13 -y,-x,-z
14 -y,x+1/4,z+1/4
15 y+1/4,-x,z+1/4
16 y+1/4,x+1/4,-z
17 -x,-z,-y
18 -x,z+1/4,y+1/4
19 x+1/4,-z,y+1/4
20 x+1/4,z+1/4,-y
21 -z,-y,-x
22 -z,y+1/4,x+1/4
23 z+1/4,-y,x+1/4
24 z+1/4,y+1/4,-x
25 -x,-y,-z
26 -x,y+1/4,z+1/4
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32 y+1/4,z+1/4,-x
33 -z,-x,-y
34 -z,x+1/4,y+1/4
35 z+1/4,-x,y+1/4
36 z+1/4,x+1/4,-y
37 y,x,z
38 y,-x+1/4,-z+1/4
39 -y+1/4,x,-z+1/4
40 -y+1/4,-x+1/4,z
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171 x+3/4,-y+1/2,z+1/4
172 x+3/4,y+3/4,-z
173 -y+1/2,-z+1/2,-x
174 -y+1/2,z+3/4,x+1/4
175 y+3/4,-z+1/2,x+1/4
176 y+3/4,z+3/4,-x
177 -z+1/2,-x+1/2,-y
178 -z+1/2,x+3/4,y+1/4
179 z+3/4,-x+1/2,y+1/4
180 z+3/4,x+3/4,-y
181 y+1/2,x+1/2,z
182 y+1/2,-x+3/4,-z+1/4
183 -y+3/4,x+1/2,-z+1/4
184 -y+3/4,-x+3/4,z
185 x+1/2,z+1/2,y
186 x+1/2,-z+3/4,-y+1/4
187 -x+3/4,z+1/2,-y+1/4
188 -x+3/4,-z+3/4,y
189 z+1/2,y+1/2,x
190 z+1/2,-y+3/4,-x+1/4
191 -z+3/4,y+1/2,-x+1/4
192 -z+3/4,-y+3/4,x

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cl1 Cl 16 c 0.00000 0.00000 0.00000 1.00000
Mg1 Mg 16 d 0.50000 0.50000 0.50000 1.00000
C1 C 32 e 0.71500 0.71500 0.71500 1.00000
Na1 Na 48 f 0.35000 0.12500 0.12500 1.00000
O1 O 96 g 0.75000 0.75000 0.65000 1.00000
```

$G7_3$ [Northupite, $Na_3MgCl(CO_3)_2$] (*obsolete*): A2BCD3E6_cF208_227_e_c_d_f_g - POSCAR

```
A2BCD3E6_cF208_227_e_c_d_f_g & a,x3,x4,x5,z5 --params=14.05,0.715,0.35,
    ↪ 0.75,0.65 & Fd-3m O_{h}^{7} #227 (cdefg) & cF208 & $G7_{3}$ &
    ↪ C2ClMgNa3O6 & Northupite & H. Shiba and T. Watanab\'{e}, {C. R.
    ↪ Acad. Sci. C 193, 1421-1423 (1931)
  1.00000000000000
  0.00000000000000  7.02500000000000  7.02500000000000
  7.02500000000000  0.00000000000000  7.02500000000000
  7.02500000000000  7.02500000000000  0.00000000000000
    C   Cl   Mg   Na   O
    8    4    4   12   24
Direct
```

0.71500000000000	0.71500000000000	0.71500000000000	C	(32e)
0.71500000000000	0.71500000000000	−1.64500000000000	C	(32e)
0.71500000000000	−1.64500000000000	0.71500000000000	C	(32e)
−1.64500000000000	0.71500000000000	0.71500000000000	C	(32e)
−0.71500000000000	−0.71500000000000	2.64500000000000	C	(32e)
−0.71500000000000	−0.71500000000000	−0.71500000000000	C	(32e)
−0.71500000000000	2.64500000000000	−0.71500000000000	C	(32e)
2.64500000000000	−0.71500000000000	−0.71500000000000	C	(32e)
0.00000000000000	0.00000000000000	0.00000000000000	Cl	(16c)
0.00000000000000	0.00000000000000	0.50000000000000	Cl	(16c)
0.00000000000000	0.50000000000000	0.00000000000000	Cl	(16c)
0.50000000000000	0.00000000000000	0.00000000000000	Cl	(16c)
0.50000000000000	0.50000000000000	0.50000000000000	Mg	(16d)
0.50000000000000	0.50000000000000	0.00000000000000	Mg	(16d)
0.50000000000000	0.00000000000000	0.50000000000000	Mg	(16d)
0.00000000000000	0.50000000000000	0.50000000000000	Mg	(16d)
−0.10000000000000	0.35000000000000	0.35000000000000	Na	(48f)
0.35000000000000	−0.10000000000000	−0.10000000000000	Na	(48f)
0.35000000000000	−0.10000000000000	0.35000000000000	Na	(48f)
−0.10000000000000	0.35000000000000	−0.10000000000000	Na	(48f)
0.35000000000000	0.35000000000000	−0.10000000000000	Na	(48f)
−0.10000000000000	−0.10000000000000	0.35000000000000	Na	(48f)
1.10000000000000	−0.35000000000000	1.10000000000000	Na	(48f)
−0.35000000000000	1.10000000000000	−0.35000000000000	Na	(48f)
−0.35000000000000	1.10000000000000	1.10000000000000	Na	(48f)
1.10000000000000	−0.35000000000000	−0.35000000000000	Na	(48f)
−0.35000000000000	−0.35000000000000	1.10000000000000	Na	(48f)
1.10000000000000	1.10000000000000	−0.35000000000000	Na	(48f)
0.65000000000000	0.65000000000000	0.85000000000000	O	(96g)
0.65000000000000	0.65000000000000	−1.65000000000000	O	(96g)
0.85000000000000	−1.65000000000000	0.65000000000000	O	(96g)
−1.65000000000000	0.85000000000000	0.65000000000000	O	(96g)
0.85000000000000	0.65000000000000	0.65000000000000	O	(96g)
−1.65000000000000	0.65000000000000	0.65000000000000	O	(96g)
0.65000000000000	0.85000000000000	−1.65000000000000	O	(96g)
0.65000000000000	−1.65000000000000	0.85000000000000	O	(96g)
0.65000000000000	0.85000000000000	0.65000000000000	O	(96g)
0.65000000000000	−1.65000000000000	0.65000000000000	O	(96g)
−1.65000000000000	0.65000000000000	0.85000000000000	O	(96g)
0.85000000000000	0.65000000000000	−1.65000000000000	O	(96g)
−0.65000000000000	−0.65000000000000	2.65000000000000	O	(96g)
−0.65000000000000	−0.65000000000000	−0.85000000000000	O	(96g)
−0.85000000000000	2.65000000000000	−0.65000000000000	O	(96g)
2.65000000000000	−0.85000000000000	−0.65000000000000	O	(96g)
−0.85000000000000	−0.65000000000000	2.65000000000000	O	(96g)
2.65000000000000	−0.65000000000000	−0.85000000000000	O	(96g)
−0.65000000000000	−0.85000000000000	−0.65000000000000	O	(96g)
−0.65000000000000	2.65000000000000	−0.65000000000000	O	(96g)
−0.65000000000000	−0.85000000000000	2.65000000000000	O	(96g)
−0.65000000000000	2.65000000000000	−0.85000000000000	O	(96g)
2.65000000000000	−0.65000000000000	−0.65000000000000	O	(96g)
−0.85000000000000	−0.65000000000000	−0.65000000000000	O	(96g)

$D6_2$ (Sb_2O_4) (*obsolete*): A2B_cF96_227_abf_cd - CIF

```
# CIF file
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_audit_creation_method FINDSYM

_chemical_name_mineral 'O2Sb'
_chemical_formula_sum 'O2 Sb'

loop_
_publ_author_name
 'G. Natta'
 'M. Baccaredda'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Kristallographie - Crystalline Materials
;
_journal_volume 85
_journal_year 1933
_journal_page_first 271
_journal_page_last 296
_publ_Section_title
;
 Tetrossido di antimonio e antimoniati
;

# Found in Strukturbericht Band III 1933-1935, 1937

_aflow_title '$D6_{2}$ (Sb$_{2}$O$_{4}$) ({\em{obsolete}}) Structure'
_aflow_proto 'A2B_cF96_227_abf_cd'
_aflow_params 'a,x_{5}'
_aflow_params_values '10.24,0.395'
_aflow_Strukturbericht '$D6_{2}$'
_aflow_Pearson 'cF96'

_symmetry_space_group_name_H-M "F 41/d -3 2/m (origin choice 2)"
_symmetry_Int_Tables_number 227

_cell_length_a     10.24000
_cell_length_b     10.24000
_cell_length_c     10.24000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y+1/4,-z+1/4
3 -x+1/4,y,-z+1/4
4 -x+1/4,-y+1/4,z
5 y,z,x
6 y,-z+1/4,-x+1/4
7 -y+1/4,z,-x+1/4
8 -y+1/4,-z+1/4,x
9 z,x,y
10 z,-x+1/4,-y+1/4
11 -z+1/4,x,-y+1/4
12 -z+1/4,-x+1/4,y
13 -y,-x,-z
14 -y,x+1/4,z+1/4
15 y+1/4,-x,z+1/4
16 y+1/4,x+1/4,-z
17 -x,-z,-y
18 -x,z+1/4,y+1/4
19 x+1/4,-z,y+1/4
20 x+1/4,z+1/4,-y
21 -z,-y,-x
22 -z,y+1/4,x+1/4
23 z+1/4,-y,x+1/4
24 z+1/4,y+1/4,-x
25 -x,-y,-z
26 -x,y+1/4,z+1/4
27 x+1/4,-y,z+1/4
28 x+1/4,y+1/4,-z
29 -y,-z,-x
30 -y,z+1/4,x+1/4
31 y+1/4,-z,x+1/4
32 y+1/4,z+1/4,-x
33 -z,-x,-y
34 -z,x+1/4,y+1/4
35 z+1/4,-x,y+1/4
36 z+1/4,x+1/4,-y
37 y,x,z
38 y,-x+1/4,-z+1/4
39 -y+1/4,x,-z+1/4
40 -y+1/4,-x+1/4,z
41 x,z,y
42 x,-z+1/4,-y+1/4
43 -x+1/4,z,-y+1/4
44 -x+1/4,-z+1/4,y
45 z,y,x
46 z,-y+1/4,-x+1/4
47 -z+1/4,y,-x+1/4
48 -z+1/4,-y+1/4,x
49 x,y+1/2,z+1/2
50 x,-y+3/4,-z+3/4
51 -x+1/4,y+1/2,-z+3/4
52 -x+1/4,-y+3/4,z+1/2
53 y,z+1/2,x+1/2
54 y,-z+3/4,-x+3/4
55 -y+1/4,z+1/2,-x+3/4
56 -y+1/4,-z+3/4,x+1/2
57 z,x+1/2,y+1/2
58 z,-x+3/4,-y+3/4
59 -z+1/4,x+1/2,-y+3/4
60 -z+1/4,-x+3/4,y+1/2
61 -y,-x+1/2,-z+1/2
62 -y,x+3/4,z+3/4
63 y+1/4,-x+1/2,z+3/4
64 y+1/4,x+3/4,-z+1/2
65 -x,-z+1/2,-y+1/2
66 -x,z+3/4,y+3/4
67 x+1/4,-z+1/2,y+3/4
68 x+1/4,z+3/4,-y+1/2
69 -z,-y+1/2,-x+1/2
70 -z,y+3/4,x+3/4
71 z+1/4,-y+1/2,x+3/4
72 z+1/4,y+3/4,-x+1/2
73 -x,-y+1/2,-z+1/2
74 -x,y+3/4,z+3/4
75 x+1/4,-y+1/2,z+3/4
76 x+1/4,y+3/4,-z+1/2
77 -y,-z+1/2,-x+1/2
78 -y,z+3/4,x+3/4
79 y+1/4,-z+1/2,x+3/4
80 y+1/4,z+3/4,-x+1/2
81 -z,-x+1/2,-y+1/2
82 -z,x+3/4,y+3/4
83 z+1/4,-x+1/2,y+3/4
84 z+1/4,x+3/4,-y+1/2
85 y,x+1/2,z+1/2
86 y,-x+3/4,-z+3/4
87 -y+1/4,x+1/2,-z+3/4
88 -y+1/4,-x+3/4,z+1/2
89 x,z+1/2,y+1/2
90 x,-z+3/4,-y+3/4
91 -x+1/4,z+1/2,-y+3/4
92 -x+1/4,-z+3/4,y+1/2
93 z,y+1/2,x+1/2
94 z,-y+3/4,-x+3/4
95 -z+1/4,y+1/2,-x+3/4
96 -z+1/4,-y+3/4,x+1/2
97 x+1/2,y,z+1/2
98 x+1/2,-y+1/4,-z+3/4
99 -x+3/4,y,-z+3/4
100 -x+3/4,-y+1/4,z+1/2
101 y+1/2,z,x+1/2
102 y+1/2,-z+1/4,-x+3/4
103 -y+3/4,z,-x+3/4
104 -y+3/4,-z+1/4,x+1/2
105 z+1/2,x,y+1/2
106 z+1/2,-x+1/4,-y+3/4
107 -z+3/4,x,-y+3/4
108 -z+3/4,-x+1/4,y+1/2
109 -y+1/2,-x,-z+1/2
```

```
110 -y+1/2,x+1/4,z+3/4
111 y+3/4,-x,z+3/4
112 y+3/4,x+1/4,-z+1/2
113 -x+1/2,-z,-y+1/2
114 -x+1/2,z+1/4,y+3/4
115 x+3/4,-z,y+3/4
116 x+3/4,z+1/4,-y+1/2
117 -z+1/2,-y,-x+1/2
118 -z+1/2,y+1/4,x+3/4
119 z+3/4,-y,x+3/4
120 z+3/4,y+1/4,-x+1/2
121 -x+1/2,-y,-z+1/2
122 -x+1/2,y+1/4,z+3/4
123 x+3/4,-y,z+3/4
124 x+3/4,y+1/4,-z+1/2
125 -y+1/2,-z,-x+1/2
126 -y+1/2,z+1/4,x+3/4
127 y+3/4,-z,x+3/4
128 y+3/4,z+1/4,-x+1/2
129 -z+1/2,-x,-y+1/2
130 -z+1/2,x+1/4,y+3/4
131 z+3/4,-x,y+3/4
132 z+3/4,x+1/4,-y+1/2
133 y+1/2,x,z+1/2
134 y+1/2,-x+1/4,-z+3/4
135 -y+3/4,x,-z+3/4
136 -y+3/4,-x+1/4,z+1/2
137 x+1/2,z,y+1/2
138 x+1/2,-z+1/4,-y+3/4
139 -x+3/4,z,-y+3/4
140 -x+3/4,-z+1/4,y+1/2
141 z+1/2,y,x+1/2
142 z+1/2,-y+1/4,-x+3/4
143 -z+3/4,y,-x+3/4
144 -z+3/4,-y+1/4,x+1/2
145 x+1/2,y+1/2,z
146 x+1/2,-y+3/4,-z+1/4
147 -x+3/4,y+1/2,-z+1/4
148 -x+3/4,-y+3/4,z
149 y+1/2,z+1/2,x
150 y+1/2,-z+3/4,-x+1/4
151 -y+3/4,z+1/2,-x+1/4
152 -y+3/4,-z+3/4,x
153 z+1/2,x+1/2,y
154 z+1/2,-x+3/4,-y+1/4
155 -z+3/4,x+1/2,-y+1/4
156 -z+3/4,-x+3/4,y
157 -y+1/2,-x+1/2,-z
158 -y+1/2,x+3/4,z+1/4
159 y+3/4,-x+1/2,z+1/4
160 y+3/4,x+3/4,-z
161 -x+1/2,-z+1/2,-y
162 -x+1/2,z+3/4,y+1/4
163 x+3/4,-z+1/2,y+1/4
164 x+3/4,z+3/4,-y
165 -z+1/2,-y+1/2,-x
166 -z+1/2,y+3/4,x+1/4
167 z+3/4,-y+1/2,x+1/4
168 z+3/4,y+3/4,-x
169 -x+1/2,-y+1/2,-z
170 -x+1/2,y+3/4,z+1/4
171 x+3/4,-y+1/2,z+1/4
172 x+3/4,y+3/4,-z
173 -y+1/2,-z+1/2,-x
174 -y+1/2,z+3/4,x+1/4
175 y+3/4,-z+1/2,x+1/4
176 y+3/4,z+3/4,-x
177 -z+1/2,-x+1/2,-y
178 -z+1/2,x+3/4,y+1/4
179 z+3/4,-x+1/2,y+1/4
180 z+3/4,x+3/4,-y
181 y+1/2,x+1/2,z
182 y+1/2,-x+3/4,-z+1/4
183 -y+3/4,x+1/2,-z+1/4
184 -y+3/4,-x+3/4,z
185 x+1/2,z+1/2,y
186 x+1/2,-z+3/4,-y+1/4
187 -x+3/4,z+1/2,-y+1/4
188 -x+3/4,-z+3/4,y
189 z+1/2,y+1/2,x
190 z+1/2,-y+3/4,-x+1/4
191 -z+3/4,y+1/2,-x+1/4
192 -z+3/4,-y+3/4,x

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
O1 O 8 a 0.12500 0.12500 0.12500 1.00000
O2 O 8 b 0.37500 0.37500 0.37500 1.00000
Sb1 Sb 16 c 0.00000 0.00000 0.00000 1.00000
Sb2 Sb 16 d 0.50000 0.50000 0.50000 1.00000
O3 O 48 f 0.39500 0.12500 0.12500 1.00000
```

$D6_2$ (Sb_2O_4) (*obsolete*): A2B_cF96_227_abf_cd - POSCAR

```
A2B_cF96_227_abf_cd & a,x5 --params=10.24,0.395 & Fd-3m O_{h}^{7} #227 (
    ↪ abcdf) & cF96 & $D6_{2}$ & O2Sb & O2Sb & G. Natta and M.
    ↪ Baccaredda, Zeitschrift f{"u}r Kristallographie - Crystalline
    ↪ Materials 85, 271-296 (1933)
   1.00000000000000
   0.00000000000000   5.12000000000000   5.12000000000000
   5.12000000000000   0.00000000000000   5.12000000000000
   5.12000000000000   5.12000000000000   0.00000000000000
     O     Sb
    16      8
Direct
   0.12500000000000   0.12500000000000   0.12500000000000   O   (8a)
   0.87500000000000   0.87500000000000   0.87500000000000   O   (8a)
   0.37500000000000   0.37500000000000   0.37500000000000   O   (8b)
   0.62500000000000   0.62500000000000   0.62500000000000   O   (8b)
  -0.14500000000000   0.39500000000000   0.39500000000000   O   (48f)
   0.39500000000000  -0.14500000000000  -0.14500000000000   O   (48f)
   0.39500000000000  -0.14500000000000   0.39500000000000   O   (48f)
  -0.14500000000000   0.39500000000000  -0.14500000000000   O   (48f)
   0.39500000000000   0.39500000000000  -0.14500000000000   O   (48f)
  -0.14500000000000  -0.14500000000000   0.39500000000000   O   (48f)
   1.14500000000000  -0.39500000000000   1.14500000000000   O   (48f)
  -0.39500000000000   1.14500000000000  -0.39500000000000   O   (48f)
  -0.39500000000000   1.14500000000000   1.14500000000000   O   (48f)
   1.14500000000000  -0.39500000000000   1.14500000000000   O   (48f)
  -0.39500000000000  -0.39500000000000   1.14500000000000   O   (48f)
   1.14500000000000   1.14500000000000  -0.39500000000000   O   (48f)
   0.00000000000000   0.00000000000000   0.00000000000000   Sb  (16c)
   0.00000000000000   0.00000000000000   0.50000000000000   Sb  (16c)
   0.00000000000000   0.50000000000000   0.00000000000000   Sb  (16c)
   0.50000000000000   0.00000000000000   0.00000000000000   Sb  (16c)
   0.50000000000000   0.50000000000000   0.50000000000000   Sb  (16d)
   0.50000000000000   0.50000000000000   0.00000000000000   Sb  (16d)
   0.50000000000000   0.00000000000000   0.50000000000000   Sb  (16d)
   0.00000000000000   0.50000000000000   0.50000000000000   Sb  (16d)
```

Senarmontite (Sb_2O_3, $D6_1$): A3B2_cF80_227_f_e - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'Senarmontite'
_chemical_formula_sum 'O3 Sb2'

loop_
_publ_author_name
 'C. Svensson'
_journal_name_full_name
;
 Acta Crystallographica Section B: Structural Science
;
_journal_volume 31
_journal_year 1975
_journal_page_first 2016
_journal_page_last 2018
_publ_Section_title
;
 Refinement of the crystal structure of cubic antimony trioxide, Sb$_{2}
    ↪ $O$_{3}$
;

_aflow_title 'Senarmontite (Sb$_{2}$O$_{3}$, $D6_{1}$) Structure'
_aflow_proto 'A3B2_cF80_227_f_e'
_aflow_params 'a,x_{1},x_{2}'
_aflow_params_values '11.1519,0.26027,0.43875'
_aflow_Strukturbericht '$D6_{1}$'
_aflow_Pearson 'cF80'

_symmetry_space_group_name_H-M "F 41/d -3 2/m (origin choice 2)"
_symmetry_Int_Tables_number 227

_cell_length_a    11.15190
_cell_length_b    11.15190
_cell_length_c    11.15190
_cell_angle_alpha 90.00000
_cell_angle_beta  90.00000
_cell_angle_gamma 90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y+1/4,-z+1/4
3 -x+1/4,y,-z+1/4
4 -x+1/4,-y+1/4,z
5 y,z,x
6 y,-z+1/4,-x+1/4
7 -y+1/4,z,-x+1/4
8 -y+1/4,-z+1/4,x
9 z,x,y
10 z,-x+1/4,-y+1/4
11 -z+1/4,x,-y+1/4
12 -z+1/4,-x+1/4,y
13 -y,-x,-z
14 -y,x+1/4,z+1/4
15 y+1/4,-x,z+1/4
16 y+1/4,x+1/4,-z
17 -x,-z,-y
18 -x,z+1/4,y+1/4
19 x+1/4,-z,y+1/4
20 x+1/4,z+1/4,-y
21 -z,-y,-x
22 -z,y+1/4,x+1/4
23 z+1/4,-y,x+1/4
24 z+1/4,y+1/4,-x
25 -x,-y,-z
26 -x,y+1/4,z+1/4
27 x+1/4,-y,z+1/4
```

```
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29 -y,-z,-x
30 -y,z+1/4,x+1/4
31 y+1/4,-z,x+1/4
32 y+1/4,z+1/4,-x
33 -z,-x,-y
34 -z,x+1/4,y+1/4
35 z+1/4,-x,y+1/4
36 z+1/4,x+1/4,-y
37 y,x,z
38 y,-x+1/4,-z+1/4
39 -y+1/4,x,-z+1/4
40 -y+1/4,-x+1/4,z
41 x,z,y
42 x,-z+1/4,-y+1/4
43 -x+1/4,z,-y+1/4
44 -x+1/4,-z+1/4,y
45 z,y,x
46 z,-y+1/4,-x+1/4
47 -z+1/4,y,-x+1/4
48 -z+1/4,-y+1/4,x
49 x,y+1/2,z+1/2
50 x,-y+3/4,-z+3/4
51 -x+1/4,y+1/2,-z+3/4
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53 y,z+1/2,x+1/2
54 y,-z+3/4,-x+3/4
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68 x+1/4,z+3/4,-y+1/2
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72 z+1/4,y+3/4,-x+1/2
73 -x,-y+1/2,-z+1/2
74 -x,y+3/4,z+3/4
75 x+1/4,-y+1/2,z+3/4
76 x+1/4,y+3/4,-z+1/2
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184 -y+3/4,-x+3/4,z
185 x+1/2,z+1/2,y
186 x+1/2,-z+3/4,-y+1/4
187 -x+3/4,z+1/2,-y+1/4
188 -x+3/4,-z+3/4,y
189 z+1/2,y+1/2,x
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O1 O 48 f 0.43875 0.12500 0.12500 1.00000
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Senarmontite (Sb_2O_3, $D6_1$): A3B2_cF80_227_f_e - POSCAR

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    ↪ h}^{7} #227 (ef) & cF80 & $D6_{1}$ & O3Sb2 & Senarmontite & C.
    ↪ Svensson, Acta Crystallogr. Sect. B Struct. Sci. 31, 2016-2018
    ↪ (1975)
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   5.57595000000000   0.00000000000000   5.57595000000000
   5.57595000000000   5.57595000000000   0.00000000000000
     O     Sb
    12      8
Direct
  -0.18875000000000   0.43875000000000   0.43875000000000   O   (48f)
   0.43875000000000  -0.18875000000000  -0.18875000000000   O   (48f)
   0.43875000000000  -0.18875000000000   0.43875000000000   O   (48f)
  -0.18875000000000   0.43875000000000  -0.18875000000000   O   (48f)
   0.43875000000000   0.43875000000000  -0.18875000000000   O   (48f)
  -0.18875000000000  -0.18875000000000   0.43875000000000   O   (48f)
   1.18875000000000  -0.43875000000000   1.18875000000000   O   (48f)
  -0.43875000000000   1.18875000000000  -0.43875000000000   O   (48f)
  -0.43875000000000   1.18875000000000   1.18875000000000   O   (48f)
   1.18875000000000  -0.43875000000000  -0.43875000000000   O   (48f)
  -0.43875000000000  -0.43875000000000   1.18875000000000   O   (48f)
   1.18875000000000   1.18875000000000  -0.43875000000000   O   (48f)
   0.26027000000000   0.26027000000000   0.26027000000000   Sb  (32e)
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   0.26027000000000  -0.28081000000000   0.26027000000000   Sb  (32e)
  -0.28081000000000   0.26027000000000   0.26027000000000   Sb  (32e)
  -0.26027000000000  -0.26027000000000   1.28081000000000   Sb  (32e)
  -0.26027000000000  -0.26027000000000  -0.26027000000000   Sb  (32e)
  -0.26027000000000   1.28081000000000  -0.26027000000000   Sb  (32e)
```

```
   1.28081000000000   −0.26027000000000   −0.26027000000000    Sb   (32e)
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$H5_6$ [Tychite, $Na_6Mg_2SO_4(CO_3)_4$)] (*obsolete*): A4B2C6D16E_cF232_227_e_d_f_eg_a - CIF

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 'T. Watanab\\'{e}'
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 Comptes Rendus de l'Acad{\'e}mie des Sciences
;
_journal_volume 193
_journal_year 1931
_journal_page_first 1421
_journal_page_last 1423
_publ_Section_title
;
 Les structures des cristaux de northupite, de northupite brom\\'{e}e et
      ↪ de tychite
;

# Found in Tychite, Na$_{6}$Mg$_{2}$(SO$_{4}$)(CO$_{3}$)$_{4}$:
      ↪ Structure analysis and Raman spectroscopic data, 2006

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_space_group_symop_operation_xyz
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4 -x+1/4,-y+1/4,z
5 y,z,x
6 y,-z+1/4,-x+1/4
7 -y+1/4,z,-x+1/4
8 -y+1/4,-z+1/4,x
9 z,x,y
10 z,-x+1/4,-y+1/4
11 -z+1/4,x,-y+1/4
12 -z+1/4,-x+1/4,y
13 -y,-x,-z
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25 -x,-y,-z
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37 y,x,z
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39 -y+1/4,x,-z+1/4
40 -y+1/4,-x+1/4,z
41 x,z,y
42 x,-z+1/4,-y+1/4
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44 -x+1/4,-z+1/4,y
45 z,y,x
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49 x,y+1/2,z+1/2
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192 -z+3/4,-y+3/4,x

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Mg1 Mg 16 d 0.50000 0.50000 0.50000 1.00000
C1 C 32 e 0.73500 0.73500 0.73500 1.00000
O1 O 32 e 0.06000 0.06000 0.06000 1.00000
Na1 Na 48 f 0.35000 0.12500 0.12500 1.00000
O2 O 96 g 0.76500 0.76500 0.66500 1.00000
```

$H5_6$ [Tychite, $Na_6Mg_2SO_4(CO_3)_4$)] (*obsolete*): A4B2C6D16E_cF232_227_e_d_f_eg_a - POSCAR

```
A4B2C6D16E_cF232_227_e_d_f_eg_a & a,x3,x4,x5,x6,z --params=13.89999,
      ↪ 0.735,0.06,0.35,0.765,0.665 & Fd-3m O_{h}^{7} #227 (ade^2fg) &
      ↪ cF232 & $H5_{6}$ & C4Mg2Na6O16S & Tychite & H. Shiba and T.
      ↪ Watanab\''{e}, {C. R. Acad. Sci. C 193, 1421-1423 (1931)
   1.00000000000000
   0.00000000000000   6.94999500000000   6.94999500000000
   6.94999500000000   0.00000000000000   6.94999500000000
   6.94999500000000   6.94999500000000   0.00000000000000
     C     Mg     Na      O      S
     8      4     12     32      2
Direct
   0.73500000000000   0.73500000000000   0.73500000000000   C  (32e)
   0.73500000000000   0.73500000000000  -1.70500000000000   C  (32e)
   0.73500000000000  -1.70500000000000   0.73500000000000   C  (32e)
  -1.70500000000000   0.73500000000000   0.73500000000000   C  (32e)
  -0.73500000000000  -0.73500000000000   2.70500000000000   C  (32e)
  -0.73500000000000  -0.73500000000000  -0.73500000000000   C  (32e)
  -0.73500000000000   2.70500000000000  -0.73500000000000   C  (32e)
   2.70500000000000  -0.73500000000000  -0.73500000000000   C  (32e)
   0.50000000000000   0.50000000000000   0.50000000000000  Mg  (16d)
   0.50000000000000   0.50000000000000   0.00000000000000  Mg  (16d)
   0.50000000000000   0.00000000000000   0.50000000000000  Mg  (16d)
   0.00000000000000   0.50000000000000   0.50000000000000  Mg  (16d)
  -0.10000000000000   0.35000000000000   0.35000000000000  Na  (48f)
   0.35000000000000  -0.10000000000000  -0.10000000000000  Na  (48f)
   0.35000000000000  -0.10000000000000   0.35000000000000  Na  (48f)
  -0.10000000000000   0.35000000000000  -0.10000000000000  Na  (48f)
   0.35000000000000   0.35000000000000  -0.10000000000000  Na  (48f)
  -0.10000000000000  -0.10000000000000   0.35000000000000  Na  (48f)
   1.10000000000000  -0.35000000000000   1.10000000000000  Na  (48f)
  -0.35000000000000   1.10000000000000  -0.35000000000000  Na  (48f)
  -0.35000000000000   1.10000000000000   1.10000000000000  Na  (48f)
   1.10000000000000  -0.35000000000000  -0.35000000000000  Na  (48f)
  -0.35000000000000  -0.35000000000000   1.10000000000000  Na  (48f)
   1.10000000000000   1.10000000000000  -0.35000000000000  Na  (48f)
   0.06000000000000   0.06000000000000   0.06000000000000   O  (32e)
   0.06000000000000   0.06000000000000   0.32000000000000   O  (32e)
   0.06000000000000   0.32000000000000   0.06000000000000   O  (32e)
   0.32000000000000   0.06000000000000   0.06000000000000   O  (32e)
  -0.06000000000000  -0.06000000000000   0.68000000000000   O  (32e)
  -0.06000000000000  -0.06000000000000  -0.06000000000000   O  (32e)
  -0.06000000000000   0.68000000000000  -0.06000000000000   O  (32e)
   0.68000000000000  -0.06000000000000  -0.06000000000000   O  (32e)
   0.66500000000000   0.66500000000000   0.86500000000000   O  (96g)
   0.66500000000000   0.66500000000000  -1.69500000000000   O  (96g)
   0.86500000000000  -1.69500000000000   0.66500000000000   O  (96g)
  -1.69500000000000   0.86500000000000   0.66500000000000   O  (96g)
   0.86500000000000   0.66500000000000   0.66500000000000   O  (96g)
  -1.69500000000000   0.66500000000000   0.66500000000000   O  (96g)
   0.66500000000000   0.86500000000000  -1.69500000000000   O  (96g)
   0.66500000000000  -1.69500000000000   0.86500000000000   O  (96g)
   0.66500000000000   0.86500000000000   0.66500000000000   O  (96g)
   0.66500000000000  -1.69500000000000   0.66500000000000   O  (96g)
  -1.69500000000000   0.66500000000000   0.86500000000000   O  (96g)
   0.86500000000000   0.66500000000000  -1.69500000000000   O  (96g)
   0.66500000000000   0.66500000000000   0.66500000000000   O  (96g)
   0.66500000000000  -1.69500000000000   0.66500000000000   O  (96g)
  -1.69500000000000   0.66500000000000   0.66500000000000   O  (96g)
   0.86500000000000   0.66500000000000   0.66500000000000   O  (96g)
   0.66500000000000   0.66500000000000   0.66500000000000   O  (96g)
  -0.86500000000000  -0.66500000000000   2.69500000000000   O  (96g)
  -0.66500000000000  -0.66500000000000  -0.86500000000000   O  (96g)
  -0.86500000000000   2.69500000000000  -0.66500000000000   O  (96g)
   2.69500000000000  -0.86500000000000  -0.66500000000000   O  (96g)
  -0.86500000000000  -0.66500000000000   2.69500000000000   O  (96g)
   2.69500000000000  -0.66500000000000  -0.86500000000000   O  (96g)
  -0.66500000000000  -0.86500000000000  -0.66500000000000   O  (96g)
  -0.66500000000000   2.69500000000000  -0.66500000000000   O  (96g)
  -0.66500000000000  -0.86500000000000   2.69500000000000   O  (96g)
  -0.66500000000000   2.69500000000000  -0.86500000000000   O  (96g)
   2.69500000000000  -0.66500000000000  -0.66500000000000   O  (96g)
  -0.86500000000000  -0.66500000000000  -0.66500000000000   O  (96g)
   0.12500000000000   0.12500000000000   0.12500000000000   S   (8a)
   0.87500000000000   0.87500000000000   0.87500000000000   S   (8a)
```

Cubic CuPt ($L1_3$ (I), $D4$): AB_cF32_227_c_d - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'CuPt'
_chemical_formula_sum 'Cu Pt'

loop_
_publ_author_name
 'C. H. Johansson'
 'J. O. Linde'
_journal_name_full_name
;
 Annalen der Physik
;
_journal_volume 387
_journal_year 1927
_journal_page_first 449
_journal_page_last 478
_publ_Section_title
;
 Gitterstruktur und elektrisches Leitverm\"{o}gen der
      ↪ Mischkristallreihen Au-Cu, Pd-Cu und Pt-Cu
;

# Found in Strukturbericht 1913-1928, 1931

_aflow_title 'Cubic CuPt ($L1_{3}$ (I), $D4$) Structure'
_aflow_proto 'AB_cF32_227_c_d'
_aflow_params 'a'
_aflow_params_values '7.23'
_aflow_Strukturbericht '$L1_{3}$ (I)'
_aflow_Pearson 'cF32'

_symmetry_space_group_name_H-M "F 41/d -3 2/m (origin choice 2)"
_symmetry_Int_Tables_number 227

_cell_length_a     7.230000
_cell_length_b     7.230000
_cell_length_c     7.230000
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y+1/4,-z+1/4
3 -x+1/4,y,-z+1/4
4 -x+1/4,-y+1/4,z
5 y,z,x
6 y,-z+1/4,-x+1/4
7 -y+1/4,z,-x+1/4
8 -y+1/4,-z+1/4,x
9 z,x,y
10 z,-x+1/4,-y+1/4
11 -z+1/4,x,-y+1/4
12 -z+1/4,-x+1/4,y
13 -y,-x,-z
14 -y,x+1/4,z+1/4
15 y+1/4,-x,z+1/4
16 y+1/4,x+1/4,-z
17 -x,-z,-y
18 -x,z+1/4,y+1/4
19 x+1/4,-z,y+1/4
20 x+1/4,z+1/4,-y
21 -z,-y,-x
22 -z,y+1/4,x+1/4
23 z+1/4,-y,x+1/4
24 z+1/4,y+1/4,-x
25 -x,-y,-z
26 -x,y+1/4,z+1/4
27 x+1/4,-y,z+1/4
28 x+1/4,y+1/4,-z
29 -y,-z,-x
30 -y,z+1/4,x+1/4
31 y+1/4,-z,x+1/4
32 y+1/4,z+1/4,-x
33 -z,-x,-y
34 -z,x+1/4,y+1/4
35 z+1/4,-x,y+1/4
36 z+1/4,x+1/4,-y
37 y,x,z
```

```
38 y,-x+1/4,-z+1/4
39 -y+1/4,x,-z+1/4
40 -y+1/4,-x+1/4,z
41 x,z,y
42 x,-z+1/4,-y+1/4
43 -x+1/4,z,-y+1/4
44 -x+1/4,-z+1/4,y
45 z,y,x
46 z,-y+1/4,-x+1/4
47 -z+1/4,y,-x+1/4
48 -z+1/4,-y+1/4,x
49 x,y+1/2,z+1/2
50 x,-y+3/4,-z+3/4
51 -x+1/4,y+1/2,-z+3/4
52 -x+1/4,-y+3/4,z+1/2
53 y,z+1/2,x+1/2
54 y,-z+3/4,-x+3/4
55 -y+1/4,z+1/2,-x+3/4
56 -y+1/4,-z+3/4,x+1/2
57 z,x+1/2,y+1/2
58 z,-x+3/4,-y+3/4
59 -z+1/4,x+1/2,-y+3/4
60 -z+1/4,-x+3/4,y+1/2
61 -y,-x+1/2,-z+1/2
62 -y,x+3/4,z+3/4
63 y+1/4,-x+1/2,z+3/4
64 y+1/4,x+3/4,-z+1/2
65 -x,-z+1/2,-y+1/2
66 -x,z+3/4,y+3/4
67 x+1/4,-z+1/2,y+3/4
68 x+1/4,z+3/4,-y+1/2
69 -z,-y+1/2,-x+1/2
70 -z,y+3/4,x+3/4
71 z+1/4,-y+1/2,x+3/4
72 z+1/4,y+3/4,-x+1/2
73 -x,-y+1/2,-z+1/2
74 -x,y+3/4,z+3/4
75 x+1/4,-y+1/2,z+3/4
76 x+1/4,y+3/4,-z+1/2
77 -y,-z+1/2,-x+1/2
78 -y,z+3/4,x+3/4
79 y+1/4,-z+1/2,x+3/4
80 y+1/4,z+3/4,-x+1/2
81 -z,-x+1/2,-y+1/2
82 -z,x+3/4,y+3/4
83 z+1/4,-x+1/2,y+3/4
84 z+1/4,x+3/4,-y+1/2
85 y,x+1/2,z+1/2
86 y,-x+3/4,-z+3/4
87 -y+1/4,x+1/2,-z+3/4
88 -y+1/4,-x+3/4,z+1/2
89 x,z+1/2,y+1/2
90 x,-z+3/4,-y+3/4
91 -x+1/4,z+1/2,-y+3/4
92 -x+1/4,-z+3/4,y+1/2
93 z,y+1/2,x+1/2
94 z,-y+3/4,-x+3/4
95 -z+1/4,y+1/2,-x+3/4
96 -z+1/4,-y+3/4,x+1/2
97 x+1/2,y,z+1/2
98 x+1/2,-y+1/4,-z+3/4
99 -x+3/4,y,-z+3/4
100 -x+3/4,-y+1/4,z+1/2
101 y+1/2,z,x+1/2
102 y+1/2,-z+1/4,-x+3/4
103 -y+3/4,z,-x+3/4
104 -y+3/4,-z+1/4,x+1/2
105 z+1/2,x,y+1/2
106 z+1/2,-x+1/4,-y+3/4
107 -z+3/4,x,-y+3/4
108 -z+3/4,-x+1/4,y+1/2
109 -y+1/2,-x,-z+1/2
110 -y+1/2,x+1/4,z+3/4
111 y+3/4,-x,z+3/4
112 y+3/4,x+1/4,-z+1/2
113 -x+1/2,-z,-y+1/2
114 -x+1/2,z+1/4,y+3/4
115 x+3/4,-z,y+3/4
116 x+3/4,z+1/4,-y+1/2
117 -z+1/2,-y,-x+1/2
118 -z+1/2,y+1/4,x+3/4
119 z+3/4,-y,x+3/4
120 z+3/4,y+1/4,-x+1/2
121 -x+1/2,-y,-z+1/2
122 -x+1/2,y+1/4,z+3/4
123 x+3/4,-y,z+3/4
124 x+3/4,y+1/4,-z+1/2
125 -y+1/2,-z,-x+1/2
126 -y+1/2,z+1/4,x+3/4
127 y+3/4,-z,x+3/4
128 y+3/4,z+1/4,-x+1/2
129 -z+1/2,-x,-y+1/2
130 -z+1/2,x+1/4,y+3/4
131 z+3/4,-x,y+3/4
132 z+3/4,x+1/4,-y+1/2
133 y+1/2,x,z+1/2
134 y+1/2,-x+1/4,-z+3/4
135 -y+3/4,x,-z+3/4
136 -y+3/4,-x+1/4,z+1/2
137 x+1/2,z,y+1/2
138 x+1/2,-z+1/4,-y+3/4
139 -x+3/4,z,-y+3/4
140 -x+3/4,-z+1/4,y+1/2
141 z+1/2,y,x+1/2
142 z+1/2,-y+1/4,-x+3/4
143 -z+3/4,y,-x+3/4
144 -z+3/4,-y+1/4,x+1/2
145 x+1/2,y+1/2,z
146 x+1/2,-y+3/4,-z+1/4
147 -x+3/4,y+1/2,-z+1/4
148 -x+3/4,-y+3/4,z
149 y+1/2,z+1/2,x
150 y+1/2,-z+3/4,-x+1/4
151 -y+3/4,z+1/2,-x+1/4
152 -y+3/4,-z+3/4,x
153 z+1/2,x+1/2,y
154 z+1/2,-x+3/4,-y+1/4
155 -z+3/4,x+1/2,-y+1/4
156 -z+3/4,-x+3/4,y
157 -y+1/2,-x+1/2,-z
158 -y+1/2,x+3/4,z+1/4
159 y+3/4,-x+1/2,z+1/4
160 y+3/4,x+3/4,-z
161 -x+1/2,-z+1/2,-y
162 -x+1/2,z+3/4,y+1/4
163 x+3/4,-z+1/2,y+1/4
164 x+3/4,z+3/4,-y
165 -z+1/2,-y+1/2,-x
166 -z+1/2,y+3/4,x+1/4
167 z+3/4,-y+1/2,x+1/4
168 z+3/4,y+3/4,-x
169 -x+1/2,-y+1/2,-z
170 -x+1/2,y+3/4,z+1/4
171 x+3/4,-y+1/2,z+1/4
172 x+3/4,y+3/4,-z
173 -y+1/2,-z+1/2,-x
174 -y+1/2,z+3/4,x+1/4
175 y+3/4,-z+1/2,x+1/4
176 y+3/4,z+3/4,-x
177 -z+1/2,-x+1/2,-y
178 -z+1/2,x+3/4,y+1/4
179 z+3/4,-x+1/2,y+1/4
180 z+3/4,x+3/4,-y
181 y+1/2,x+1/2,z
182 y+1/2,-x+3/4,-z+1/4
183 -y+3/4,x+1/2,-z+1/4
184 -y+3/4,-x+3/4,z
185 x+1/2,z+1/2,y
186 x+1/2,-z+3/4,-y+1/4
187 -x+3/4,z+1/2,-y+1/4
188 -x+3/4,-z+3/4,y
189 z+1/2,y+1/2,x
190 z+1/2,-y+3/4,-x+1/4
191 -z+3/4,y+1/2,-x+1/4
192 -z+3/4,-y+3/4,x

loop_
_atom_site_label
_atom_site_type_symbol
_atom_site_symmetry_multiplicity
_atom_site_Wyckoff_label
_atom_site_fract_x
_atom_site_fract_y
_atom_site_fract_z
_atom_site_occupancy
Cu1 Cu 16 c 0.00000 0.00000 0.00000 1.00000
Pt1 Pt 16 d 0.50000 0.50000 0.50000 1.00000
```

Cubic CuPt ($L1_3$ (I), $D4$): AB_cF32_227_c_d - POSCAR

```
AB_cF32_227_c_d & a --params=7.23 & Fd-3m O_{h}^{7} #227 (cd) & cF32 &
    ↪ $L1_{3}$ (I) & CuPt & CuPt & C. H. Johansson and J. O. Linde,
    ↪ Ann. Phys. 387, 449-478 (1927)
   1.00000000000000
   0.00000000000000   3.61500000000000   3.61500000000000
   3.61500000000000   0.00000000000000   3.61500000000000
   3.61500000000000   3.61500000000000   0.00000000000000
    Cu    Pt
     4     4
Direct
   0.00000000000000   0.00000000000000   0.00000000000000   Cu   (16c)
   0.00000000000000   0.00000000000000   0.50000000000000   Cu   (16c)
   0.00000000000000   0.50000000000000   0.00000000000000   Cu   (16c)
   0.50000000000000   0.00000000000000   0.00000000000000   Cu   (16c)
   0.50000000000000   0.50000000000000   0.50000000000000   Pt   (16d)
   0.50000000000000   0.50000000000000   0.00000000000000   Pt   (16d)
   0.50000000000000   0.00000000000000   0.50000000000000   Pt   (16d)
   0.00000000000000   0.50000000000000   0.50000000000000   Pt   (16d)
```

α-AgI ($B23$): A21B_cI44_229_bdh_a - CIF

```
# CIF file
data_findsym-output
_audit_creation_method FINDSYM

_chemical_name_mineral 'AgI'
_chemical_formula_sum 'Ag21 I'

loop_
_publ_author_name
 'L. W. Strock'
_journal_name_full_name
;
 Zeitschrift f{\"u}r Physikalische Chemie B
;
_journal_volume 25
_journal_year 1934
_journal_page_first 441
_journal_page_last 459
_publ_Section_title
```

```
;
 Kristallstruktur des Hochtemperatur-Jodsilbers $\alpha$-AgJ
;

# Found in Crystal Structure and Phase Transition of Some Metallic
      ↪ Halides IV On the Anomalous Structure of $\alpha$-AgI, 1957

_aflow_title '$\alpha$-AgI ($B23$) Structure'
_aflow_proto 'A21B_cI44_229_bdh_a'
_aflow_params 'a,y_{4}'
_aflow_params_values '5.034,0.375'
_aflow_Strukturbericht '$B23$'
_aflow_Pearson 'cI44'

_symmetry_space_group_name_H-M "I 4/m -3 2/m"
_symmetry_Int_Tables_number 229

_cell_length_a     5.03400
_cell_length_b     5.03400
_cell_length_c     5.03400
_cell_angle_alpha  90.00000
_cell_angle_beta   90.00000
_cell_angle_gamma  90.00000

loop_
_space_group_symop_id
_space_group_symop_operation_xyz
1 x,y,z
2 x,-y,-z
3 -x,y,-z
4 -x,-y,z
5 y,z,x
6 y,-z,-x
7 -y,z,-x
8 -y,-z,x
9 z,x,y
10 z,-x,-y
11 -z,x,-y
12 -z,-x,y
13 -y,-x,-z
14 -y,x,z
15 y,-x,z
16 y,x,-z
17 -x,-z,-y
18 -x,z,y
19 x,-z,y
20 x,z,-y
21 -z,-y,-x
22 -z,y,x
23 z,-y,x
24 z,y,-x
25 -x,-y,-z
26 -x,y,z
27 x,-y,z
28 x,y,-z
29 -y,-z,-x
30 -y,z,x
31 y,-z,x
32 y,z,-x
33 -z,-x,-y
34 -z,x,y
35 z,-x,y
36 z,x,-y
37 y,x,z
38 y,-x,-z
39 -y,x,-z
40 -y,-x,z
41 x,z,y
42 x,-z,-y
43 -x,z,-y
44 -x,-z,y
45 z,y,x
46 z,-y,-x
47 -z,y,-x
48 -z,-y,x
49 x+1/2,y+1/2,z+1/2
50 x+1/2,-y+1/2,-z+1/2
51 -x+1/2,y+1/2,-z+1/2
52 -x+1/2,-y+1/2,z+1/2
53 y+1/2,z+1/2,x+1/2
54 y+1/2,-z+1/2,-x+1/2
55 -y+1/2,z+1/2,-x+1/2
56 -y+1/2,-z+1/2,x+1/2
57 z+1/2,x+1/2,y+1/2
58 z+1/2,-x+1/2,-y+1/2
59 -z+1/2,x+1/2,-y+1/2
60 -z+1/2,-x+1/2,y+1/2
61 -y+1/2,-x+1/2,-z+1/2
62 -y+1/2,x+1/2,z+1/2
63 y+1/2,-x+1/2,z+1/2
64 y+1/2,x+1/2,-z+1/2
65 -x+1/2,-z+1/2,-y+1/2
66 -x+1/2,z+1/2,y+1/2
67 x+1/2,-z+1/2,y+1/2
68 x+1/2,z+1/2,-y+1/2
69 -z+1/2,-y+1/2,-x+1/2
70 -z+1/2,y+1/2,x+1/2
71 z+1/2,-y+1/2,x+1/2
72 z+1/2,y+1/2,-x+1/2
73 -x+1/2,-y+1/2,-z+1/2
74 -x+1/2,y+1/2,z+1/2
75 x+1/2,-y+1/2,z+1/2
76 x+1/2,y+1/2,-z+1/2
77 -y+1/2,-z+1/2,-x+1/2
78 -y+1/2,z+1/2,x+1/2
79 y+1/2,-z+1/2,x+1/2
80 y+1/2,z+1/2,-x+1/2
81 -z+1/2,-x+1/2,-y+1/2
82 -z+1/2,x+1/2,y+1/2
83 z+1/2,-x+1/2,y+1/2
84 z+1/2,x+1/2,-y+1/2
85 y+1/2,x+1/2,z+1/2
86 y+1/2,-x+1/2,-z+1/2
87 -y+1/2,x+1/2,-z+1/2
88 -y+1/2,-x+1/2,z+1/2
89 x+1/2,z+1/2,y+1/2
90 x+1/2,-z+1/2,-y+1/2
91 -x+1/2,z+1/2,-y+1/2
92 -x+1/2,-z+1/2,y+1/2
93 z+1/2,y+1/2,x+1/2
94 z+1/2,-y+1/2,-x+1/2
95 -z+1/2,y+1/2,-x+1/2
96 -z+1/2,-y+1/2,x+1/2

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_atom_site_fract_z
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Ag2 Ag 12 d 0.25000 0.00000 0.50000 0.04762
Ag3 Ag 24 h 0.00000 0.37500 0.37500 0.04762
```

α-AgI ($B23$): A21B_cI44_229_bdh_a - POSCAR

```
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      ↪ abdh) & cI44 & $B23$ & AgI & AgI & L. W. Strock, Z. Physik.
      ↪ Chem. B 25, 441-459 (1934)
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  -2.51700000000000    2.51700000000000    2.51700000000000
   2.51700000000000   -2.51700000000000    2.51700000000000
   2.51700000000000    2.51700000000000   -2.51700000000000
    Ag   I
    21   1
Direct
   0.00000000000000    0.50000000000000    0.50000000000000   Ag   (6b)
   0.50000000000000    0.00000000000000    0.50000000000000   Ag   (6b)
   0.50000000000000    0.50000000000000    0.00000000000000   Ag   (6b)
   0.50000000000000    0.75000000000000    0.25000000000000   Ag  (12d)
   0.50000000000000    0.25000000000000    0.75000000000000   Ag  (12d)
   0.25000000000000    0.50000000000000    0.75000000000000   Ag  (12d)
   0.75000000000000    0.50000000000000    0.25000000000000   Ag  (12d)
   0.75000000000000    0.25000000000000    0.50000000000000   Ag  (12d)
   0.25000000000000    0.75000000000000    0.50000000000000   Ag  (12d)
   0.75000000000000    0.37500000000000    0.37500000000000   Ag  (24h)
   0.00000000000000    0.37500000000000   -0.37500000000000   Ag  (24h)
   0.00000000000000   -0.37500000000000    0.37500000000000   Ag  (24h)
  -0.75000000000000   -0.37500000000000   -0.37500000000000   Ag  (24h)
   0.37500000000000    0.75000000000000    0.37500000000000   Ag  (24h)
  -0.37500000000000    0.00000000000000    0.37500000000000   Ag  (24h)
   0.37500000000000    0.00000000000000   -0.37500000000000   Ag  (24h)
  -0.37500000000000   -0.75000000000000   -0.37500000000000   Ag  (24h)
   0.37500000000000    0.37500000000000    0.75000000000000   Ag  (24h)
   0.37500000000000   -0.37500000000000    0.00000000000000   Ag  (24h)
  -0.37500000000000    0.37500000000000    0.00000000000000   Ag  (24h)
  -0.37500000000000   -0.37500000000000   -0.75000000000000   Ag  (24h)
   0.00000000000000    0.00000000000000    0.00000000000000    I   (2a)
```

$Ca_3Al_2(OH)_{12}$ ($J2_3$): A2B3C12D12_cI232_230_a_c_h_h - CIF

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;
_journal_volume 324
_journal_year 1986
_journal_page_first 124
_journal_page_last 126
_publ_Section_title
;
 Tricalciumaluminathexahydrat, Ca$_{3}$[Al(OH)$_{6}$]$_{2}$, Bindungsl
      ↪ \"{a}ngen und -valenzen aus R\"{o}ntgeneinkristallmessungen
;

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_aflow_Strukturbericht '$J2_{3}$'
_aflow_Pearson 'cI232'

_symmetry_space_group_name_H-M "I 41/a -3 2/d"
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4 -x,-y+1/2,z
5 y,z,x
6 y,-z,-x+1/2
7 -y+1/2,z,-x
8 -y,-z+1/2,x
9 z,x,y
10 z,-x,-y+1/2
11 -z+1/2,x,-y
12 -z,-x+1/2,y
13 -y+1/4,-x+1/4,-z+1/4
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19 x+1/4,-z+1/4,y+3/4
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21 -z+1/4,-y+1/4,-x+1/4
22 -z+1/4,y+3/4,x+1/4
23 z+1/4,-y+1/4,x+3/4
24 z+3/4,y+1/4,-x+1/4
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44 -x+3/4,-z+1/4,y+1/4
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66 -x+3/4,z+1/4,y+3/4
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74 -x+1/2,y+1/2,z
75 x,-y+1/2,z+1/2
76 x+1/2,y,-z+1/2
77 -y+1/2,-z+1/2,-x+1/2
78 -y+1/2,z+1/2,x
79 y,-z+1/2,x+1/2
80 y+1/2,z,-x+1/2
81 -z+1/2,-x+1/2,-y+1/2
82 -z+1/2,x+1/2,y
83 z,-x+1/2,y+1/2
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85 y+3/4,x+3/4,z+3/4
86 y+3/4,-x+1/4,-z+3/4
87 -y+3/4,x+3/4,-z+1/4
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91 -x+3/4,z+3/4,-y+1/4
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93 z+3/4,y+3/4,x+3/4
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Ca1 Ca 24 c 0.12500 0.00000 0.25000 1.00000
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O1 O 96 h -0.0282 0.05310 0.13960 1.00000
```

$Ca_3Al_2(OH)_{12}$ ($J2_3$): A2B3C12D12_cI232_230_a_c_h_h - POSCAR

```
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    ↪ 0.091,0.054,0.163,-0.0282,0.0531,0.1396 & Ia-3d O_{h}^{10} #230
    ↪ (ach^2) & cI232 & $J2_{3}$ & Al2Ca3H12O12 & Al2Ca3H12O12 & H.
    ↪ Bartl, Fresen. Z. Anal. Chem. 324, 124-126 (1986)
   1.00000000000000
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   6.28750000000000   -6.28750000000000    6.28750000000000
   6.28750000000000    6.28750000000000   -6.28750000000000
    Al    Ca    H    O
     8    12    48    48
Direct
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    0.50000000000000    0.00000000000000    0.50000000000000   Al   (16a)
    0.00000000000000    0.50000000000000    0.50000000000000   Al   (16a)
    0.50000000000000    0.50000000000000    0.00000000000000   Al   (16a)
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    0.60900000000000    0.25400000000000    0.53700000000000   H    (96h)
   -0.10900000000000    0.42800000000000    0.64500000000000   H    (96h)
    0.28300000000000    0.24600000000000   -0.14500000000000   H    (96h)
   -0.03700000000000    0.21700000000000    0.07200000000000   H    (96h)
    0.53700000000000    0.60900000000000    0.25400000000000   H    (96h)
    0.64500000000000   -0.10900000000000    0.42800000000000   H    (96h)
   -0.14500000000000    0.28300000000000    0.24600000000000   H    (96h)
    0.07200000000000   -0.03700000000000    0.21700000000000   H    (96h)
    0.25400000000000    0.53700000000000    0.60900000000000   H    (96h)
    0.42800000000000    0.64500000000000   -0.10900000000000   H    (96h)
    0.24600000000000   -0.14500000000000    0.28300000000000   H    (96h)
    0.24600000000000   -0.10900000000000   -0.03700000000000   H    (96h)
    0.42800000000000    0.28300000000000    0.53700000000000   H    (96h)
    0.25400000000000    0.21700000000000    0.64500000000000   H    (96h)
    0.07200000000000    0.60900000000000   -0.14500000000000   H    (96h)
    0.60900000000000   -0.14500000000000    0.07200000000000   H    (96h)
    0.21700000000000    0.64500000000000    0.25400000000000   H    (96h)
    0.28300000000000    0.53700000000000    0.42800000000000   H    (96h)
   -0.10900000000000   -0.03700000000000    0.24600000000000   H    (96h)
    0.64500000000000    0.25400000000000    0.21700000000000   H    (96h)
   -0.14500000000000    0.07200000000000    0.60900000000000   H    (96h)
   -0.03700000000000    0.24600000000000   -0.10900000000000   H    (96h)
    0.53700000000000    0.42800000000000    0.28300000000000   H    (96h)
   -0.21700000000000   -0.07200000000000    0.03700000000000   H    (96h)
    0.39100000000000   -0.25400000000000    0.46300000000000   H    (96h)
    0.10900000000000    0.57200000000000    0.35500000000000   H    (96h)
    0.71700000000000    0.75400000000000    0.14500000000000   H    (96h)
    0.03700000000000   -0.21700000000000   -0.07200000000000   H    (96h)
    0.46300000000000    0.39100000000000   -0.25400000000000   H    (96h)
    0.35500000000000    0.10900000000000    0.57200000000000   H    (96h)
    0.14500000000000    0.71700000000000    0.75400000000000   H    (96h)
   -0.07200000000000    0.03700000000000   -0.21700000000000   H    (96h)
   -0.25400000000000    0.46300000000000    0.39100000000000   H    (96h)
    0.57200000000000    0.35500000000000    0.10900000000000   H    (96h)
    0.75400000000000    0.14500000000000    0.71700000000000   H    (96h)
    0.75400000000000    0.10900000000000    0.03700000000000   H    (96h)
    0.57200000000000    0.71700000000000    0.46300000000000   H    (96h)
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   -0.07200000000000    0.39100000000000    0.14500000000000   H    (96h)
    0.39100000000000    0.14500000000000   -0.07200000000000   H    (96h)
   -0.21700000000000    0.35500000000000   -0.25400000000000   H    (96h)
    0.71700000000000    0.46300000000000    0.57200000000000   H    (96h)
    0.10900000000000    0.03700000000000    0.75400000000000   H    (96h)
    0.35500000000000   -0.25400000000000   -0.21700000000000   H    (96h)
    0.14500000000000   -0.07200000000000    0.39100000000000   H    (96h)
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    0.19270000000000    0.11140000000000    0.02490000000000   O    (96h)
    0.58650000000000    0.16780000000000    0.47510000000000   O    (96h)
   -0.08650000000000    0.38860000000000    0.58130000000000   O    (96h)
    0.30730000000000    0.33220000000000   -0.08130000000000   O    (96h)
    0.02490000000000    0.19270000000000    0.11140000000000   O    (96h)
    0.47510000000000    0.58650000000000    0.16780000000000   O    (96h)
    0.58130000000000   -0.08650000000000    0.38860000000000   O    (96h)
   -0.08130000000000    0.30730000000000    0.33220000000000   O    (96h)
    0.11140000000000    0.02490000000000    0.19270000000000   O    (96h)
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0.38860000000000	0.58130000000000	−0.08650000000000	O	(96h)
0.33220000000000	−0.08130000000000	0.30730000000000	O	(96h)
0.33220000000000	−0.08650000000000	0.02490000000000	O	(96h)
0.38860000000000	0.30730000000000	0.47510000000000	O	(96h)
0.16780000000000	0.19270000000000	0.58130000000000	O	(96h)
0.11140000000000	0.58650000000000	−0.08130000000000	O	(96h)
0.58650000000000	−0.08130000000000	0.11140000000000	O	(96h)
0.19270000000000	0.58130000000000	0.16780000000000	O	(96h)
0.30730000000000	0.47510000000000	0.38860000000000	O	(96h)
−0.08650000000000	0.02490000000000	0.33220000000000	O	(96h)
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−0.08130000000000	0.11140000000000	0.58650000000000	O	(96h)
0.02490000000000	0.33220000000000	−0.08650000000000	O	(96h)
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−0.19270000000000	−0.11140000000000	−0.02490000000000	O	(96h)
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0.69270000000000	0.66780000000000	0.08130000000000	O	(96h)
−0.02490000000000	−0.19270000000000	−0.11140000000000	O	(96h)
0.52490000000000	0.41350000000000	−0.16780000000000	O	(96h)
0.41870000000000	0.08650000000000	0.61140000000000	O	(96h)
0.08130000000000	0.69270000000000	0.66780000000000	O	(96h)
−0.11140000000000	−0.02490000000000	−0.19270000000000	O	(96h)
−0.16780000000000	0.52490000000000	0.41350000000000	O	(96h)
0.61140000000000	0.41870000000000	0.08650000000000	O	(96h)
0.66780000000000	0.08130000000000	0.69270000000000	O	(96h)
0.66780000000000	0.08650000000000	−0.02490000000000	O	(96h)
0.61140000000000	0.69270000000000	0.52490000000000	O	(96h)
−0.16780000000000	−0.19270000000000	0.41870000000000	O	(96h)
−0.11140000000000	0.41350000000000	0.08130000000000	O	(96h)
0.41350000000000	0.08130000000000	−0.11140000000000	O	(96h)
−0.19270000000000	0.41870000000000	−0.16780000000000	O	(96h)
0.69270000000000	0.52490000000000	0.61140000000000	O	(96h)
0.08650000000000	−0.02490000000000	0.66780000000000	O	(96h)
0.41870000000000	−0.16780000000000	−0.19270000000000	O	(96h)
0.08130000000000	−0.11140000000000	0.41350000000000	O	(96h)
−0.02490000000000	0.66780000000000	0.08650000000000	O	(96h)
0.52490000000000	0.61140000000000	0.69270000000000	O	(96h)

Prototype Index

1. $B30$ (MgZn?): AB_oI48_44_6d_ab2cde 520
2. $C17$ (Fe_2B) (*obsolete*): AB2_tI12_121_ab_i 884
3. $C2$ ($Ba,Ca)CO_3$: ABC3_mC10_5_b_a_ac 92
4. $C26_a$ (NO_2) (*obsolete*): AB2_cI36_199_b_c 1282
5. $C27$ (CdI_2) (*questionable*): AB2_hP6_186_b_ab 1213
6. $C53$ ($SrBr_2$) (*obsolete*): A2B_oP12_62_2c_c 654
7. $D0_{10}$ (WO_3) (*obsolete*): A3B_oP16_57_a2d_d 566
8. $D0_{13}$ ($AlCl_3$) (*obsolete*): AB3_hP4_164_b_ad 1107
9. $D0_{15}$ ($AlCl_3$) (*obsolete*): AB3_mC16_5_c_3c 87
10. $D0_6$ (Tysonite, LaF_3) (*obsolete*): A3B_hP24_193_ack_g 1238
11. $D0_7$ (CrO_3) (*obsolete*): AB3_oC16_20_a_bc 409
12. $D2_2$ ($MgZn_5$?) (*problematic*): AB5_mC48_12_2i_ac5i2j 175
13. $D6_2$ (Sb_2O_4) (*obsolete*): A2B_cF96_227_abf_cd 1478
14. $D8_7$ (Shcherbinaite, V_2O_5) (*obsolete*): A5B2_oP14_31_a2b_b 440
15. $E2_3$ ($LiIO_3$) (*obsolete*): ABC3_hP10_182_c_b_g 1202
16. $E3_1$ (β-Ag_2HgI_4) (*obsolete*): A2BC4_tP7_111_f_a_n 861
17. $E6_1$ ($Sr(OH)_2(H_2O)_8$) (*obsolete*): A8B2C_tP11_123_r_f_a 906
18. $E6_2$ [$SrO_2(H_2O)_8$] (*possibly obsolete*): A8B2C_tP11_123_r_h_a 908
19. $F5_{11}$ (KNO_2) (*obsolete*): ABC2_mC8_8_a_a_b 117
20. $F5_4$ (NH_4ClO_2) (*obsolete*): ABC2_tP8_100_b_a_c 851
21. $F6_1$ (Chalcopyrite, $CuFeS_2$) (*obsolete*): ABC2_tP4_115_a_c_g 872
22. $G7_3$ [Northupite, $Na_3MgCl(CO_3)_2$] (*obsolete*): A2BCD3E6_cF208_227_e_c_d_f_g 1475
23. $G7_5$ ($PbCO_3 \cdot PbCl_2$, Phosgenite) (*obsolete*): AB2C3D2_tP16_90_c_f_ce_e 839
24. $H5_6$ [Tychite, $Na_6Mg_2SO_4(CO_3)_4$)] (*obsolete*): A4B2C6D16E_cF232_227_e_d_f_eg_a 1484
25. $H5_9$ [Autunite, $Ca(UO_2)_2(PO_4)_2 \cdot 10\frac{1}{2}H_2O$] (*obsolete*)[⊗‡‡]: AB2C2_tI10_139_a_d_e 994
26. $H6_4$ [$Ni(NO_3)_2(NH_3)_6$] (*obsolete*)[⊗⊗]: A2B6CD6_cP60_205_c_d_a_d 1302
27. $I1_3$ ($SrCl_2 \cdot (H_2O)_6$) (*obsolete*)[§§]: A2B6C_hP9_162_d_k_a 1083
28. $L1_a$ (disputed $CuPt_3$): AB7_cF32_225_b_ad 1451
29. $S0_4$ (Staurolite, $Fe(OH)_2Al_4Si_2O_{10}$) (*obsolete*): A4BC12D2_oC76_63_eg_c_f3gh_g 746
30. $S3_4$ (II) (Catapleiite, $Na_2Zr(SiO_3)_3 \cdot H_2O$) (*obsolete*): A3B2C9D3E_hP36_194_g_f_hk_h_a 1254
31. α-AgI ($B23$): A21B_cI44_229_bdh_a 1490
32. α-Alum [$KAl(SO_4)_2 \cdot 12H_2O$, $H4_{13}$]: AB24CD28E2_cP224_205_a_4d_b_2c4d_c 1335
33. α-BaB_2O_4 (Low-Temperature): A2BC4_hR42_161_2b_b_4b 1078
34. α-Carnegieite ($NaAlSiO_4$, $S6_5$): ABC4D_cP28_198_a_a_ab_a 1280
35. α-$Ho_2Si_2O_7$: A2B7C2_aP44_2_4i_14i_4i 47
36. α-ICl[§]: AB_mP16_14_2e_2e 299
37. α-$LiIO_3$: ABC3_hP10_173_b_a_c 1178
38. α-PbO_2[×]: A2B_oP12_60_d_c 602
39. α-Potassium Nitrate (KNO_3) I[°]: ABC3_oP20_62_c_c_cd 720
40. α-Potassium Nitrate (KNO_3) II: ABC3_oC80_36_2ab_2ab_2a5b 478
41. α-V_3S: AB3_tI32_121_g_f2i 888
42. α-WO_3: A3B_tP16_130_cf_c 957

[⊗] Li_2CN_2 and $H5_9$ [Autunite, $Ca(UO_2)_2(PO_4)_2 \cdot 10\frac{1}{2}H_2O$] (*obsolete*) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[‡‡] $TlCo_2S_2$, $ThCr_2Si_2$, Li_2CN_2, and $H5_9$ [Autunite, $Ca(UO_2)_2(PO_4)_2 \cdot 10\frac{1}{2}H_2O$] (*obsolete*) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

[⊗⊗] $Zn(BrO_3)_2 \cdot 6H_2O$ ($J1_{10}$) and $H6_4$ [$Ni(NO_3)_2(NH_3)_6$] (*obsolete*) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

[§§] $I1_3$ ($SrCl_2 \cdot (H_2O)_6$) (*obsolete*) and Rosiaite ($PbSb_2O_6$) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

[§] α-ICl and LiAs have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[×] ζ-Fe_2N and α-PbO_2 have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[°] α-Potassium Nitrate (KNO_3) I, NH_4NO_3 III ($G0_{10}$), and Aragonite ($CaCO_3$, $G0_2$) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

43. α-$Zn_2V_2O_7$: A7B2C2_mC44_15_e3f_f_f 354
44. β-Alum [$Al(NH_3CH_3)_2(SO_4)_2{\cdot}12H_2O$, $H4_{14}$]: AB2C36D2E20F2_cP252_205_a_c_6d_c_c3d_c ... 1344
45. β-Alumina (Al_2O_3, $D5_6$): A2B3_hP60_194_3fk_cdef2k 1251
46. β-Arabinose [$(CH_2O)_{20}$]: AB2C_oP80_19_5a_10a_5a 401
47. β-B_2H_6[*]: AB3_mP16_14_e_3e 264
48. β-BaB_2O_4 (High-Temperature): A2BC4_hR42_167_f_ac_2f 1151
49. β-Ga (*obsolete*): A_mC4_15_e 381
50. β-Ga_2O_3: A2B3_mC20_12_2i_3i 151
51. β-$LiIO_3$: ABC3_tP40_86_g_g_3g 822
52. β-Potassium Nitrate (KNO_3): ABC6_hR8_166_a_b_h 1138
53. β-Si_3N_4: A4B3_hP14_176_ch_h 1187
54. δ-CuTi ($L2_a$): AB_tP2_123_a_d 912
55. δ-Ni_3Sn_4 ($D7_a$): A3B4_mC14_12_ai_2i 159
56. δ-WO_3: A3B_aP32_2_12i_4i 54
57. ϵ-1,2,3,4,5,6-Hexachlorocyclohexane (C_6Cl_6): AB_mP24_14_3e_3e 303
58. η-NiSi (B_d): AB_oP8_62_c_c 738
59. γ-Alum [$AlNa(SO_4)_2{\cdot}12H_2O$, $H4_{15}$]: AB24CD20E2_cP192_205_a_4d_b_c3d_c 1327
60. γ-Fe_4N ($L'1_0$): A4B_cP5_221_bc_a 1414
61. γ-Ga_2O_3: A11B4_cF120_227_acdf_e 1455
62. γ-$LiIO_3$: ABC3_oP20_33_a_a_3a 462
63. γ-Potassium Nitrate (KNO_3)[□]: ABC3_hR5_160_a_a_b 1076
64. γ-TeO_2: A2B_oP12_18_2c_c 386
65. γ-WO_3: A3B_mP32_14_6e_2e 226
66. γ-$Y_2Si_2O_7$[‖]: A4BC_mP24_14_4e_e_e 235
67. ζ-Fe_2N[×]: A2B_oP12_60_d_c 600
68. ζ-Nb_2O_5 (B-Nb_2O_5): A2B5_mC28_15_f_e2f 327
69. y-$Y_2Si_2O_7$: A7B2C2_mP22_11_3e2f_2e_ab 135
70. $(CdSO_4)_3{\cdot}8H_2O$ ($H4_{20}$): A3B16C20D3_mC168_15_ef_8f_10f_ef 342
71. $(NH_4)_3AlF_6$ ($J2_1$): AB30C16D3_cF200_225_a_ej_2f_bc 1448
72. $(TiCl_4{\cdot}POCl_3)_2$: A7BCD_oP80_61_7c_c_c_c 616
73. 12-phosphotungstic acid [$H_3PW_{12}O_{40}{\cdot}5H_2O$ ($H4_{16}$)]: A5B40CD12_cP116_224_cd_e3k_a_k 1434
74. Adamite [$Zn_2(AsO_4)(OH)$, $H2_7$]: ABC5D2_oP36_58_g_g_3gh_eg 586
75. Ag_2O_3[**]: A2B3_oF40_43_b_ab 497
76. Ag_2PbO_2: A2B2C_mC20_15_ad_f_e 312
77. $Ag_2SO_4{\cdot}4NH_3$ ($H4_{17}$): A2B12C4D4E_tP46_114_d_3e_e_e_a 869
78. $AgMnO_4$ ($H0_9$)[‖]: ABC4_mP24_14_e_e_4e 289
79. $AgNO_2$ ($F5_{12}$): ABC2_oI8_44_a_a_d 518
80. $Ag[Co(NH_3)_2(NO_2)_4]$ ($J1_9$): ABC4D2E8_tP32_126_a_b_h_e_k 926
81. $Al_{13}Fe_4$: A13B4_mC102_12_dg8i5j_4ij 141
82. Al_2Mo_3C: A2BC3_cP24_213_c_a_d 1358
83. $Al(PO_3)_3$ ($G5_2$): AB9C3_cI208_220_c_3e_e 1405
84. AlN (cF40): AB_cF40_216_de_ce 1382
85. AlN (cI16): AB_cI16_217_c_c 1387
86. AlN (cI24): AB_cI24_220_a_b 1412
87. $AlNbO_4$: ABC4_mC24_12_i_i_4i 191
88. $AlPO_4$ "low cristobalite type": AB4C_oC24_20_b_2c_a 411
89. Albite ($NaAlSi_3O_8$, $S6_8$): ABC8D3_aP26_2_i_i_8i_3i 68
90. Alluaudite [$NaMnFe_2(PO_4)_3$]: A2BCD12E3_mC76_15_f_e_b_6f_ef 334
91. Ammonium Chlorite (NH_4ClO_2): AB4CD2_tP16_113_c_f_a_e 863
92. Ammonium Persulfate [$(NH_4)_2S_2O_8$, $K4_1$][†‖]: AB4C_mP24_14_e_4e_e 274
93. Analcime ($NaAlSi_2O_6{\cdot}H_2O$, $S6_1$): A2B2C3D12E4_tI184_142_f_f_be_3g_g 1019
94. Andalusite (Al_2SiO_5, $S0_2$): A2B5C_oP32_58_eg_3gh_g 576
95. Anhydrous $KAuBr_4$: AB4C_mP24_14_ab_4e_e ... 271
96. Anthophyllite ($Mg_5Fe_2Si_8O_{22}(OH)_2$, $S4_4$): A2B5C22D2E8_oP156_62_d_c2d_2c10d_2c_4d 638
97. Apophyllite ($KCa_4Si_8O_{20}F{\cdot}8H_2O$, $S5_2$): A4BC16DE28F8_tP116_128_h_a_2i_b_g3i_i 937
98. Aragonite ($CaCO_3$, $G0_2$)[○]: ABC3_oP20_62_c_c_cd 726
99. Arcanite (K_2SO_4, $H1_6$): A2B4C_oP28_62_2c_2cd_c 635
100. Archerite (KH_2PO_4): A2BC4D_oF64_43_b_a_2b_a 507
101. Arsenopyrite (FeAsS, $E0_7$): ABC_mP12_14_e_e_e 297

[*] β-B_2H_6 and B_2H_6 ($P2_1/c$) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[□] $KBrO_3$ ($G0_7$) and γ-Potassium Nitrate (KNO_3) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[‖] γ-$Y_2Si_2O_7$, Ammonium Persulfate [$(NH_4)_2S_2O_8$, $K4_1$], Monazite ($LaPO_4$), and $AgMnO_4$ ($H0_9$) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

[**] Ag_2O_3 and Zr_2Al_3 have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

[†] Ammonium Persulfate [$(NH_4)_2S_2O_8$, $K4_1$] and Monazite ($LaPO_4$) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

102. Atacamite ($Cu_2(OH)_3Cl$): AB2C3D3_oP36_62_c_ac_cd_cd 704
103. Au_2Nb_3: A2B3_tI10_139_e_ae 976
104. $AuCsCl_3$ ($K7_6$): AB3C_tI20_139_ab_eh_d 996
105. Autunite $\{Ca[(UO_2)(PO_4)]_2(H_2O)_{11}\}$: AB22C23D2E2_oP200_62_c_11d_3c10d_d_d 695
106. Azurite [$Cu_3(CO_3)_2(OH)_2$, $G7_4$]: A2B3C2D8_mP30_14_e_ce_e_4e 205
107. $B_{13}C_2$ "B_4C" ($D1_g$): A13B2_hR15_166_b2h_c .. 1111
108. B_2H_6 ($P2_1/c$)[*]: AB3_mP16_14_e_3e 266
109. B_4SrO_7: A4B7C_oP24_31_2b_a3b_a 438
110. $BaAl_2O_4$ ($H2_8$): A2BC6_hP18_182_f_b_gh 1200
111. $BaCd_{11}$: AB11_tI48_141_a_bdi 1016
112. $BaNi(CN)_4{\cdot}4H_2O$ ($H4_{22}$): AB4C4D4E_mC56_15_e_2f_2f_2f_a 363
113. $BaNiSn_3$: ABC3_tI10_107_a_a_ab 859
114. Bararite (Trigonal $(NH_4)_2SiF_6$, $J1_6$)[××]: A6B2C_hP9_164_i_d_a 1101
115. Barytocalcite ($BaCa(CO_3)_2$): AB2CD6_mP20_11_e_2e_e_2e2f 137
116. Base-centered orthorhombic $Sr_4Ru_3O_{10}$: A10B3C4_oC68_64_2dfg_ad_2d 769
117. Bassanite [$CaSO_4(H_2O)_{0.5}$, $H4_7$]: A2B2C9D2_mC90_5_ab2c_3c_b13c_3c 81
118. Bastnäsite [$CeF(CO_3)$]: ABCD3_hP36_190_h_g_af_hi 1230
119. $BeSO_4{\cdot}4H_2O$ ($H4_3$): AB8C8D_tI72_120_c_2i_2i_b 879
120. Berthierite ($FeSb_2S_4$, $E3_3$): AB4C2_oP28_62_c_4c_2c 709
121. Bertrandite ($Be_4Si_2O_7(OH)_2$, $S4_6$): A4B7C2D2_oC60_36_2b_a3b_2a_b 473
122. Bi_2GeO_5: A2BC5_oC32_36_b_a_a2b 468
123. $Bi_3Ru_3O_{11}$: A3B11C3_cP68_201_be_efh_g 1285
124. Bischofite ($MgCl_2{\cdot}6H_2O$, $J1_7$): A2B12CD6_mC42_12_i_2i2j_a_ij 146
125. Blossite (α-$Cu_2V_2O_7$): A2B7C2_oF88_43_b_a3b_b 504
126. Boric Acid (H_3BO_3, $G5_1$): AB3C3_aP28_2_2i_6i_6i 62
127. Bromocarnallite ($KMg(H_2O)_6(Cl,Br)_3$, $E2_6$): A3B6CD_tP44_85_bcg_3g_ac_e 807
128. Brucite [$Mg(OH)_2$][¶¶]: A2BC2_hP5_164_d_a_d .. 1097
129. $C_{19}Sc_{15}$: A19B15_tP68_114_bc4e_ac3e 865
130. COCl: ABC_oP24_61_c_c_c 627
131. Ca_2RuO_4: A2B4C_oP28_61_c_2c_a 611
132. Ca_2UO_5: A2B5C_mP32_14_2e_5e_ab 218
133. $Ca_3Al_2(OH)_{12}$ ($J2_3$): A2B3C12D12_cI232_230_a_c_h_h 1492
134. Ca_3UO_6: A3B6C_mP20_4_3a_6a_a 79
135. CaB_2O_4 (III): A2BC4_oP84_33_6a_3a_12a 450
136. CaB_2O_4 (IV): A2BC4_cP84_205_d_ac_2d 1309
137. CaB_2O_4 I ($E3_2$): A2BC4_oP28_60_d_c_2d 597
138. $CaBe_2Ge_2$: A2BC2_tP10_129_ac_c_bc 943
139. CaC_2-I ($C11_a$): A2B_tI6_139_e_a 978
140. CaC_2-III: A2B_mC12_12_2i_i 155
141. $CaCu_4P_2$[°°]: AB4C2_hR7_166_a_2c_c 1129
142. $CaO_2(H_2O)_8$: AB8C2_tP22_124_a_n_h 914
143. $CaSi_2$ ($C12$): AB2_hR6_166_c_2c 1124
144. $CaUO_4$: AB4C_hR6_166_b_2c_a 1131
145. Calaverite ($AuTe_2$): AB2_mP12_7_2a_4a 107
146. Calciborite (CaB_2O_4 II): A2BC4_oP56_56_2e_e_4e 562
147. Carnallite [$Mg(H_2O)_6KCl_3$]: A3B12CDE6_oP276_52_d4e_18e_ce_de_2d8e 532
148. Catapleiite ($Na_2ZrSi_3O_9{\cdot}2H_2O$): A2B3C9D3E_mC144_15_2f_bcdef_9f_3f_ae 314
149. Cd_3As_2: A2B3_tI160_142_deg_3g 1025
150. Cd(OH)Cl ($E0_3$): ABCD_hP8_186_b_b_a_a 1218
151. Ce_2O_2S[¶¶]: A2B2C_hP5_164_d_d_a 1095
152. $CeCu_2$: AB2_oI12_74_e_h 797
153. Cervantite (α-Sb_2O_4): A2B_oP24_33_4a_2a 455
154. Chabazite ($Ca_{1.4}Sr_{0.3}Al_{3.8}Si_{8.3}O_{24}{\cdot}13H_2O$, $S3_4$ (I)): A5B21C24D12_hR62_166_a2c_ehi_fg2h_i 1118
155. Chalcanthite ($CuSO_4{\cdot}5H_2O$, $H4_{10}$): AB10C9D_aP42_2_ae_10i_9i_i 58
156. Chalcocyanite ($CuSO_4$): AB4C_oP24_62_a_2cd_c 711
157. Chiolite ($Na_5Al_3F_{14}$, $K7_5$): A3B14C5_tP44_128_ac_ehi_bg 934
158. Chlorite [$Mg_3(Mg_2Al)(Si_3Al)O_{10}(OH)_8$, $S5_5$]: A3B5C4D2_mC112_15_a3ef_5f_4f_2f 347
159. Chrysotile ($H_4Mg_3Si_2O_9$, $S4_5$): AB6C11D6E4_mC112_12_e_gi2j_i5j_2i2j_2j 178
160. Chrysotile ($Mg_3Si_2O_5(OH)_4$): A3B5C4D2_mC56_9_3a_5a_4a_2a 122
161. Clinocervantite (β-Sb_2O_4): A2B_mC24_15_2f_ce 340
162. Co_2Al_9 ($D8_d$): A9B2_mP22_14_a4e_e 252

[××] $K_2Pt(SCN)_6$ ($H6_3$), Bararite (Trigonal $(NH_4)_2SiF_6$, $J1_6$), and K_2GeF_6 ($J1_{13}$) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

[¶¶] Ce_2O_2S and Brucite [$Mg(OH)_2$] have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

[°°] $MnBi_2Te_4$ and $CaCu_4P_2$ have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

163. $Co_2B_2O_5$: A2B2C5_aP18_2_2i_2i_5i 42
164. $Co_3(SeO_3)_3{\cdot}H_2O$: A3B2C10D3_aP36_2_ah2i_2i_10i_3i 51
165. Co_9S_8 ($D8_9$): A9B8_cF68_225_af_ce 1446
166. Colquiriite ($LiCaAlF_6$): ABC6D_hP18_163_d_b_i_c 1090
167. Columbite ($FeNb_2O_6$, $E5_1$): AB2C6_oP36_60_c_d_3d 608
168. Copper (I) Azide (CuN_3): AB3_tI32_88_d_cf 835
169. Copper (II) Azide [$Cu(N_3)_2$]: AB6_oP28_62_c_6c 716
170. Cr_5O_{12}: A5B12_oP68_60_c2d_6d 604
171. Cr-233 Quasi-One-Dimensional Superconductor ($K_2Cr_3As_3$): A3B3C2_hP16_187_jk_jk_ck 1220
172. $CrCl_3(H_2O)_6$ ($J2_2$): A3BC6_hR20_167_e_b_f ... 1156
173. Crancrinite ($Na_6Ca_2Al_6Si_6O_{24}(CO_3)_2$, $S3_3$ (I)): A3BCD3E15F3_hP52_173_c_b_b_c_5c_c 1168
174. Cronstedtite {$Fe(Fe,Si)[(OH)_2,O]O_3$, $S5_7$}: AB3C2D_hR7_160_a_b_2a_a 1070
175. Cryolite (Na_3AlF_6, $J2_6$): AB6C3_mP20_14_a_3e_de 282
176. $Cs_{11}O_3$: A11B3_mP56_14_11e_3e 201
177. Cs_2Sb: A2B_oP24_62_4c_2c 656
178. Cs_2Se: A2B_oF24_43_b_a 510
179. $Cs_3As_2Cl_9$ ($K7_3$): A2B9C3_hP14_150_d_eg_ad .. 1052
180. Cs_3CoCl_5 ($K3_1$): A5BC3_tI36_140_cl_b_ah 1012
181. $Cs_3Cr_2Cl_9$: A9B2C3_hP28_194_hk_f_bf 1259
182. $Cs_3Tl_2Cl_9$ ($K7_2$): A9B3C2_hR28_167_ef_e_c 1164
183. $Cs_6W_{11}O_{36}$: A6B36C11_mC212_9_6a_36a_11a ... 125
184. Cs_7O: A7B_hP24_187_ai2j2kn_j 1222
185. CsB_4O_6F: A4BCD6_oP48_33_4a_a_a_6a 457
186. $CsFeS_2$ (100 K): ABC2_oI16_71_g_i_eh 789
187. CsO: AB_oI8_71_g_i 791
188. $CsSO_3$ ($K1_2$): AB3C_hP20_190_ac_i_f 1227
189. $Cu_2Pb(SeO_3)_2Br_2$: A2B2C6DE2_oC52_63_g_e_fh_c_f 740
190. $Cu_3[Fe(CN)_6]_2{\cdot}xH_2O$ ($J2_5$, $x \approx 3$): A6B9CD2E6_cF96_225_e_bf_a_c_e 1443
191. Cu(OH)Cl: ABCD_mP16_14_e_e_e_e 295
192. Cubic Cu_2OSeO_3: A2B4C_cP56_198_ab_2a2b_2a 1274
193. Cubic CuPt ($L1_3$ (I), $D4$): AB_cF32_227_c_d ... 1488
194. Danburite ($CaB_2Si_2O_8$, $S6_3$): A2BC8D2_oP52_62_d_c_2c3d_d 651
195. Diamminetriamidodizinc Chloride ($[Zn_2(NH_3)_2(NH_2)_3]Cl$): AB12C5D2_oP40_18_a_6c_b2c_c 388
196. Diaspore (AlOOH, $E0_2$): ABC2_oP16_62_c_c_2c .718
197. Diopside [$CaMg(SiO_3)_2$, $S4_1$]: ABC6D2_mC40_15_e_e_3f_f 378
198. Dodecatungstophosphoric Acid Hexahydrate [$H_3PW_{12}O_{40}{\cdot}6H_2O$]: A27B52CD12_cP184_224_dl_eh3k_a_k 1420
199. Dolomite [$MgCa(CO_3)_2$, $G1_1$]: A2BCD6_hR10_148_c_a_b_f 1039
200. Double Perovskite (Ba_2MnWO_6): A2BC6D_cF40_225_c_a_e_b 1441
201. Enstatite ($MgSiO_3$, $S4_3$): AB3C_oP80_61_2c_6c_2c 623
202. Epididymite ($BeHNaO_8Si_3$, $S4_7$): ABCD8E3_oP112_62_d_2c_d_4c6d_3d 728
203. Eriochalcite ($CuCl_2{\cdot}2H_2O$, $C45$): A2BC4D2_oP18_53_h_a_i_e 544
204. $EuIn_2P_2$: AB2C2_hP10_194_a_f_f 1263
205. Eudidymite ($BeHNaO_8Si_3$): A2B4C2D17E6_mC124_15_f_2f_f_e8f_3f 322
206. Eulytine ($Bi_4(SiO_4)_3$, $S1_5$): A4B12C3_cI76_220_c_e_a 1397
207. $Fe_2(CO)_9$ ($F4_1$): A9B2C9_hP40_176_hi_f_hi 1192
208. Fe_2N (approximate, $L'3_0$)[∂∂]: AB_hP4_194_c_a .. 1272
209. Fe_3PO_7: A3B7C_hR11_160_b_a2b_a 1068
210. Fe_8N ($D2_g$): A8B_tI18_139_deh_a 990
211. FeF_3 ($D0_{12}$): A3B_hR8_167_e_b 1159
212. Ferroelectric $NH_4H_2PO_4$: A6BC4D_oP48_19_6a_a_4a_a 398
213. Ferroelectric $NaNO_2$ ($F5_5$): ABC2_oI8_44_a_a_c 516
214. Fluorapatite [$Ca_5F(PO_4)_3$, $H5_7$]: A5BC12D3_hP42_176_fh_a_2hi_h 1189
215. Ga_2Mg_5 ($D8_g$): A2B5_oI28_72_j_bfj 793
216. $GaMo_4S_8$: AB4C8_cF52_216_a_e_2e 1378
217. Gd_2SiO_5 (RE_2SiO_5 X1): A2B5C_mP32_14_2e_5e_e 221
218. Gwihabaite [NH_4NO_3 (V)]: A4B2C3_tP72_77_8d_ab2c2d_6d 801
219. Gypsum ($CaSO_4{\cdot}2H_2O$, $H4_6$): AB4C6D_mC48_15_e_2f_3f_e 366
220. $H_3PW_{12}O_{40}{\cdot}29H_2O$ ($H4_{21}$): A29B40CD12_cF656_227_ae2fg_e3g_b_g 1467
221. $H_3PW_{12}O_{40}{\cdot}3H_2O$: A3B40CD12_cP112_224_d_e3k_a_k 1429
222. Hambergite [$Be_2BO_3(OH)$, $G7_2$]: AB2CD4_oP64_61_c_2c_c_4c 620
223. Hauyne [$(Na_{0.5}Ca_{0.3}K_{0.2})_8(Al_6Si_6O_{24})(SO_4)_{1.5}$, $S6_9$]: A3B4C4D4E16F4G3_cP76_218_c_e_e_e_ei_e_d .. 1389
224. Hemimorphite ($Zn_4Si_2O_7(OH)_2{\cdot}H_2O$, $S2_2$): A2B5CD2_oI40_44_2c_abcde_d_e 514

[∂∂] $LiZn_2$ (C_k) and Fe_2N (approximate, $L'3_0$) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

225. Hexagonal Delafossite ($CuAlO_2$): ABC2_hP8_194_a_c_f 1268
226. Hexagonal WO_3: A3B_hP12_191_gl_f 1235
227. Hg_2O_2NaI: A2BCD2_hP18_180_f_c_b_i 1198
228. Hg_2TiCu Inverse Heusler: AB2C_cF16_216_b_ad_c 1376
229. $HgCl_2{\cdot}2HgO$: A2B3C2_mP14_14_e_ae_e 208
230. High-Temperature Cryolite (Na_3AlF_6): AB6C3_oI20_71_a_in_cj 787
231. High-Temperature Cubic $KClO_4$ ($H0_5$): ABC4_cF24_216_b_a_e 1380
232. High-Temperature Mo_8O_{23}: A8B23_mP62_13_4g_c11g 195
233. $HoMn_2O_5$: AB2C5_oP32_55_g_fh_eghi 559
234. $HoSb_2$: AB2_oC6_21_a_k 415
235. Huanzalaite ($MgWO_4$, $H0_6$): AB4C_mP12_13_f_2g_e 199
236. In_4Se_3: A4B3_oP28_58_4g_3g 584
237. InS: AB_oP8_58_g_g 589
238. Jacutingaite (Pt_2HgSe_3): AB2C3_hP12_164_d_ae_i 1105
239. $K_2CuCl_4{\cdot}2H_2O$ ($H4_1$): A4BC4D2E2_tP26_136_fg_a_j_d_e 964
240. K_2GeF_6 ($J1_{13}$)[××]: A6BC2_hP9_164_i_a_d 1103
241. $K_2HgCl_4{\cdot}H_2O$ ($E3_4$): A4BCD2_oP32_55_ghi_f_e_gh 553
242. K_2NbF_7 ($K6_2$): A7B2C_mP40_14_7e_2e_e 242
243. $K_2Ni(CN)_4$: A4B2C4D_mP22_14_2e_e_2e_a 230
244. K_2NiF_4: A4B2C_tI14_139_ce_e_a 982
245. $K_2OsO_2Cl_4$ ($J1_5$): A4B2C2D_tI18_139_h_d_e_a 980
246. $K_2Pt(SCN)_6$ ($H6_3$)[××]: A2BC6_hP9_164_d_a_i 1099
247. $K_2Pt(SCN)_6{\cdot}2H_2O$: A6B4C2D6E2FG6_mP54_14_3e_2e_e_3e_e_a_3e 238
248. K_2PtCl_4 ($H1_5$): A4B2C_tP7_123_j_e_a 904
249. $K_2S_2O_5$ ($K0_1$): A2B5C2_mP18_11_2e_e2f_2e 131
250. $K_2S_3O_6$ ($K5_1$): A2B6C3_oP44_62_2c_2c2d_3c 648
251. $K_2Sn(OH)_6$ ($H6_2$): A6B2C6D_hR15_148_f_c_f_a 1041
252. $K_2SnCl_4{\cdot}H_2O$: A4BC2D_oP32_62_2cd_c_d_c 679
253. $K_2SnCl_4{\cdot}H_2O$ ($E3_5$): A4BC2D_oP32_62_2cd_b_2c_a 676
254. $K_2Ti_2O_5$: A2B5C2_mC18_12_i_a2i_i 153
255. $K_3Co(NO_2)_6$ ($J2_4$): AB3C6D12_cF88_202_a_bc_e_h 1289
256. K_3CrO_8: AB3C8_tI24_121_a_bd_2i 886
257. $K_3TlCl_6{\cdot}2H_2O$ ($J3_1$): A6B2C3D_tI168_139_egikl2m_ejn_bh2n_acf 984
258. $K_3W_2Cl_9$ ($K7_1$): A9B3C2_hP28_176_hi_af_f 1195
259. $K_4[Mo(CN)_8]{\cdot}2H_2O$ ($F2_1$): A8B4C4DE8F2_oP108_62_4c2d_2d_2cd_c_4c2d_d 687
260. K(SH) ($B22$): AB_hR2_166_a_b 1140
261. $KAuBr_4{\cdot}2H_2O$ ($H4_{19}$): AB4C2D_mP32_14_e_4e_2e_e 268
262. $KBe_2BO_3F_2$: AB2C2DE3_hR9_155_b_c_c_a_e 1063
263. $KBrO_3$ ($G0_7$)[□]: ABC3_hR5_160_a_a_b 1074
264. $KFeS_2$ ($F5_a$): ABC2_mC16_15_e_e_f 376
265. KH_2PO_4 ($H2_2$): A4BC4D_tI40_122_e_b_e_a 893
266. KHF_2 ($F5_2$): A2BC_tI16_140_h_d_a 1002
267. $KICl_4{\cdot}H_2O$ ($H0_{10}$): A4BCD_mP28_14_4e_e_e_e 232
268. KNO_2 III[‡]: ABC2_mP16_14_e_e_2e 285
269. KSO_3 ($K1_1$): AB3C_hP30_150_ef_3g_c2d 1057
270. Kesterite [$Cu_2(Zn,Fe)SnS_4$]: A2BCD4_tI16_82_ac_b_d_g 805
271. Kotoite ($Mg_3(BO_3)_2$): A2B3C6_oP22_58_g_af_gh 574
272. Kyanite (Al_2SiO_5, $S0_1$): A2B5C_aP32_2_4i_10i_2i 44
273. La_3BWO_9 ($P3$): AB3C9D_hP28_143_2a_2d_6d_bc 1030
274. La_3BWO_9 ($P6_3$): AB3C9D_hP28_173_a_c_3c_b 1175
275. La_3CuSiS_7: AB3C7D_hP24_173_a_c_b2c_b 1172
276. $LaFe_4P_{12}$: A4BC12_cI34_204_c_a_g 1292
277. LaH_{10} High-T_c Superconductor: A10B_cF44_225_cf_b 1439
278. LaOAgS: ABCD_tP8_129_b_c_a_c 950
279. Lepidocrocite (γ-FeO(OH), $E0_4$): AB2C2_oC20_63_c_f_2c 753
280. Li_2CN_2[⊗‡‡]: AB2C2_tI10_139_a_d_e 992
281. Li_2PrO_3: A2B3C_oC12_65_h_bh_a 775
282. $Li_2SO_4{\cdot}H_2O$ ($H4_8$): A2B2C5D_mP20_4_2a_2a_5a_a 77
283. Li_7TaO_6: A8B6C_hR15_148_cf_f_a 1047
284. LiAs[§]: AB_mP16_14_2e_2e 301
285. $LiClO_4{\cdot}3H_2O$ ($H4_{18}$): AB6CD7_hP30_186_b_d_a_b2c 1215
286. $LiCuVO_4$: ABC4D_oI28_74_a_d_hi_e 799
287. $LiGaO_2$: ABC2_oP16_33_a_a_2a 460
288. $LiKSO_4$ ($H1_4$): ABC4D_hP14_173_a_b_bc_b 1180
289. $LiNb_6O_{15}F$: ABC6D15_oP46_51_f_d_2e2i_aef4i2j 529
290. $LiOH{\cdot}H_2O$ ($B36$): A3BC2_mC24_12_ij_h_gi 161
291. $LiZn_2$ (C_k)[∂∂]: AB_hP4_194_a_c 1270
292. Low-Temperature $(NH_3CH_3)Al(SO_4)_2{\cdot}12H_2O$: ABC30DE20F2_oP220_29_a_a_30a_a_20a_2a 422
293. Low-Temperature $GaMo_4S_8$: AB4C8_hR13_160_a_ab_2a2b 1072
294. Low-Temperature Mo_8O_{23}: A8B23_mP124_7_16a_46a 100
295. Lu_2CoGa_3: AB3C2_hP24_194_f_k_bh 1265
296. Lueshite ($NaNbO_3$): ABC3_oP40_57_cd_e_cd2e 571

[‡] KNO_2 III and Manganite (γ-MnO(OH), $E0_6$) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

297. Maghemite (γ-Fe_2O_3, $D5_7$): A2B3_cP60_212_bcd_ace 1354
298. Magnetoplumbite ($PbFe_{12}O_{19}$): A12B19C_hP64_194_ab2fk_efh2k_d 1246
299. Manganese-leonite 110 K [$K_2Mn(SO_4)_2{\cdot}4H_2O$]: A8B2CD12E2_mP100_14_8e_2e_ad_12e_2e 246
300. Manganese-leonite 185 K [$K_2Mn(SO_4)_2{\cdot}4H_2O$]: A8B2CD12E2_mC200_15_8f_2f_ce_2e11f_2f 357
301. Manganese-leonite [$K_2Mn(SO_4)_2{\cdot}4H_2O$, $H4_{23}$]: A8B2CD15E2_mC112_12_2i3j_j_ad_g4i5j_2i 167
302. Manganite (γ-MnO(OH), $E0_6$)[‡]: ABC2_mP16_14_e_e_2e 287
303. Marialite Scapolite [$Na_4Cl(AlSi_3)_3O_{24}$, $S6_4$]: AB4C24D12_tI82_87_a_h_2h2i_hi 825
304. Mayenite ($12CaO{\cdot}7Al_2O_3$, $K7_4$, C12A7): A7B12C19_cI152_220_bc_2d_ace 1400
305. Mercury (II) Azide [$Hg(N_3)_2$]: AB6_oP28_29_a_6a 420
306. Mercury Cyanide [$Hg(CN)_2$, $F1_1$]: A2BC2_tI40_122_e_d_e 890
307. Meta-autunite (I) [$Ca(UO_2)_2(PO_4)_2{\cdot}6H_2O$, $H5_{10}$]: AB4C6DE_tP26_129_c_j_2ci_a_c 945
308. Mg_2Cu (C_b): AB2_oF48_70_g_fg 785
309. $Mg_3Cr_2Al_{18}$: A18B2C3_cF184_227_fg_d_ac 1461
310. Mg_3P_2 ($D5_5$): A3B2_cP10_224_d_b 1427
311. Mg_3Ru_2: A3B2_cP20_213_d_c 1361
312. $Mg(ClO_4)_2{\cdot}6H_2O$ ($H4_{11}$): A2B6CD8_oP34_31_2a_2a2b_a_4a2b 435
313. $Mg(NH_3)_2Cl_2$ ($E1_3$): A2B8CD2_oC26_65_h_r_a_i 777
314. $MgCuAl_2$ ($E1_a$): A2BC_oC16_63_f_c_c 744
315. Mn_3As ($D0_d$): AB3_oC16_63_c_3c 759
316. $MnBi_2Te_4$[°°]: A2BC4_hR7_166_c_a_2c 1114
317. MnCuP: ABC_oP12_62_c_c_c 736
318. $MnF_{2-x}(OH)_x$: A2BC2_oP10_34_c_a_c 464
319. $MnPS_3$: ABC3_mC20_12_g_i_ij 189
320. $Mo_{17}O_{47}$: A17B47_oP128_32_a8c_a23c 442
321. Mo_4P_3: A4B3_oP56_62_8c_6c 673
322. MoP_2: AB2_oC12_36_a_2a 476
323. $MoPO_5$: AB5C_tP14_85_c_cg_b 810
324. Monazite ($LaPO_4$)[†‖]: AB4C_mP24_14_e_4e_e 277
325. Monoclinic Co_4Al_{13}: A13B4_mC102_8_17a11b_8a2b 109
326. Monoclinic Cu_2OSeO_3: A2B4C_mP28_14_abe_4e_e 213
327. Monoclinic $FeTlSe_2$: AB2C_mC16_12_g_2i_i 171
328. Morenosite ($NiSO_4{\cdot}7H_2O$, $H4_{12}$): A14BC11D_oP108_19_14a_a_11a_a 391
329. Murataite [$(Y,Na)_6(Zn,Fe)_5Ti_{12}O_{29}(O,F)_{10}F_4$]: A16B40C12D6E5_cF316_216_eh_e2g2h_h_f_be .. 1367
330. Muscovite ($KH_2Al_3Si_3O_{12}$, $S5_1$): A2BC10D2E4_mC76_15_f_e_5f_f_2f 329
331. NH_4Br ($B25$): AB4C_tP12_129_c_i_a 948
332. NH_4CdCl_3 ($E2_4$): AB3C_oP20_62_c_3c_c 707
333. NH_4ClBrI ($F5_{14}$): ABCD_oP16_62_c_c_c_c 734
334. $NH_4H_2PO_2$ ($F5_7$): A2BC2D_oC24_67_m_a_n_g .. 781
335. $NH_4H_2PO_4$: A8BC4D_tI56_122_2e_b_e_a 896
336. NH_4HF_2 ($F5_8$): A2BC_oP16_53_eh_ab_g 546
337. NH_4HgCl_3 ($E2_5$): A3BC_tP5_123_cg_a_d 902
338. NH_4I_3 ($D0_{16}$): A3B_oP16_62_3c_c 663
339. NH_4NO_3 I ($G0_8$): AB_cP2_221_a_b 1418
340. NH_4NO_3 II ($G0_9$): ABC3_tP10_100_b_a_bc 853
341. NH_4NO_3 III ($G0_{10}$)[°]: ABC3_oP20_62_c_c_cd 723
342. NH_4NO_3 IV ($G0_{11}$): A4B2C3_oP18_59_ef_ab_af . 593
343. $NH_4Pb_2Br_5$ ($K3_4$): A5BC2_tI32_140_bl_a_h 1010
344. NO_2 (Modern, $C26$): AB2_cI36_204_d_g 1296
345. $Na_{0.74}CoO_2$: AB2C2_hP10_194_a_bc_f 1261
346. $Na_2Ca_6Si_4O_{15}$: A6B2C15D4_mP54_7_6a_2a_15a_4a 96
347. Na_2CaSiO_4 ($S6_6$): AB2C4D_cP32_198_a_2a_ab_a 1277
348. Na_2CrO_4 ($H1_8$): AB2C4_oC28_63_c_bc_fg 755
349. $Na_2Mo_2O_7$: A2B2C7_oC88_64_ef_df_3f2g 772
350. Na_2PrO_3: A2B3C_mC48_15_aef_3f_2e 319
351. Na_2SO_3 ($G3_2$): A2B3C_hP12_147_abd_g_d 1037
352. $Na_4Ge_9O_{20}$: A9B4C20_tI132_88_a2f_f_5f 830
353. $NaAlCl_4$: AB4C_oP24_19_a_4a_a 405
354. $NaC_5H_{11}O_8S$: A5B11CD8E_aP26_1_5a_11a_a_8a_a 35
355. $NaCr(SO_4)_2{\cdot}12H_2O$ Alum: AB12CD8E2_cP96_205_a_2d_b_cd_c 1322
356. $NaMn_7O_{12}$: A7BC12_cI40_204_bc_a_g 1294
357. $NaNb_6O_{15}F$: ABC6D15_oC46_38_b_b_2a2d_2ab4d2e 485
358. $NaNbO_3$: ABC3_oP40_17_abcd_2e_abcd4e 383
359. NaP: AB_oP16_19_2a_2a 407
360. NaS_2: AB2_tI48_122_cd_2e 899
361. $NaSb(OH)_6$ ($J1_{11}$): AB6C_tP32_86_d_3g_c 819
362. $NaSbF_4(OH)_2$ ($J1_{12}$): A6BC_hP16_163_i_b_c ... 1087
363. $NaSbF_6$: A6BC_cP32_205_d_b_a 1316
364. Nacrite [$Al_2Si_2O_5(OH)_4$, $S5_4$]: A2B4C9D2_mC68_9_2a_4a_9a_2a 119
365. Nahcolite ($NaHCO_3$, $G0_{12}$): ABCD3_mP24_14_e_e_e_3e 292
366. Natrolite ($Na_2Al_2Si_3O_{10}{\cdot}2H_2O$, $S6_{10}$): A2B4C2D12E3_oF184_43_b_2b_b_6b_ab 499
367. $Nb_2Pd_3Se_8$: A2B3C8_oP26_55_h_ag_2g2h 551
368. $Nb_2Zr_6O_{17}$: A2B17C6_oI100_46_ab_b8c_3c 523
369. Nb_3O_7F: A3B8_oC22_65_ag_bd2gh 779
370. $NbAs_2$: A2B_mC12_5_2c_c 85
371. $NbTe_2$: AB2_mC18_12_ai_3i 173
372. $Nd_2Fe_{14}B$: AB14C2_tP68_136_f_ce2j2k_fg 966
373. $Nd_4Re_2O_{11}$: A4B11C2_tP68_86_2g_ab5g_g 815
374. $Nd(BrO_3)_3{\cdot}9H_2O$ ($G2_2$): A3B9CD9_hP44_186_c_3c_b_cd 1207
375. Nevskite (BiSe): AB_hP12_164_c2d_c2d 1109

376. Ni_3Si_2: A3B2_oC80_36_4a4b_2a3b 470
377. $Ni(H_2O)_6SnCl_6$ ($I6_1$): A6B6CD_hR14_148_f_f_b_a 1044
378. $Ni(NO_3)_2(H_2O)_6$: A12B2CD12_aP54_2_12i_2i_i_12i 38
379. Norbergite [$Mg(F,OH)_2 \cdot Mg_2SiO_4$, $S0_7$]: A2B3C4D_oP40_62_d_cd_2cd_c 632
380. O(OH)Y: ABC_mP6_11_e_e_e 139
381. Original β-WO_3 (*obsolete*): A3B_oP32_62_ab4c_2c 665
382. Orpiment (As_2S_3, $D5_f$): A2B3_mP20_14_2e_3e .. 210
383. Orthorhombic Co_4Al_{13}: A13B4_oP102_31_17a11b_8a2b 430
384. Orthorhombic CrO_3: AB3_oC16_40_b_a2b 491
385. Orthorhombic $Sr_4Ru_3O_{10}$: A10B3C4_oP68_55_2e2fgh2i_adef_2e2f 548
386. Os_4Al_{13}: A13B4_mC34_12_b6i_2i 144
387. P_4Se_3: A4B3_oP112_62_8c4d_4c4d 668
388. $PNCl_2$ ($E1_4$): A2BC_tP32_86_2g_g_g 812
389. Paralstonite ($BaCa(CO_3)_2$): AB2CD6_hP30_150_e_c2d_f_3g 1054
390. Pararealgar (AsS)[¶]: AB_mP32_14_4e_4e 306
391. Paratellurite (α-TeO_2): A2B_tP12_92_b_a 846
392. Parawollastonite ($CaSiO_3$, $S3_3$ (II)): AB3C_mP60_14_3e_9e_3e 260
393. Parkerite ($Ni_3Bi_2S_2$): AB2C_oP8_51_e_be_f 527
394. $Pb(NO_3)_2$ ($G2_1$): A2B6C_cP36_205_c_d_a 1306
395. Pd_5Pu_3: A5B3_oC32_63_cfg_ce 749
396. $Pd(NH_3)_4Cl_2 \cdot H_2O$ ($H4_9$): A2BC4D_tP16_127_h_d_i_a 929
397. Phase II $Cd_2Re_2O_7$: A2B7C2_tI44_119_i_bdefgh_i 874
398. Phase III $Cd_2Re_2O_7$: A2B7C2_tI44_98_f_bcde_f 848
399. Phosgenite [$Pb_2Cl_2(CO_3)$]: AB2C3D2_tP32_127_g_eh_gk_k 931
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405. Predicted Phase IV $Cd_2Re_2O_7$: A2B7C2_oF88_22_k_bdefghij_k 417
406. Proposed 300 GPa HfH_{10}: A10B_hP22_194_bhj_c 1243
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409. Pt_2Sn_3 ($D5_b$): A2B3_hP10_194_f_bf 1249
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413. $Rb_2CaCu_6(PO_4)_4O_2$: AB6C18D4E2_mC62_5_a_2b2c_9c_2c_c 89
414. $Rb_2Mo_2O_7$: A2B7C2_oC88_40_abc_2b6c_a3b 488
415. $RbNO_3$ (IV): AB3C_hP45_144_3a_9a_3a 1033
416. Re_3B: AB3_oC16_63_c_cf 761
417. ReB_3: A3B_hP8_194_af_c 1257
418. Realgar (AsS, B_l)[¶]: AB_mP32_14_4e_4e 309
419. Retgersite (α-$NiSO_4 \cdot 6H_2O$, $H4_5$): A12BC10D_tP96_92_6b_a_5b_a 841
420. $Rh_{20}Si_{13}$: A10B7_hP34_176_c3h_b2h 1182
421. $RhCl_2(NH_3)_5Cl$ ($J1_8$): A3B15C5D_oP96_62_cd_3c6d_3cd_c 658
422. Rhombohedral $CuTi_2S_4$: AB4C2_hR28_166_2c_2c2h_abh 1126
423. Rhombohedral Delafossite ($CuFeO_2$): ABC2_hR4_166_a_b_c 1136
424. Rinneite ($K_3NaFeCl_6$): A6BC3D_hR22_167_f_b_e_a 1161
425. Rosiaite ($PbSb_2O_6$)[§§]: A6BC2_hP9_162_k_a_d .. 1085
426. $Ru_{11}B_8$: A8B11_oP38_55_g3h_a3g2h 556
427. RuB_2: A2B_oP6_59_f_a 591
428. Rynersonite (Orthorhombic $CaTa_2O_6$): AB6C2_oP36_62_c_2c2d_d 713
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433. $SbCl_5 \cdot POCl_3$: A8BCD_oP44_62_4c2d_c_c_c 692
434. SbI_3S_{24}: A3B24C_hR28_160_b_2b3c_a 1065
435. Scheelite ($CaWO_4$, $H0_4$): AB4C_tI24_88_b_f_a .. 837

[¶]Pararealgar (AsS) and Realgar (AsS, B_l) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[††]Pseudobrookite (Fe_2TiO_5, $E4_1$) and Ta_2NiS_5 have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

436. Senarmontite (Sb_2O_3, $D6_1$): A3B2_cF80_227_f_e 1481
437. Shandite ($Ni_3Pb_2S_2$): A3B2C2_hR7_166_d_ab_c 1116
438. Shcherbinaite (V_2O_5) (*revised*): A5B2_oP14_59_a2f_f 595
439. Si_{24} Clathrate: A_oC24_63_3f 767
440. Si_2N_2O: A2BC2_oC20_36_b_a_b 466
441. SiAs: AB_mC24_12_3i_3i 193
442. Sillimanite (Al_2SiO_5, $S0_3$): A2B5C_oP32_62_bc_3cd_c 645
443. $Sm_{11}Cd_{45}$: A45B11_cF448_216_bd4efg5h_ac2eh . 1371
444. SnI_4 ($D1_1$): A4B_cP40_205_cd_c 1313
445. Sodalite [$Na_4(AlSiO_4)_3Cl$, $S6_2$]: A3BC4D12E3_cP46_218_d_a_e_i_c 1394
446. Sr_2MnTeO_6: AB6C2D_mP20_14_a_3e_e_d 280
447. Sr_2NiTeO_6: AB6C2D_mC40_12_ad_gh4i_j_bc 182
448. Sr_2NiWO_6: AB6C2D_tI20_87_a_eh_d_b 828
449. $Sr_3Ti_2O_7$: A7B3C2_tI24_139_aeg_be_e 988
450. $Sr_4Ti_3O_{10}$: A10B4C3_tI34_139_c2eg_2e_ae 970
451. $Sr(OH)_2(H_2O)_8$: A18B10C_tP116_130_2c4g_2c2g_a 952
452. $SrCl_2{\cdot}(H_2O)_6$: A2B12C6D_hP21_150_d_2g_ef_a 1049
453. $SrCu_2(BO_3)_2$: A2B2C6D_tI44_121_i_i_ij_c 882
454. $SrUO_4$: A4BC_oP24_57_cde_d_a 569
455. Staurolite ($Al_5Fe_2O_{10}(OH)_2Si_2$): A5B2C10D2E2_mC84_12_acghj_bdi_5j_2i_j 163
456. Steklite [$KAl(SO_4)_2$, $H3_2$]: ABC8D2_hP12_150_b_a_dg_d 1060
457. Sulphohalite [$Na_6ClF(SO_4)_2$, $H5_8$]: ABC6D8E2_cF72_225_b_a_e_f_c 1453
458. Swedenborgite ($NaBe_4SbO_7$, $E9_2$): A4BC7D_hP26_186_ac_b_a2c_b 1210
459. Ta_2NiS_5[††]: AB5C2_oC32_63_c_c2f_f 763
460. Ta_2NiSe_5: AB5C2_mC32_15_e_e2f_f 369
461. Ta_2PdSe_6: AB6C2_mC18_12_a_3i_i 184
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463. Ta_3Ti_5 (BCC SQS-16): A3B5_oC32_38_abce_abcdf 483
464. Ta_5Ti_{11} (BCC SQS-16): A5B11_mP16_6_2abc_2a3b3c 94
465. TaTi (BCC SQS-16): AB_aP16_2_4i_4i 71
466. $TaTi_3$ (BCC SQS-16): AB3_mC32_8_4a_12a 113
467. $TaTi_3$ (BCC SQS-16): AB3_mC32_8_4a_4a4b 115
468. $TaTi_7$ (BCC SQS-16): AB7_hR16_166_c_c2h ... 1133
469. Tellurite (β-TeO_2, $C52$): A2B_oP24_61_2c_c 614
470. Tennantite ($Cu_{12}As_4S_{13}$): A4B24C13_cI82_217_c_deg_ag 1384
471. Tetragonal $TlFeS_2$: AB2C_tI8_119_c_e_a 877
472. Th_7S_{12} ($D8_k$): A3B2_hP20_176_2h_ah 1185
473. ThC_2 (C_g): A2B_mC12_15_f_e 338
474. $ThCr_2Si_2$[∂‡‡]: A2B2C_tI10_139_d_e_a 974
475. $ThFe_2SiC$: AB2CD_oC20_63_b_f_c_c 757
476. Thenardite [Na_2SO_4 (V), $H1_7$]: A2B4C_oF56_70_g_h_a 783
477. Ti_5Ga_4: A4B5_hP18_193_bg_dg 1241
478. $TiBe_{12}$ (approximate, $D2_a$): A12B_hP13_191_cdei_a 1233
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483. Tolbachite ($CuCl_2$): A2B_mC6_12_i_a 157
484. Topaz ($Al_2SiO_4F_2$, $S0_5$): A2B2C4D_oP36_62_d_d_2cd_c 629
485. Tremolite ($Ca_2Mg_5Si_8O_{22}(OH)_2$, $S4_2$): A2B2C5D24E8_mC82_12_h_i_agh_2i5j_2j 148
486. Tutton salt [$Cu(NH_4)_2(SO_4)_2{\cdot}6H_2O$, $H4_4$]: AB20C2D14E2_mP78_14_a_10e_e_7e_e 255
487. U_6Mn ($D2_c$): AB6_tI28_140_a_hk 1014
488. V_3AsC: ABC3_oC20_63_c_b_cf 765
489. V_4SiSb_2: A2BC4_tI28_140_h_a_k 1000
490. $VOSO_4$: A5BC_oP28_62_3cd_c_c 682
491. VSe_2O_6: A6B2C_tP72_103_abc5d_2d_abc 855
492. Vesuvianite ($Ca_{10}Al_4(Mg,Fe)_2Si_9O_{34}(OH)_4$, $S2_3$): A4B10C2D34E4F9_tP252_126_k_ce2k_f_h8k_k_d2k 916
493. $W_2O_3(PO_4)_2$: A11B2C2_mP60_4_22a_4a_4a 73
494. Wülfingite (ϵ-$Zn(OH)_2$, $C31$): A2B2C_oP20_19_2a_2a_a 396
495. Wollastonite ($CaSiO_3$): AB3C_aP30_2_3i_9i_3i 65
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497. $Zn_{22}Zr$: A22B_cF184_227_cdfg_a 1464
498. $Zn_2Mo_3O_8$: A3B8C2_hP26_186_c_ab2c_2b 1204
499. $Zn(BrO_3)_2{\cdot}6H_2O$ ($J1_{10}$)[⊗⊗]: A2B6C6D_cP60_205_c_d_d_a 1299
500. $Zn(NH_3)_2Cl_2$ ($E1_2$): A2B6C2D_oI44_74_h_ij_i_e . 795
501. $Zr_{21}Re_{25}$: A25B21_hR92_167_b2e3f_e3f 1142
502. Zr_2Al_3[**]: A3B2_oF40_43_ab_b 512
503. Zr_3Al_2: A2B3_tP20_136_j_dfg 960
504. $ZrFe_4Si_2$: A4B2C_tP14_136_i_g_b 962
505. ZrNiAl: ABC_hP9_189_g_ad_f 1225
506. ZrP_2O_7 High-Temperature ($K6_1$): A7B2C_cP40_205_bd_c_a 1319
507. $ZrSe_3$: A3B_mP8_11_3e_e 133
508. $ZrTe_5$: A5B_oC24_63_c2f_c 751

[∂] $TlCo_2S_2$ and $ThCr_2Si_2$ have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

509. Zunyite [$Al_{13}(OH,F)_{18}Si_5O_{20}Cl$, $S0_8$]: A13BC18D20E5_cF228_216_dh_b_fh_2eh_ce 1363
510. "Martensite Type" FeC_x ($x \leq 0.06$) ($L'2_0$): AB_tI4_139_b_a 998

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    14.1. $Bi_3Ru_3O_{11}$: A3B11C3_cP68_201_be_efh_g ... 1285
15. **cP76**
    15.1. Hauyne [$(Na_{0.5}Ca_{0.3}K_{0.2})_8(Al_6Si_6O_{24})(SO_4)_{1.5}$, $S6_9$]: A3B4C4D4E16F4G3_cP76_218_c_e_e_e_ei_e_d 1389
16. **cP84**
    16.1. CaB_2O_4 (IV): A2BC4_cP84_205_d_ac_2d ... 1309
17. **cP96**
    17.1. $NaCr(SO_4)_2\cdot12H_2O$ Alum: AB12CD8E2_cP96_205_a_2d_b_cd_c ... 1322
18. **cP112**
    18.1. $H_3PW_{12}O_{40}\cdot3H_2O$: A3B40CD12_cP112_224_d_e3k_a_k ... 1429
19. **cP116**
    19.1. 12-phosphotungstic acid [$H_3PW_{12}O_{40}\cdot5H_2O$ ($H4_{16}$)]: A5B40CD12_cP116_224_cd_e3k_a_k ... 1434
20. **cP184**
    20.1. Dodecatungstophosphoric Acid Hexahydrate [$H_3PW_{12}O_{40}\cdot6H_2O$]: A27B52CD12_cP184_224_dl_eh3k_a_k ... 1420
21. **cP192**
    21.1. γ-Alum [$AlNa(SO_4)_2\cdot12H_2O$, $H4_{15}$]: AB24CD20E2_cP192_205_a_4d_b_c3d_c ... 1327
22. **cP224**
    22.1. α-Alum [$KAl(SO_4)_2\cdot12H_2O$, $H4_{13}$]: AB24CD28E2_cP224_205_a_4d_b_2c4d_c ... 1335

[⊗⊗] $Zn(BrO_3)_2\cdot6H_2O$ ($J1_{10}$) and $H6_4$ [$Ni(NO_3)_2(NH_3)_6$] (*obsolete*) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

23. **cP252**
    23.1. β-Alum [$Al(NH_3CH_3)_2(SO_4)_2{\cdot}12H_2O$, $H4_{14}$]: AB2C36D2E20F2_cP252_205_a_c_6d_c_c3d_c 1344

hP

1. **hP4**
    1.1. $D0_{13}$ ($AlCl_3$) (*obsolete*): AB3_hP4_164_b_ad 1107
    1.2. $LiZn_2$ (C_k)[∂∂]: AB_hP4_194_a_c 1270
    1.3. Fe_2N (approximate, $L'3_0$)[∂∂]: AB_hP4_194_c_a 1272
2. **hP5**
    2.1. Ce_2O_2S[¶¶]: A2B2C_hP5_164_d_d_a 1095
    2.2. Brucite [$Mg(OH)_2$][¶¶]: A2BC2_hP5_164_d_a_d 1097
3. **hP6**
    3.1. $C27$ (CdI_2) (*questionable*): AB2_hP6_186_b_ab 1213
4. **hP8**
    4.1. Cd(OH)Cl ($E0_3$): ABCD_hP8_186_b_b_a_a 1218
    4.2. ReB_3: A3B_hP8_194_af_c 1257
    4.3. Hexagonal Delafossite ($CuAlO_2$): ABC2_hP8_194_a_c_f 1268
5. **hP9**
    5.1. $I1_3$ ($SrCl_2{\cdot}(H_2O)_6$) (*obsolete*)[§§]: A2B6C_hP9_162_d_k_a 1083
    5.2. Rosiaite ($PbSb_2O_6$)[§§]: A6BC2_hP9_162_k_a_d 1085
    5.3. $K_2Pt(SCN)_6$ ($H6_3$)[××]: A2BC6_hP9_164_d_a_i 1099
    5.4. Bararite (Trigonal $(NH_4)_2SiF_6$, $J1_6$)[××]: A6B2C_hP9_164_i_d_a 1101
    5.5. K_2GeF_6 ($J1_{13}$)[××]: A6BC2_hP9_164_i_a_d 1103
    5.6. ZrNiAl: ABC_hP9_189_g_ad_f 1225
6. **hP10**
    6.1. α-$LiIO_3$: ABC3_hP10_173_b_a_c 1178
    6.2. $E2_3$ ($LiIO_3$) (*obsolete*): ABC3_hP10_182_c_b_g 1202
    6.3. Pt_2Sn_3 ($D5_b$): A2B3_hP10_194_f_bf 1249
    6.4. $Na_{0.74}CoO_2$: AB2C2_hP10_194_a_bc_f 1261
    6.5. $EuIn_2P_2$: AB2C2_hP10_194_a_f_f 1263
7. **hP12**
    7.1. Na_2SO_3 ($G3_2$): A2B3C_hP12_147_abd_g_d 1037
    7.2. Steklite [$KAl(SO_4)_2$, $H3_2$]: ABC8D2_hP12_150_b_a_dg_d 1060
    7.3. Jacutingaite (Pt_2HgSe_3): AB2C3_hP12_164_d_ae_i 1105
    7.4. Nevskite (BiSe): AB_hP12_164_c2d_c2d 1109
    7.5. Hexagonal WO_3: A3B_hP12_191_gl_f 1235
8. **hP13**
    8.1. $TiBe_{12}$ (approximate, $D2_a$): A12B_hP13_191_cdei_a 1233
9. **hP14**
    9.1. $Cs_3As_2Cl_9$ ($K7_3$): A2B9C3_hP14_150_d_eg_ad 1052
    9.2. $LiKSO_4$ ($H1_4$): ABC4D_hP14_173_a_b_bc_b 1180
    9.3. β-Si_3N_4: A4B3_hP14_176_ch_h 1187
10. **hP16**
    10.1. $NaSbF_4(OH)_2$ ($J1_{12}$): A6BC_hP16_163_i_b_c 1087
    10.2. Cr-233 Quasi-One-Dimensional Superconductor ($K_2Cr_3As_3$): A3B3C2_hP16_187_jk_jk_ck 1220
11. **hP18**
    11.1. Colquiriite ($LiCaAlF_6$): ABC6D_hP18_163_d_b_i_c 1090
    11.2. Hg_2O_2NaI: A2BCD2_hP18_180_f_c_b_i 1198
    11.3. $BaAl_2O_4$ ($H2_8$): A2BC6_hP18_182_f_b_gh 1200
    11.4. Ti_5Ga_4: A4B5_hP18_193_bg_dg 1241
12. **hP19**
    12.1. Predicted Li_2MgH_{16} 300 GPa: A16B2C_hP19_164_2d2i_d_b 1093
13. **hP20**
    13.1. Th_7S_{12} ($D8_k$): A3B2_hP20_176_2h_ah 1185
    13.2. $CsSO_3$ ($K1_2$): AB3C_hP20_190_ac_i_f 1227
14. **hP21**
    14.1. $SrCl_2{\cdot}(H_2O)_6$: A2B12C6D_hP21_150_d_2g_ef_a 1049
15. **hP22**
    15.1. Proposed 300 GPa HfH_{10}: A10B_hP22_194_bhj_c 1243
16. **hP24**
    16.1. La_3CuSiS_7: AB3C7D_hP24_173_a_c_b2c_b 1172
    16.2. Cs_7O: A7B_hP24_187_ai2j2kn_j 1222
    16.3. $D0_6$ (Tysonite, LaF_3) (*obsolete*): A3B_hP24_193_ack_g 1238
    16.4. Lu_2CoGa_3: AB3C2_hP24_194_f_k_bh 1265
17. **hP26**
    17.1. $Zn_2Mo_3O_8$: A3B8C2_hP26_186_c_ab2c_2b 1204
    17.2. Swedenborgite ($NaBe_4SbO_7$, $E9_2$): A4BC7D_hP26_186_ac_b_a2c_b 1210
18. **hP28**

[∂∂] $LiZn_2$ (C_k) and Fe_2N (approximate, $L'3_0$) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

[¶¶] Ce_2O_2S and Brucite [$Mg(OH)_2$] have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

[§§] $I1_3$ ($SrCl_2{\cdot}(H_2O)_6$) (*obsolete*) and Rosiaite ($PbSb_2O_6$) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

[××] $K_2Pt(SCN)_6$ ($H6_3$), Bararite (Trigonal $(NH_4)_2SiF_6$, $J1_6$), and K_2GeF_6 ($J1_{13}$) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

18.1. La_3BWO_9 ($P3$): AB3C9D_hP28_143_2a_2d_6d_bc 1030
18.2. La_3BWO_9 ($P6_3$): AB3C9D_hP28_173_a_c_3c_b 1175
18.3. $K_3W_2Cl_9$ ($K7_1$): A9B3C2_hP28_176_hi_af_f 1195
18.4. $Cs_3Cr_2Cl_9$: A9B2C3_hP28_194_hk_f_bf ... 1259

19. **hP30**
19.1. Paralstonite ($BaCa(CO_3)_2$): AB2CD6_hP30_150_e_c2d_f_3g 1054
19.2. KSO_3 ($K1_1$): AB3C_hP30_150_ef_3g_c2d .. 1057
19.3. $LiClO_4{\cdot}3H_2O$ ($H4_{18}$): AB6CD7_hP30_186_b_d_a_b2c 1215

20. **hP34**
20.1. $Rh_{20}Si_{13}$: A10B7_hP34_176_c3h_b2h 1182

21. **hP36**
21.1. Bastnäsite [$CeF(CO_3)$]: ABCD3_hP36_190_h_g_af_hi 1230
21.2. $S3_4$ (II) (Catapleiite, $Na_2Zr(SiO_3)_3{\cdot}\ H_2O$) (*obsolete*): A3B2C9D3E_hP36_194_g_f_hk_h_a 1254

22. **hP40**
22.1. $Fe_2(CO)_9$ ($F4_1$): A9B2C9_hP40_176_hi_f_hi 1192

23. **hP42**
23.1. Fluorapatite [$Ca_5F(PO_4)_3$, $H5_7$]: A5BC12D3_hP42_176_fh_a_2hi_h 1189

24. **hP44**
24.1. $Nd(BrO_3)_3{\cdot}9H_2O$ ($G2_2$): A3B9CD9_hP44_186_c_3c_b_cd 1207

25. **hP45**
25.1. $RbNO_3$ (IV): AB3C_hP45_144_3a_9a_3a .. 1033

26. **hP52**
26.1. Crancrinite ($Na_6Ca_2Al_6Si_6O_{24}(CO_3)_2$, $S3_3$ (I)): A3BCD3E15F3_hP52_173_c_b_b_c_5c_c ... 1168

27. **hP60**
27.1. β-Alumina (Al_2O_3, $D5_6$): A2B3_hP60_194_3fk_cdef2k 1251

28. **hP64**
28.1. Magnetoplumbite ($PbFe_{12}O_{19}$): A12B19C_hP64_194_ab2fk_efh2k_d 1246

hR .....

1. **hR2**
1.1. K(SH) ($B22$): AB_hR2_166_a_b 1140

2. **hR4**
2.1. Rhombohedral Delafossite ($CuFeO_2$): ABC2_hR4_166_a_b_c 1136

3. **hR5**
3.1. $KBrO_3$ ($G0_7$)[□]: ABC3_hR5_160_a_a_b 1074
3.2. γ-Potassium Nitrate (KNO_3)[□]: ABC3_hR5_160_a_a_b 1076

4. **hR6**
4.1. $CaSi_2$ ($C12$): AB2_hR6_166_c_2c 1124
4.2. $CaUO_4$: AB4C_hR6_166_b_2c_a 1131

5. **hR7**
5.1. Cronstedtite {$Fe(Fe,Si)[(OH)_2,O]O_3$, $S5_7$}: AB3C2D_hR7_160_a_b_2a_a 1070
5.2. $MnBi_2Te_4$[○○]: A2BC4_hR7_166_c_a_2c 1114
5.3. Shandite ($Ni_3Pb_2S_2$): A3B2C2_hR7_166_d_ab_c 1116
5.4. $CaCu_4P_2$[○○]: AB4C2_hR7_166_a_2c_c 1129

6. **hR8**
6.1. β-Potassium Nitrate (KNO_3): ABC6_hR8_166_a_b_h 1138
6.2. FeF_3 ($D0_{12}$): A3B_hR8_167_e_b 1159

7. **hR9**
7.1. $KBe_2BO_3F_2$: AB2C2DE3_hR9_155_b_c_c_a_e 1063

8. **hR10**
8.1. Dolomite [$MgCa(CO_3)_2$, $G1_1$]: A2BCD6_hR10_148_c_a_b_f 1039

9. **hR11**
9.1. Fe_3PO_7: A3B7C_hR11_160_b_a2b_a 1068

10. **hR13**
10.1. Low-Temperature $GaMo_4S_8$: AB4C8_hR13_160_a_ab_2a2b 1072

11. **hR14**
11.1. $Ni(H_2O)_6SnCl_6$ ($I6_1$): A6B6CD_hR14_148_f_f_b_a 1044

12. **hR15**
12.1. $K_2Sn(OH)_6$ ($H6_2$): A6B2C6D_hR15_148_f_c_f_a 1041
12.2. Li_7TaO_6: A8B6C_hR15_148_cf_f_a 1047
12.3. $B_{13}C_2$ "B_4C" ($D1_g$): A13B2_hR15_166_b2h_c 1111

13. **hR16**
13.1. $TaTi_7$ (BCC SQS-16): AB7_hR16_166_c_c2h 1133

14. **hR20**
14.1. $CrCl_3(H_2O)_6$ ($J2_2$): A3BC6_hR20_167_e_b_f 1156

15. **hR22**
15.1. Rinneite ($K_3NaFeCl_6$): A6BC3D_hR22_167_f_b_e_a 1161

16. **hR28**
16.1. SbI_3S_{24}: A3B24C_hR28_160_b_2b3c_a 1065

[□] $KBrO_3$ ($G0_7$) and γ-Potassium Nitrate (KNO_3) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[○○] $MnBi_2Te_4$ and $CaCu_4P_2$ have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

16.2. Rhombohedral $CuTi_2S_4$: AB4C2_hR28_166_2c_2c2h_abh 1126
16.3. $Cs_3Tl_2Cl_9$ ($K7_2$): A9B3C2_hR28_167_ef_e_c 1164
17. **hR42**
17.1. α-BaB_2O_4 (Low-Temperature): A2BC4_hR42_161_2b_b_4b 1078
17.2. β-BaB_2O_4 (High-Temperature): A2BC4_hR42_167_f_ac_2f 1151
18. **hR62**
18.1. Chabazite ($Ca_{1.4}Sr_{0.3}Al_{3.8}Si_{8.3}O_{24}{\cdot}13H_2O$, $S3_4$ (I)): A5B21C24D12_hR62_166_a2c_ehi_fg2h_i .. 1118
19. **hR92**
19.1. $Zr_{21}Re_{25}$: A25B21_hR92_167_b2e3f_e3f ... 1142
mC .....
1. **mC4**
1.1. β-Ga (*obsolete*): A_mC4_15_e 381
2. **mC6**
2.1. Tolbachite ($CuCl_2$): A2B_mC6_12_i_a 157
3. **mC8**
3.1. $F5_{11}$ (KNO_2) (*obsolete*): ABC2_mC8_8_a_a_b 117
4. **mC10**
4.1. $C2$ $(Ba,Ca)CO_3$: ABC3_mC10_5_b_a_ac 92
5. **mC12**
5.1. $NbAs_2$: A2B_mC12_5_2c_c 85
5.2. CaC_2-III: A2B_mC12_12_2i_i 155
5.3. ThC_2 (C_g): A2B_mC12_15_f_e 338
6. **mC14**
6.1. δ-Ni_3Sn_4 ($D7_a$): A3B4_mC14_12_ai_2i 159
7. **mC16**
7.1. $D0_{15}$ ($AlCl_3$) (*obsolete*): AB3_mC16_5_c_3c 87
7.2. Monoclinic $FeTlSe_2$: AB2C_mC16_12_g_2i_i 171
7.3. $KFeS_2$ ($F5_a$): ABC2_mC16_15_e_e_f 376
8. **mC18**
8.1. $K_2Ti_2O_5$: A2B5C2_mC18_12_i_a2i_i 153
8.2. $NbTe_2$: AB2_mC18_12_ai_3i 173
8.3. Ta_2PdSe_6: AB6C2_mC18_12_a_3i_i 184
9. **mC20**
9.1. β-Ga_2O_3: A2B3_mC20_12_2i_3i 151
9.2. $MnPS_3$: ABC3_mC20_12_g_i_ij 189
9.3. Ag_2PbO_2: A2B2C_mC20_15_ad_f_e 312
10. **mC24**
10.1. $LiOH{\cdot}H_2O$ ($B36$): A3BC2_mC24_12_ij_h_gi 161
10.2. $AlNbO_4$: ABC4_mC24_12_i_i_4i 191
10.3. SiAs: AB_mC24_12_3i_3i 193
10.4. Clinocervantite (β-Sb_2O_4): A2B_mC24_15_2f_ce 340
11. **mC28**
11.1. ζ-Nb_2O_5 (B-Nb_2O_5): A2B5_mC28_15_f_e2f 327
12. **mC32**
12.1. $TaTi_3$ (BCC SQS-16): AB3_mC32_8_4a_12a 113
12.2. $TaTi_3$ (BCC SQS-16): AB3_mC32_8_4a_4a4b 115
12.3. Ta_2NiSe_5: AB5C2_mC32_15_e_e2f_f 369
12.4. Titanite ($CaTiSiO_5$, $S0_6$): AB5CD_mC32_15_e_e2f_e_b 374
13. **mC34**
13.1. Os_4Al_{13}: A13B4_mC34_12_b6i_2i 144
14. **mC36**
14.1. $Rb_2C_2O_4{\cdot}H_2O$: A2BC4D2_mC36_15_f_e_2f_f 332
15. **mC40**
15.1. Sr_2NiTeO_6: AB6C2D_mC40_12_ad_gh4i_j_bc 182
15.2. Diopside [$CaMg(SiO_3)_2$, $S4_1$]: ABC6D2_mC40_15_e_e_3f_f 378
16. **mC42**
16.1. Bischofite ($MgCl_2{\cdot}6H_2O$, $J1_7$): A2B12CD6_mC42_12_i_2i2j_a_ij 146
17. **mC44**
17.1. α-$Zn_2V_2O_7$: A7B2C2_mC44_15_e3f_f_f 354
18. **mC48**
18.1. $D2_2$ ($MgZn_5$?) (*problematic*): AB5_mC48_12_2i_ac5i2j 175
18.2. Na_2PrO_3: A2B3C_mC48_15_aef_3f_2e 319
18.3. Gypsum ($CaSO_4{\cdot}2H_2O$, $H4_6$): AB4C6D_mC48_15_e_2f_3f_e 366
19. **mC52**
19.1. Sanidine ($KAlSi_3O_8$, $S6_7$): AB8C4_mC52_12_i_gi3j_2j 186
20. **mC56**
20.1. Chrysotile ($Mg_3Si_2O_5(OH)_4$): A3B5C4D2_mC56_9_3a_5a_4a_2a 122
20.2. $BaNi(CN)_4{\cdot}4H_2O$ ($H4_{22}$): AB4C4D4E_mC56_15_e_2f_2f_2f_a 363
21. **mC62**
21.1. $Rb_2CaCu_6(PO_4)_4O_2$: AB6C18D4E2_mC62_5_a_2b2c_9c_2c_c 89
22. **mC64**
22.1. Y_2SiO_5 (RE_2SiO_5 X2): A5BC2_mC64_15_5f_f_2f 351
23. **mC68**
23.1. Nacrite [$Al_2Si_2O_5(OH)_4$, $S5_4$]: A2B4C9D2_mC68_9_2a_4a_9a_2a 119
24. **mC72**
24.1. Pyrophyllite [$AlSi_2O_5(OH)$, $S5_6$]: AB5CD2_mC72_15_f_5f_f_2f 371
25. **mC76**

25.1. Muscovite ($KH_2Al_3Si_3O_{12}$, $S5_1$): A2BC10D2E4_mC76_15_f_e_5f_f_2f 329
25.2. Alluaudite [$NaMnFe_2(PO_4)_3$]: A2BCD12E3_mC76_15_f_e_b_6f_ef 334

26. **mC82**
26.1. Tremolite ($Ca_2Mg_5Si_8O_{22}(OH)_2$, $S4_2$): A2B2C5D24E8_mC82_12_h_i_agh_2i5j_2j ... 148

27. **mC84**
27.1. Staurolite ($Al_5Fe_2O_{10}(OH)_2Si_2$): A5B2C10D2E2_mC84_12_acghj_bdi_5j_2i_j . 163

28. **mC90**
28.1. Bassanite [$CaSO_4(H_2O)_{0.5}$, $H4_7$]: A2B2C9D2_mC90_5_ab2c_3c_b13c_3c 81

29. **mC102**
29.1. Monoclinic Co_4Al_{13}: A13B4_mC102_8_17a11b_8a2b 109
29.2. $Al_{13}Fe_4$: A13B4_mC102_12_dg8i5j_4ij 141

30. **mC112**
30.1. Manganese-leonite [$K_2Mn(SO_4)_2{\cdot}4H_2O$, $H4_{23}$]: A8B2CD15E2_mC112_12_2i3j_j_ad_g4i5j_2i . 167
30.2. Chrysotile ($H_4Mg_3Si_2O_9$, $S4_5$): AB6C11D6E4_mC112_12_e_gi2j_i5j_2i2j_2j . 178
30.3. Chlorite [$Mg_3(Mg_2Al)(Si_3Al)O_{10}(OH)_8$, $S5_5$]: A3B5C4D2_mC112_15_a3ef_5f_4f_2f 347

31. **mC124**
31.1. Eudidymite ($BeHNaO_8Si_3$): A2B4C2D17E6_mC124_15_f_2f_f_e8f_3f 322

32. **mC144**
32.1. Catapleiite ($Na_2ZrSi_3O_9{\cdot}2H_2O$): A2B3C9D3E_mC144_15_2f_bcdef_9f_3f_ae .. 314

33. **mC168**
33.1. $(CdSO_4)_3{\cdot}8H_2O$ ($H4_{20}$): A3B16C20D3_mC168_15_ef_8f_10f_ef 342

34. **mC200**
34.1. Manganese-leonite 185 K [$K_2Mn(SO_4)_2{\cdot}4H_2O$]: A8B2CD12E2_mC200_15_8f_2f_ce_2e11f_2f 357

35. **mC212**
35.1. $Cs_6W_{11}O_{36}$: A6B36C11_mC212_9_6a_36a_11a 125

mP

1. **mP6**
1.1. O(OH)Y: ABC_mP6_11_e_e_e 139

2. **mP8**
2.1. $ZrSe_3$: A3B_mP8_11_3e_e 133

3. **mP12**
3.1. Calaverite ($AuTe_2$): AB2_mP12_7_2a_4a ... 107
3.2. Huanzalaite ($MgWO_4$, $H0_6$): AB4C_mP12_13_f_2g_e 199
3.3. Arsenopyrite (FeAsS, $E0_7$): ABC_mP12_14_e_e_e 297

4. **mP14**
4.1. $HgCl_2{\cdot}2HgO$: A2B3C2_mP14_14_e_ae_e 208

5. **mP16**
5.1. Ta_5Ti_{11} (BCC SQS-16): A5B11_mP16_6_2abc_2a3b3c 94
5.2. β-B_2H_6[*]: AB3_mP16_14_e_3e 264
5.3. B_2H_6 ($P2_1/c$)[*]: AB3_mP16_14_e_3e 266
5.4. KNO_2 III[‡]: ABC2_mP16_14_e_e_2e 285
5.5. Manganite (γ-MnO(OH), $E0_6$)[‡]: ABC2_mP16_14_e_e_2e 287
5.6. Cu(OH)Cl: ABCD_mP16_14_e_e_e_e 295
5.7. α-ICl[§]: AB_mP16_14_2e_2e 299
5.8. LiAs[§]: AB_mP16_14_2e_2e 301

6. **mP18**
6.1. $K_2S_2O_5$ ($K0_1$): A2B5C2_mP18_11_2e_e2f_2e 131

7. **mP20**
7.1. $Li_2SO_4{\cdot}H_2O$ ($H4_8$): A2B2C5D_mP20_4_2a_2a_5a_a 77
7.2. Ca_3UO_6: A3B6C_mP20_4_3a_6a_a 79
7.3. Barytocalcite ($BaCa(CO_3)_2$): AB2CD6_mP20_11_e_2e_e_2e2f 137
7.4. Orpiment (As_2S_3, $D5_f$): A2B3_mP20_14_2e_3e 210
7.5. Sanguite ($KCuCl_3$): A3BC_mP20_14_3e_e_e 224
7.6. Sr_2MnTeO_6: AB6C2D_mP20_14_a_3e_e_d .. 280
7.7. Cryolite (Na_3AlF_6, $J2_6$): AB6C3_mP20_14_a_3e_de 282

8. **mP22**
8.1. y-$Y_2Si_2O_7$: A7B2C2_mP22_11_3e2f_2e_ab .. 135
8.2. $Sb_4O_5Cl_2$: A2B5C4_mP22_14_e_c2e_2e 216
8.3. $K_2Ni(CN)_4$: A4B2C4D_mP22_14_2e_e_2e_a 230
8.4. Co_2Al_9 ($D8_d$): A9B2_mP22_14_a4e_e 252

9. **mP24**
9.1. γ-$Y_2Si_2O_7$[‖]: A4BC_mP24_14_4e_e_e 235
9.2. Anhydrous $KAuBr_4$: AB4C_mP24_14_ab_4e_e 271
9.3. Ammonium Persulfate [$(NH_4)_2S_2O_8$, $K4_1$][†‖]: AB4C_mP24_14_e_4e_e 274
9.4. Monazite ($LaPO_4$)[†‖]: AB4C_mP24_14_e_4e_e 277
9.5. $AgMnO_4$ ($H0_9$)[‖]: ABC4_mP24_14_e_e_4e .. 289

[*] β-B_2H_6 and B_2H_6 ($P2_1/c$) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[‡] KNO_2 III and Manganite (γ-MnO(OH), $E0_6$) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[§] α-ICl and LiAs have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[‖] γ-$Y_2Si_2O_7$, Ammonium Persulfate [$(NH_4)_2S_2O_8$, $K4_1$], Monazite ($LaPO_4$), and $AgMnO_4$ ($H0_9$) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

[†] Ammonium Persulfate [$(NH_4)_2S_2O_8$, $K4_1$] and Monazite ($LaPO_4$) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

9.6. Nahcolite ($NaHCO_3$, $G0_{12}$): ABCD3_mP24_14_e_e_e_3e 292
9.7. ϵ-1,2,3,4,5,6-Hexachlorocyclohexane (C_6Cl_6): AB_mP24_14_3e_3e 303

10. **mP28**
10.1. Monoclinic Cu_2OSeO_3: A2B4C_mP28_14_abe_4e_e 213
10.2. $KICl_4{\cdot}H_2O$ ($H0_{10}$): A4BCD_mP28_14_4e_e_e_e 232

11. **mP30**
11.1. Azurite [$Cu_3(CO_3)_2(OH)_2$, $G7_4$]: A2B3C2D8_mP30_14_e_ce_e_4e 205

12. **mP32**
12.1. Ca_2UO_5: A2B5C_mP32_14_2e_5e_ab 218
12.2. Gd_2SiO_5 (RE_2SiO_5 X1): A2B5C_mP32_14_2e_5e_e 221
12.3. γ-WO_3: A3B_mP32_14_6e_2e 226
12.4. $KAuBr_4{\cdot}2H_2O$ ($H4_{19}$): AB4C2D_mP32_14_e_4e_2e_e 268
12.5. Pararealgar (AsS)[¶]: AB_mP32_14_4e_4e 306
12.6. Realgar (AsS, B_l)[¶]: AB_mP32_14_4e_4e 309

13. **mP40**
13.1. K_2NbF_7 ($K6_2$): A7B2C_mP40_14_7e_2e_e . . 242

14. **mP54**
14.1. $Na_2Ca_6Si_4O_{15}$: A6B2C15D4_mP54_7_6a_2a_15a_4a 96
14.2. $K_2Pt(SCN)_6{\cdot}2H_2O$: A6B4C2D6E2FG6_mP54_14_3e_2e_e_3e_e_a_3e 238

15. **mP56**
15.1. $Cs_{11}O_3$: A11B3_mP56_14_11e_3e 201

16. **mP60**
16.1. $W_2O_3(PO_4)_2$: A11B2C2_mP60_4_22a_4a_4a . 73
16.2. Parawollastonite ($CaSiO_3$, $S3_3$ (II)): AB3C_mP60_14_3e_9e_3e 260

17. **mP62**
17.1. High-Temperature Mo_8O_{23}: A8B23_mP62_13_4g_c11g 195

18. **mP78**
18.1. Tutton salt [$Cu(NH_4)_2(SO_4)_2{\cdot}6H_2O$, $H4_4$]: AB20C2D14E2_mP78_14_a_10e_e_7e_e 255

19. **mP100**
19.1. Manganese-leonite 110 K [$K_2Mn(SO_4)_2{\cdot}4H_2O$]: A8B2CD12E2_mP100_14_8e_2e_ad_12e_2e . . 246

20. **mP124**
20.1. Low-Temperature Mo_8O_{23}: A8B23_mP124_7_16a_46a 100

oC . . . . .

1. **oC6**
1.1. $HoSb_2$: AB2_oC6_21_a_k 415

2. **oC12**
2.1. MoP_2: AB2_oC12_36_a_2a 476
2.2. Li_2PrO_3: A2B3C_oC12_65_h_bh_a 775

3. **oC16**
3.1. $D0_7$ (CrO_3) (*obsolete*): AB3_oC16_20_a_bc 409
3.2. Orthorhombic CrO_3: AB3_oC16_40_b_a2b . . 491
3.3. $MgCuAl_2$ ($E1_a$): A2BC_oC16_63_f_c_c 744
3.4. Mn_3As ($D0_d$): AB3_oC16_63_c_3c 759
3.5. Re_3B: AB3_oC16_63_c_cf 761

4. **oC20**
4.1. Si_2N_2O: A2BC2_oC20_36_b_a_b 466
4.2. Lepidocrocite (γ-FeO(OH), $E0_4$): AB2C2_oC20_63_c_f_2c 753
4.3. $ThFe_2SiC$: AB2CD_oC20_63_b_f_c_c 757
4.4. V_3AsC: ABC3_oC20_63_c_b_cf 765

5. **oC22**
5.1. Nb_3O_7F: A3B8_oC22_65_ag_bd2gh 779

6. **oC24**
6.1. $AlPO_4$ "low cristobalite type": AB4C_oC24_20_b_2c_a 411
6.2. $ZrTe_5$: A5B_oC24_63_c2f_c 751
6.3. Si_{24} Clathrate: A_oC24_63_3f 767
6.4. $NH_4H_2PO_2$ ($F5_7$): A2BC2D_oC24_67_m_a_n_g 781

7. **oC26**
7.1. $Mg(NH_3)_2Cl_2$ ($E1_3$): A2B8CD2_oC26_65_h_r_a_i 777

8. **oC28**
8.1. Na_2CrO_4 ($H1_8$): AB2C4_oC28_63_c_bc_fg . 755

9. **oC32**
9.1. Tl_2AlF_5 ($K3_3$): AB5C2_oC32_20_b_a2bc_c . 413
9.2. Bi_2GeO_5: A2BC5_oC32_36_b_a_a2b 468
9.3. Ta_3Ti_{13} (BCC SQS-16): A3B13_oC32_38_ac_a2bcdef 481
9.4. Ta_3Ti_5 (BCC SQS-16): A3B5_oC32_38_abce_abcdf 483
9.5. Pseudobrookite (Fe_2TiO_5, $E4_1$)[††]: A2B5C_oC32_63_f_c2f_c 742
9.6. Pd_5Pu_3: A5B3_oC32_63_cfg_ce 749
9.7. Ta_2NiS_5[††]: AB5C2_oC32_63_c_c2f_f 763

10. **oC46**
10.1. $NaNb_6O_{15}F$: ABC6D15_oC46_38_b_b_2a2d_2ab4d2e 485

11. **oC52**
11.1. $Cu_2Pb(SeO_3)_2Br_2$: A2B2C6DE2_oC52_63_g_e_fh_c_f 740

12. **oC60**

[¶]Pararealgar (AsS) and Realgar (AsS, B_l) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[††]Pseudobrookite (Fe_2TiO_5, $E4_1$) and Ta_2NiS_5 have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

12.1. Bertrandite ($Be_4Si_2O_7(OH)_2$, $S4_6$): A4B7C2D2_oC60_36_2b_a3b_2a_b 473
13. **oC68**
13.1. Base-centered orthorhombic $Sr_4Ru_3O_{10}$: A10B3C4_oC68_64_2dfg_ad_2d 769
14. **oC76**
14.1. $S0_4$ (Staurolite, $Fe(OH)_2Al_4Si_2O_{10}$) (*obsolete*): A4BC12D2_oC76_63_eg_c_f3gh_g 746
15. **oC80**
15.1. Ni_3Si_2: A3B2_oC80_36_4a4b_2a3b 470
15.2. α-Potassium Nitrate (KNO_3) II: ABC3_oC80_36_2ab_2ab_2a5b 478
16. **oC88**
16.1. $Rb_2Mo_2O_7$: A2B7C2_oC88_40_abc_2b6c_a3b 488
16.2. $Na_2Mo_2O_7$: A2B2C7_oC88_64_ef_df_3f2g .. 772
17. **oC104**
17.1. Santite ($KB_5O_8{\cdot}4H_2O$, $K3_5$): A5B8CD12_oC104_41_a2b_4b_a_6b 493

oF .....
1. **oF24**
1.1. Cs_2Se: A2B_oF24_43_b_a 510
2. **oF40**
2.1. Ag_2O_3**: A2B3_oF40_43_b_ab 497
2.2. Zr_2Al_3**: A3B2_oF40_43_ab_b 512
3. **oF48**
3.1. Mg_2Cu (C_b): AB2_oF48_70_g_fg 785
4. **oF56**
4.1. Thenardite [Na_2SO_4 (V), $H1_7$]: A2B4C_oF56_70_g_h_a 783
5. **oF64**
5.1. Archerite (KH_2PO_4): A2BC4D_oF64_43_b_a_2b_a 507
6. **oF88**
6.1. Predicted Phase IV $Cd_2Re_2O_7$: A2B7C2_oF88_22_k_bdefghij_k 417
6.2. Blossite (α-$Cu_2V_2O_7$): A2B7C2_oF88_43_b_a3b_b 504
7. **oF184**
7.1. Natrolite ($Na_2Al_2Si_3O_{10}{\cdot}2H_2O$, $S6_{10}$): A2B4C2D12E3_oF184_43_b_2b_b_6b_ab 499

oI .....
1. **oI8**
1.1. Ferroelectric $NaNO_2$ ($F5_5$): ABC2_oI8_44_a_a_c 516
1.2. $AgNO_2$ ($F5_{12}$): ABC2_oI8_44_a_a_d 518
1.3. CsO: AB_oI8_71_g_i 791
2. **oI12**
2.1. $CeCu_2$: AB2_oI12_74_e_h 797
3. **oI16**
3.1. $CsFeS_2$ (100 K): ABC2_oI16_71_g_i_eh 789
4. **oI20**
4.1. High-Temperature Cryolite (Na_3AlF_6): AB6C3_oI20_71_a_in_cj 787
5. **oI28**
5.1. Ga_2Mg_5 ($D8_g$): A2B5_oI28_72_j_bfj 793
5.2. $LiCuVO_4$: ABC4D_oI28_74_a_d_hi_e 799
6. **oI40**
6.1. Hemimorphite ($Zn_4Si_2O_7(OH)_2{\cdot}H_2O$, $S2_2$): A2B5CD2_oI40_44_2c_abcde_d_e 514
7. **oI44**
7.1. $Zn(NH_3)_2Cl_2$ ($E1_2$): A2B6C2D_oI44_74_h_ij_i_e 795
8. **oI48**
8.1. $B30$ (MgZn?): AB_oI48_44_6d_ab2cde 520
9. **oI100**
9.1. $Nb_2Zr_6O_{17}$: A2B17C6_oI100_46_ab_b8c_3c . 523

oP .....
1. **oP6**
1.1. RuB_2: A2B_oP6_59_f_a 591
2. **oP8**
2.1. Parkerite ($Ni_3Bi_2S_2$): AB2C_oP8_51_e_be_f 527
2.2. InS: AB_oP8_58_g_g 589
2.3. η-NiSi (B_d): AB_oP8_62_c_c 738
3. **oP10**
3.1. $MnF_{2-x}(OH)_x$: A2BC2_oP10_34_c_a_c 464
4. **oP12**
4.1. γ-TeO_2: A2B_oP12_18_2c_c 386
4.2. ζ-Fe_2N×: A2B_oP12_60_d_c 600
4.3. α-PbO_2×: A2B_oP12_60_d_c 602
4.4. $C53$ ($SrBr_2$) (*obsolete*): A2B_oP12_62_2c_c 654
4.5. MnCuP: ABC_oP12_62_c_c_c 736
5. **oP14**
5.1. $D8_7$ (Shcherbinaite, V_2O_5) (*obsolete*): A5B2_oP14_31_a2b_b 440
5.2. Shcherbinaite (V_2O_5) (*revised*): A5B2_oP14_59_a2f_f 595
6. **oP16**
6.1. NaP: AB_oP16_19_2a_2a 407
6.2. $LiGaO_2$: ABC2_oP16_33_a_a_2a 460
6.3. NH_4HF_2 ($F5_8$): A2BC_oP16_53_eh_ab_g ... 546
6.4. $D0_{10}$ (WO_3) (*obsolete*): A3B_oP16_57_a2d_d 566
6.5. NH_4I_3 ($D0_{16}$): A3B_oP16_62_3c_c 663
6.6. Diaspore ($AlOOH$, $E0_2$): ABC2_oP16_62_c_c_2c 718
6.7. NH_4ClBrI ($F5_{14}$): ABCD_oP16_62_c_c_c_c . 734
7. **oP18**

**Ag_2O_3 and Zr_2Al_3 have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

×ζ-Fe_2N and α-PbO_2 have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

7.1. Eriochalcite ($CuCl_2{\cdot}2H_2O$, $C45$): A2BC4D2_oP18_53_h_a_i_e 544
7.2. NH_4NO_3 IV ($G0_{11}$): A4B2C3_oP18_59_ef_ab_af 593

8. **oP20**
8.1. Wülfingite (ϵ-$Zn(OH)_2$, $C31$): A2B2C_oP20_19_2a_2a_a 396
8.2. γ-$LiIO_3$: ABC3_oP20_33_a_a_3a 462
8.3. NH_4CdCl_3 ($E2_4$): AB3C_oP20_62_c_3c_c 707
8.4. α-Potassium Nitrate (KNO_3) I°: ABC3_oP20_62_c_c_cd 720
8.5. NH_4NO_3 III ($G0_{10}$)°: ABC3_oP20_62_c_c_cd 723
8.6. Aragonite ($CaCO_3$, $G0_2$)°: ABC3_oP20_62_c_c_cd 726

9. **oP22**
9.1. Kotoite ($Mg_3(BO_3)_2$): A2B3C6_oP22_58_g_af_gh 574

10. **oP24**
10.1. $NaAlCl_4$: AB4C_oP24_19_a_4a_a 405
10.2. B_4SrO_7: A4B7C_oP24_31_2b_a3b_a 438
10.3. Cervantite (α-Sb_2O_4): A2B_oP24_33_4a_2a 455
10.4. $SrUO_4$: A4BC_oP24_57_cde_d_a 569
10.5. Tellurite (β-TeO_2, $C52$): A2B_oP24_61_2c_c 614
10.6. COCl: ABC_oP24_61_c_c_c 627
10.7. Cs_2Sb: A2B_oP24_62_4c_2c 656
10.8. Chalcocyanite ($CuSO_4$): AB4C_oP24_62_a_2cd_c 711

11. **oP26**
11.1. $Nb_2Pd_3Se_8$: A2B3C8_oP26_55_h_ag_2g2h 551

12. **oP28**
12.1. Mercury (II) Azide [$Hg(N_3)_2$]: AB6_oP28_29_a_6a 420
12.2. In_4Se_3: A4B3_oP28_58_4g_3g 584
12.3. CaB_2O_4 I ($E3_2$): A2BC4_oP28_60_d_c_2d 597
12.4. Ca_2RuO_4: A2B4C_oP28_61_c_2c_a 611
12.5. Arcanite (K_2SO_4, $H1_6$): A2B4C_oP28_62_2c_2cd_c 635
12.6. $VOSO_4$: A5BC_oP28_62_3cd_c_c 682
12.7. Berthierite ($FeSb_2S_4$, $E3_3$): AB4C2_oP28_62_c_4c_2c 709
12.8. Copper (II) Azide [$Cu(N_3)_2$]: AB6_oP28_62_c_6c 716

13. **oP32**
13.1. $K_2HgCl_4{\cdot}H_2O$ ($E3_4$): A4BCD2_oP32_55_ghi_f_e_gh 553
13.2. $HoMn_2O_5$: AB2C5_oP32_55_g_fh_eghi 559
13.3. Andalusite (Al_2SiO_5, $S0_2$): A2B5C_oP32_58_eg_3gh_g 576
13.4. Sillimanite (Al_2SiO_5, $S0_3$): A2B5C_oP32_62_bc_3cd_c 645
13.5. Original β-WO_3 (*obsolete*): A3B_oP32_62_ab4c_2c 665
13.6. $K_2SnCl_4{\cdot}H_2O$ ($E3_5$): A4BC2D_oP32_62_2cd_b_2c_a 676
13.7. $K_2SnCl_4{\cdot}H_2O$: A4BC2D_oP32_62_2cd_c_d_c 679

14. **oP34**
14.1. $Mg(ClO_4)_2{\cdot}6H_2O$ ($H4_{11}$): A2B6CD8_oP34_31_2a_2a2b_a_4a2b 435

15. **oP36**
15.1. Adamite [$Zn_2(AsO_4)(OH)$, $H2_7$]: ABC5D2_oP36_58_g_g_3gh_eg 586
15.2. Columbite ($FeNb_2O_6$, $E5_1$): AB2C6_oP36_60_c_d_3d 608
15.3. Topaz ($Al_2SiO_4F_2$, $S0_5$): A2B2C4D_oP36_62_d_d_2cd_c 629
15.4. Atacamite ($Cu_2(OH)_3Cl$): AB2C3D3_oP36_62_c_ac_cd_cd 704
15.5. Rynersonite (Orthorhombic $CaTa_2O_6$): AB6C2_oP36_62_c_2c2d_d 713

16. **oP38**
16.1. $Ru_{11}B_8$: A8B11_oP38_55_g3h_a3g2h 556

17. **oP40**
17.1. $NaNbO_3$: ABC3_oP40_17_abcd_2e_abcd4e 383
17.2. Diamminetriamidodizinc Chloride ($[Zn_2(NH_3)_2(NH_2)_3]Cl$): AB12C5D2_oP40_18_a_6c_b2c_c 388
17.3. Lueshite ($NaNbO_3$): ABC3_oP40_57_cd_e_cd2e 571
17.4. Norbergite [$Mg(F,OH)_2 \cdot Mg_2SiO_4$, $S0_7$]: A2B3C4D_oP40_62_d_cd_2cd_c 632

18. **oP44**
18.1. Possible δ-$Gd_2Si_2O_7$: A2B7C2_oP44_33_2a_7a_2a 447
18.2. $K_2S_3O_6$ ($K5_1$): A2B6C3_oP44_62_2c_2c2d_3c 648
18.3. Possible δ-$Y_2Si_2O_7$: A7B2C2_oP44_62_3c2d_2c_d 684
18.4. $SbCl_5{\cdot}POCl_3$: A8BCD_oP44_62_4c2d_c_c_c 692

19. **oP46**
19.1. $LiNb_6O_{15}F$: ABC6D15_oP46_51_f_d_2e2i_aef4i2j 529

20. **oP48**
20.1. Ferroelectric $NH_4H_2PO_4$: A6BC4D_oP48_19_6a_a_4a_a 398
20.2. CsB_4O_6F: A4BCD6_oP48_33_4a_a_a_6a 457

21. **oP52**
21.1. Danburite ($CaB_2Si_2O_8$, $S6_3$): A2BC8D2_oP52_62_d_c_2c3d_d 651

22. **oP56**
22.1. Calciborite (CaB_2O_4 II): A2BC4_oP56_56_2e_e_4e 562
22.2. Mo_4P_3: A4B3_oP56_62_8c_6c 673

23. **oP64**

°α-Potassium Nitrate (KNO_3) I, NH_4NO_3 III ($G0_{10}$), and Aragonite ($CaCO_3$, $G0_2$) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

23.1. Hambergite [$Be_2BO_3(OH)$, $G7_2$]: AB2CD4_oP64_61_c_2c_c_4c 620

24. **oP68**
24.1. Orthorhombic $Sr_4Ru_3O_{10}$: A10B3C4_oP68_55_2e2fgh2i_adef_2e2f 548
24.2. Cr_5O_{12}: A5B12_oP68_60_c2d_6d 604

25. **oP80**
25.1. β-Arabinose [$(CH_2O)_{20}$]: AB2C_oP80_19_5a_10a_5a 401
25.2. $(TiCl_4{\cdot}POCl_3)_2$: A7BCD_oP80_61_7c_c_c_c 616
25.3. Enstatite ($MgSiO_3$, $S4_3$): AB3C_oP80_61_2c_6c_2c 623

26. **oP82**
26.1. Protoanthophyllite ($H_2Mg_7Si_8O_{24}$): A2B7C24D8_oP82_58_g_ae2f_2g5h_2h 579

27. **oP84**
27.1. CaB_2O_4 (III): A2BC4_oP84_33_6a_3a_12a 450

28. **oP96**
28.1. $RhCl_2(NH_3)_5Cl$ ($J1_8$): A3B15C5D_oP96_62_cd_3c6d_3cd_c 658

29. **oP102**
29.1. Orthorhombic Co_4Al_{13}: A13B4_oP102_31_17a11b_8a2b 430

30. **oP108**
30.1. Morenosite ($NiSO_4{\cdot}7H_2O$, $H4_{12}$): A14BC11D_oP108_19_14a_a_11a_a 391
30.2. $K_4[Mo(CN)_8]{\cdot}2H_2O$ ($F2_1$): A8B4C4DE8F2_oP108_62_4c2d_2d_2cd_c_4c2d_d 687

31. **oP112**
31.1. P_4Se_3: A4B3_oP112_62_8c4d_4c4d 668
31.2. Epididymite ($BeHNaO_8Si_3$, $S4_7$): ABCD8E3_oP112_62_d_2c_d_4c6d_3d 728

32. **oP128**
32.1. $Mo_{17}O_{47}$: A17B47_oP128_32_a8c_a23c 442

33. **oP156**
33.1. Anthophyllite ($Mg_5Fe_2Si_8O_{22}(OH)_2$, $S4_4$): A2B5C22D2E8_oP156_62_d_c2d_2c10d_2c_4d 638

34. **oP200**
34.1. Autunite $\{Ca[(UO_2)(PO_4)]_2(H_2O)_{11}\}$: AB22C23D2E2_oP200_62_c_11d_3c10d_d_d . 695

35. **oP220**
35.1. Low-Temperature $(NH_3CH_3)Al(SO_4)_2{\cdot}12H_2O$: ABC30DE20F2_oP220_29_a_a_30a_a_20a_2a 422

36. **oP276**
36.1. Carnallite [$Mg(H_2O)_6KCl_3$]: A3B12CDE6_oP276_52_d4e_18e_ce_de_2d8e 532

tI

1. **tI4**
1.1. "Martensite Type" FeC_x ($x \leq 0.06$) ($L'2_0$): AB_tI4_139_b_a 998

2. **tI6**
2.1. CaC_2-I ($C11_a$): A2B_tI6_139_e_a 978

3. **tI8**
3.1. Tetragonal $TlFeS_2$: AB2C_tI8_119_c_e_a 877

4. **tI10**
4.1. $BaNiSn_3$: ABC3_tI10_107_a_a_ab 859
4.2. $TlCo_2S_2$[∂‡‡]: A2B2C_tI10_139_d_e_a 972
4.3. $ThCr_2Si_2$[∂‡‡]: A2B2C_tI10_139_d_e_a 974
4.4. Au_2Nb_3: A2B3_tI10_139_e_ae 976
4.5. Li_2CN_2[⊗‡‡]: AB2C2_tI10_139_a_d_e 992
4.6. $H5_9$ [Autunite, $Ca(UO_2)_2(PO_4)_2{\cdot}10\frac{1}{2}H_2O$] (*obsolete*)[⊗‡‡]: AB2C2_tI10_139_a_d_e 994

5. **tI12**
5.1. $C17$ (Fe_2B) (*obsolete*): AB2_tI12_121_ab_i 884

6. **tI14**
6.1. K_2NiF_4: A4B2C_tI14_139_ce_e_a 982

7. **tI16**
7.1. Kesterite [$Cu_2(Zn,Fe)SnS_4$]: A2BCD4_tI16_82_ac_b_d_g 805
7.2. KHF_2 ($F5_2$): A2BC_tI16_140_h_d_a 1002

8. **tI18**
8.1. $K_2OsO_2Cl_4$ ($J1_5$): A4B2C2D_tI18_139_h_d_e_a 980
8.2. Fe_8N ($D2_g$): A8B_tI18_139_deh_a 990

9. **tI20**
9.1. Sr_2NiWO_6: AB6C2D_tI20_87_a_eh_d_b 828
9.2. $AuCsCl_3$ ($K7_6$): AB3C_tI20_139_ab_eh_d ...996

10. **tI24**
10.1. Scheelite ($CaWO_4$, $H0_4$): AB4C_tI24_88_b_f_a 837
10.2. K_3CrO_8: AB3C8_tI24_121_a_bd_2i 886
10.3. $Sr_3Ti_2O_7$: A7B3C2_tI24_139_aeg_be_e 988

11. **tI28**
11.1. V_4SiSb_2: A2BC4_tI28_140_h_a_k 1000
11.2. U_6Mn ($D2_c$): AB6_tI28_140_a_hk 1014

12. **tI32**
12.1. Copper (I) Azide (CuN_3): AB3_tI32_88_d_cf 835
12.2. α-V_3S: AB3_tI32_121_g_f2i 888
12.3. $NH_4Pb_2Br_5$ ($K3_4$): A5BC2_tI32_140_bl_a_h 1010

[∂] $TlCo_2S_2$ and $ThCr_2Si_2$ have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[‡‡] $TlCo_2S_2$, $ThCr_2Si_2$, Li_2CN_2, and $H5_9$ [Autunite, $Ca(UO_2)_2(PO_4)_2{\cdot}10\frac{1}{2}H_2O$] (*obsolete*) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

[⊗] Li_2CN_2 and $H5_9$ [Autunite, $Ca(UO_2)_2(PO_4)_2{\cdot}10\frac{1}{2}H_2O$] (*obsolete*) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

13. **tI34**
13.1. $Sr_4Ti_3O_{10}$: A10B4C3_tI34_139_c2eg_2e_ae 970
14. **tI36**
14.1. Cs_3CoCl_5 ($K3_1$): A5BC3_tI36_140_cl_b_ah 1012
15. **tI40**
15.1. Mercury Cyanide [$Hg(CN)_2$, $F1_1$]: A2BC2_tI40_122_e_d_e 890
15.2. KH_2PO_4 ($H2_2$): A4BC4D_tI40_122_e_b_e_a 893
16. **tI44**
16.1. Phase III $Cd_2Re_2O_7$: A2B7C2_tI44_98_f_bcde_f 848
16.2. Phase II $Cd_2Re_2O_7$: A2B7C2_tI44_119_i_bdefgh_i 874
16.3. $SrCu_2(BO_3)_2$: A2B2C6D_tI44_121_i_i_ij_c 882
17. **tI48**
17.1. NaS_2: AB2_tI48_122_cd_2e 899
17.2. $BaCd_{11}$: AB11_tI48_141_a_bdi 1016
18. **tI56**
18.1. $NH_4H_2PO_4$: A8BC4D_tI56_122_2e_b_e_a ... 896
19. **tI72**
19.1. $BeSO_4{\cdot}4H_2O$ ($H4_3$): AB8C8D_tI72_120_c_2i_2i_b 879
20. **tI82**
20.1. Marialite Scapolite [$Na_4Cl(AlSi_3)_3O_{24}$, $S6_4$]: AB4C24D12_tI82_87_a_h_2h2i_hi 825
21. **tI132**
21.1. $Na_4Ge_9O_{20}$: A9B4C20_tI132_88_a2f_f_5f ... 830
22. **tI160**
22.1. Cd_3As_2: A2B3_tI160_142_deg_3g 1025
23. **tI168**
23.1. $K_3TlCl_6{\cdot}2H_2O$ ($J3_1$): A6B2C3D_tI168_139_egikl2m_ejn_bh2n_acf . 984
24. **tI184**
24.1. Analcime ($NaAlSi_2O_6{\cdot}H_2O$, $S6_1$): A2B2C3D12E4_tI184_142_f_f_be_3g_g 1019
25. **tI204**
25.1. $Pu_{31}Rh_{20}$: A31B20_tI204_140_b2gh3m_ac2fh3l 1004

tP
1. **tP2**
1.1. δ-CuTi ($L2_a$): AB_tP2_123_a_d 912
2. **tP4**
2.1. $F6_1$ (Chalcopyrite, $CuFeS_2$) (*obsolete*): ABC2_tP4_115_a_c_g 872
3. **tP5**
3.1. NH_4HgCl_3 ($E2_5$): A3BC_tP5_123_cg_a_d .. 902
4. **tP6**
4.1. $TlAlF_4$ ($H0_8$): AB4C_tP6_123_d_eh_a 910
5. **tP7**
5.1. $E3_1$ (β-Ag_2HgI_4) (*obsolete*): A2BC4_tP7_111_f_a_n 861
5.2. K_2PtCl_4 ($H1_5$): A4B2C_tP7_123_j_e_a 904
6. **tP8**
6.1. $F5_4$ (NH_4ClO_2) (*obsolete*): ABC2_tP8_100_b_a_c 851
6.2. LaOAgS: ABCD_tP8_129_b_c_a_c 950
7. **tP10**
7.1. NH_4NO_3 II ($G0_9$): ABC3_tP10_100_b_a_bc . 853
7.2. $CaBe_2Ge_2$: A2BC2_tP10_129_ac_c_bc 943
8. **tP11**
8.1. $E6_1$ ($Sr(OH)_2(H_2O)_8$) (*obsolete*): A8B2C_tP11_123_r_f_a 906
8.2. $E6_2$ [$SrO_2(H_2O)_8$] (*possibly obsolete*): A8B2C_tP11_123_r_h_a 908
9. **tP12**
9.1. Paratellurite (α-TeO_2): A2B_tP12_92_b_a 846
9.2. NH_4Br ($B25$): AB4C_tP12_129_c_i_a 948
10. **tP14**
10.1. $MoPO_5$: AB5C_tP14_85_c_cg_b 810
10.2. $ZrFe_4Si_2$: A4B2C_tP14_136_i_g_b 962
11. **tP16**
11.1. $G7_5$ ($PbCO_3 \cdot PbCl_2$, Phosgenite) (*obsolete*): AB2C3D2_tP16_90_c_f_ce_e 839
11.2. Ammonium Chlorite (NH_4ClO_2): AB4CD2_tP16_113_c_f_a_e 863
11.3. $Pd(NH_3)_4Cl_2{\cdot}H_2O$ ($H4_9$): A2BC4D_tP16_127_h_d_i_a 929
11.4. α-WO_3: A3B_tP16_130_cf_c 957
12. **tP20**
12.1. Zr_3Al_2: A2B3_tP20_136_j_dfg 960
13. **tP22**
13.1. $CaO_2(H_2O)_8$: AB8C2_tP22_124_a_n_h 914
14. **tP26**
14.1. Meta-autunite (I) [$Ca(UO_2)_2(PO_4)_2{\cdot}6H_2O$, $H5_{10}$]: AB4C6DE_tP26_129_c_j_2ci_a_c 945
14.2. $K_2CuCl_4{\cdot}2H_2O$ ($H4_1$): A4BC4D2E2_tP26_136_fg_a_j_d_e 964
15. **tP32**
15.1. $PNCl_2$ ($E1_4$): A2BC_tP32_86_2g_g_g 812
15.2. $NaSb(OH)_6$ ($J1_{11}$): AB6C_tP32_86_d_3g_c . 819
15.3. $Ag[Co(NH_3)_2(NO_2)_4]$ ($J1_9$): ABC4D2E8_tP32_126_a_b_h_e_k 926
15.4. Phosgenite [$Pb_2Cl_2(CO_3)$]: AB2C3D2_tP32_127_g_eh_gk_k 931
16. **tP40**
16.1. β-$LiIO_3$: ABC3_tP40_86_g_g_3g 822
17. **tP44**
17.1. Bromocarnallite ($KMg(H_2O)_6(Cl,Br)_3$, $E2_6$): A3B6CD_tP44_85_bcg_3g_ac_e 807
17.2. Chiolite ($Na_5Al_3F_{14}$, $K7_5$): A3B14C5_tP44_128_ac_ehi_bg 934
18. **tP46**

18.1. $Ag_2SO_4{\cdot}4NH_3$ ($H4_{17}$): A2B12C4D4E_tP46_114_d_3e_e_e_a 869

19. **tP68**

19.1. $Nd_4Re_2O_{11}$: A4B11C2_tP68_86_2g_ab5g_g . 815

19.2. $C_{19}Sc_{15}$: A19B15_tP68_114_bc4e_ac3e 865

19.3. $Nd_2Fe_{14}B$: AB14C2_tP68_136_f_ce2j2k_fg .. 966

20. **tP72**

20.1. Gwihabaite [NH_4NO_3 (V)]: A4B2C3_tP72_77_8d_ab2c2d_6d 801

20.2. VSe_2O_6: A6B2C_tP72_103_abc5d_2d_abc ... 855

21. **tP96**

21.1. Retgersite (α-$NiSO_4{\cdot}6H_2O$, $H4_5$): A12BC10D_tP96_92_6b_a_5b_a 841

22. **tP116**

22.1. Apophyllite ($KCa_4Si_8O_{20}F{\cdot}8H_2O$, $S5_2$): A4BC16DE28F8_tP116_128_h_a_2i_b_g3i_i . 937

22.2. $Sr(OH)_2(H_2O)_8$: A18B10C_tP116_130_2c4g_2c2g_a 952

23. **tP252**

23.1. Vesuvianite ($Ca_{10}Al_4(Mg,Fe)_2Si_9O_{34}(OH)_4$, $S2_3$): A4B10C2D34E4F9_tP252_126_k_ce2k_f_h8k_k_d2k 916

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1. K(SH) (*B22*): AB_hR2_166_a_b 1140

B23 .....

1. α-AgI (*B23*): A21B_cI44_229_bdh_a 1490

B25 .....

1. NH_4Br (*B25*): AB4C_tP12_129_c_i_a 948

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1. *B30* (MgZn?): AB_oI48_44_6d_ab2cde 520

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1. $LiOH{\cdot}H_2O$ (*B36*): A3BC2_mC24_12_ij_h_gi 161

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1. CaC_2-I ($C11_a$): A2B_tI6_139_e_a 978

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1. $CaSi_2$ (*C12*): AB2_hR6_166_c_2c 1124

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1. *C17* (Fe_2B) (*obsolete*): AB2_tI12_121_ab_i 884

C26 .....

1. NO_2 (Modern, *C26*): AB2_cI36_204_d_g 1296

$C26_a$ .....

1. $C26_a$ (NO_2) (*obsolete*): AB2_cI36_199_b_c 1282

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1. *C27* (CdI_2) (*questionable*): AB2_hP6_186_b_ab 1213

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1. Wülfingite (ϵ-$Zn(OH)_2$, *C31*): A2B2C_oP20_19_2a_2a_a 396

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1. Eriochalcite ($CuCl_2{\cdot}2H_2O$, *C45*): A2BC4D2_oP18_53_h_a_i_e 544

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1. Tellurite (β-TeO_2, *C52*): A2B_oP24_61_2c_c 614

C53 .....

1. *C53* ($SrBr_2$) (*obsolete*): A2B_oP12_62_2c_c 654

C_b .....

1. Mg_2Cu (C_b): AB2_oF48_70_g_fg 785

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1. $LiZn_2$ (C_k)[∂∂]: AB_hP4_194_a_c 1270

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1. $D0_6$ (Tysonite, LaF_3) (*obsolete*): A3B_hP24_193_ack_g 1238

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1. $D0_7$ (CrO_3) (*obsolete*): AB3_oC16_20_a_bc 409

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1. $D0_{10}$ (WO_3) (*obsolete*): A3B_oP16_57_a2d_d 566

$D0_{12}$ .....

1. FeF_3 ($D0_{12}$): A3B_hR8_167_e_b 1159

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1. $D0_{13}$ ($AlCl_3$) (*obsolete*): AB3_hP4_164_b_ad 1107

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1. NH_4I_3 ($D0_{16}$): A3B_oP16_62_3c_c 663

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1. Mn_3As ($D0_d$): AB3_oC16_63_c_3c 759

$D1_1$ .....

1. SnI_4 ($D1_1$): A4B_cP40_205_cd_c 1313

$D1_g$ .....

[¶]Pararealgar (AsS) and Realgar (AsS, B_l) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[∂∂]$LiZn_2$ (C_k) and Fe_2N (approximate, $L'3_0$) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

1. $B_{13}C_2$ "B_4C" ($D1_g$): A13B2_hR15_166_b2h_c .. 1111

$D2_2$

1. $D2_2$ ($MgZn_5$?) (*problematic*): AB5_mC48_12_2i_ac5i2j 175

$D2_a$

1. $TiBe_{12}$ (approximate, $D2_a$): A12B_hP13_191_cdei_a 1233

$D2_c$

1. U_6Mn ($D2_c$): AB6_tI28_140_a_hk 1014

$D2_g$

1. Fe_8N ($D2_g$): A8B_tI18_139_deh_a 990

$D5_5$

1. Mg_3P_2 ($D5_5$): A3B2_cP10_224_d_b 1427

$D5_6$

1. β-Alumina (Al_2O_3, $D5_6$): A2B3_hP60_194_3fk_cdef2k 1251

$D5_7$

1. Maghemite (γ-Fe_2O_3, $D5_7$): A2B3_cP60_212_bcd_ace 1354

$D5_b$

1. Pt_2Sn_3 ($D5_b$): A2B3_hP10_194_f_bf 1249

$D5_f$

1. Orpiment (As_2S_3, $D5_f$): A2B3_mP20_14_2e_3e .. 210

$D6_1$

1. Senarmontite (Sb_2O_3, $D6_1$): A3B2_cF80_227_f_e 1481

$D6_2$

1. $D6_2$ (Sb_2O_4) (*obsolete*): A2B_cF96_227_abf_cd 1478

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1. δ-Ni_3Sn_4 ($D7_a$): A3B4_mC14_12_ai_2i 159

$D8_7$

1. $D8_7$ (Shcherbinaite, V_2O_5) (*obsolete*): A5B2_oP14_31_a2b_b 440

$D8_9$

1. Co_9S_8 ($D8_9$): A9B8_cF68_225_af_ce 1446

$D8_d$

1. Co_2Al_9 ($D8_d$): A9B2_mP22_14_a4e_e 252

$D8_g$

1. Ga_2Mg_5 ($D8_g$): A2B5_oI28_72_j_bfj 793

$D8_k$

1. Th_7S_{12} ($D8_k$): A3B2_hP20_176_2h_ah 1185

$E0_2$

1. Diaspore (AlOOH, $E0_2$): ABC2_oP16_62_c_c_2c . 718

$E0_3$

1. Cd(OH)Cl ($E0_3$): ABCD_hP8_186_b_b_a_a 1218

$E0_4$

1. Lepidocrocite (γ-FeO(OH), $E0_4$): AB2C2_oC20_63_c_f_2c 753

$E0_6$

1. Manganite (γ-MnO(OH), $E0_6$)[‡]: ABC2_mP16_14_e_e_2e 287

$E0_7$

1. Arsenopyrite (FeAsS, $E0_7$): ABC_mP12_14_e_e_e 297

$E1_2$

1. $Zn(NH_3)_2Cl_2$ ($E1_2$): A2B6C2D_oI44_74_h_ij_i_e . 795

$E1_3$

1. $Mg(NH_3)_2Cl_2$ ($E1_3$): A2B8CD2_oC26_65_h_r_a_i 777

$E1_4$

1. $PNCl_2$ ($E1_4$): A2BC_tP32_86_2g_g_g 812

$E1_a$

1. $MgCuAl_2$ ($E1_a$): A2BC_oC16_63_f_c_c 744

$E2_3$

1. $E2_3$ ($LiIO_3$) (*obsolete*): ABC3_hP10_182_c_b_g 1202

$E2_4$

1. NH_4CdCl_3 ($E2_4$): AB3C_oP20_62_c_3c_c 707

$E2_5$

1. NH_4HgCl_3 ($E2_5$): A3BC_tP5_123_cg_a_d 902

$E2_6$

1. Bromocarnallite ($KMg(H_2O)_6(Cl,Br)_3$, $E2_6$): A3B6CD_tP44_85_bcg_3g_ac_e 807

$E3_1$

1. $E3_1$ (β-Ag_2HgI_4) (*obsolete*): A2BC4_tP7_111_f_a_n 861

$E3_2$

1. CaB_2O_4 I ($E3_2$): A2BC4_oP28_60_d_c_2d 597

$E3_3$

1. Berthierite ($FeSb_2S_4$, $E3_3$): AB4C2_oP28_62_c_4c_2c 709

$E3_4$

1. $K_2HgCl_4{\cdot}H_2O$ ($E3_4$): A4BCD2_oP32_55_ghi_f_e_gh 553

$E3_5$

1. $K_2SnCl_4{\cdot}H_2O$ ($E3_5$): A4BC2D_oP32_62_2cd_b_2c_a 676

$E4_1$

1. Pseudobrookite (Fe_2TiO_5, $E4_1$)[††]: A2B5C_oC32_63_f_c2f_c 742

$E5_1$

1. Columbite ($FeNb_2O_6$, $E5_1$): AB2C6_oP36_60_c_d_3d 608

$E6_1$

[‡] KNO_2 III and Manganite (γ-MnO(OH), $E0_6$) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[††] Pseudobrookite (Fe_2TiO_5, $E4_1$) and Ta_2NiS_5 have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

1. $E6_1$ ($Sr(OH)_2(H_2O)_8$) (*obsolete*):
A8B2C_tP11_123_r_f_a 906
$E6_2$ .....
1. $E6_2$ [$SrO_2(H_2O)_8$] (*possibly obsolete*):
A8B2C_tP11_123_r_h_a 908
$E9_2$ .....
1. Swedenborgite ($NaBe_4SbO_7$, $E9_2$):
A4BC7D_hP26_186_ac_b_a2c_b 1210
$F1_1$ .....
1. Mercury Cyanide [$Hg(CN)_2$, $F1_1$]:
A2BC2_tI40_122_e_d_e 890
$F2_1$ .....
1. $K_4[Mo(CN)_8]\cdot 2H_2O$ ($F2_1$):
A8B4C4DE8F2_oP108_62_4c2d_2d_2cd_c_4c2d_d . 687
$F4_1$ .....
1. $Fe_2(CO)_9$ ($F4_1$): A9B2C9_hP40_176_hi_f_hi 1192
$F5_2$ .....
1. KHF_2 ($F5_2$): A2BC_tI16_140_h_d_a 1002
$F5_4$ .....
1. $F5_4$ (NH_4ClO_2) (*obsolete*):
ABC2_tP8_100_b_a_c 851
$F5_5$ .....
1. Ferroelectric $NaNO_2$ ($F5_5$):
ABC2_oI8_44_a_a_c 516
$F5_7$ .....
1. $NH_4H_2PO_2$ ($F5_7$): A2BC2D_oC24_67_m_a_n_g .. 781
$F5_8$ .....
1. NH_4HF_2 ($F5_8$): A2BC_oP16_53_eh_ab_g 546
$F5_{11}$ .....
1. $F5_{11}$ (KNO_2) (*obsolete*):
ABC2_mC8_8_a_a_b 117
$F5_{12}$ .....
1. $AgNO_2$ ($F5_{12}$): ABC2_oI8_44_a_a_d 518
$F5_{14}$ .....
1. NH_4ClBrI ($F5_{14}$): ABCD_oP16_62_c_c_c_c 734
$F5_a$ .....
1. $KFeS_2$ ($F5_a$): ABC2_mC16_15_e_e_f 376
$F6_1$ .....
1. $F6_1$ (Chalcopyrite, $CuFeS_2$) (*obsolete*):
ABC2_tP4_115_a_c_g 872
$G0_2$ .....
1. Aragonite ($CaCO_3$, $G0_2$)°:
ABC3_oP20_62_c_c_cd 726
$G0_7$ .....
1. $KBrO_3$ ($G0_7$)□: ABC3_hR5_160_a_a_b 1074
$G0_8$ .....
1. NH_4NO_3 I ($G0_8$): AB_cP2_221_a_b 1418
$G0_9$ .....
1. NH_4NO_3 II ($G0_9$): ABC3_tP10_100_b_a_bc 853
$G0_{10}$ .....
1. NH_4NO_3 III ($G0_{10}$)°: ABC3_oP20_62_c_c_cd 723
$G0_{11}$ .....
1. NH_4NO_3 IV ($G0_{11}$): A4B2C3_oP18_59_ef_ab_af . 593
$G0_{12}$ .....
1. Nahcolite ($NaHCO_3$, $G0_{12}$):
ABCD3_mP24_14_e_e_e_3e 292
$G1_1$ .....
1. Dolomite [$MgCa(CO_3)_2$, $G1_1$]:
A2BCD6_hR10_148_c_a_b_f 1039
$G2_1$ .....
1. $Pb(NO_3)_2$ ($G2_1$): A2B6C_cP36_205_c_d_a 1306
$G2_2$ .....
1. $Nd(BrO_3)_3\cdot 9H_2O$ ($G2_2$):
A3B9CD9_hP44_186_c_3c_b_cd 1207
$G3_2$ .....
1. Na_2SO_3 ($G3_2$): A2B3C_hP12_147_abd_g_d 1037
$G5_1$ .....
1. Boric Acid (H_3BO_3, $G5_1$):
AB3C3_aP28_2_2i_6i_6i 62
$G5_2$ .....
1. $Al(PO_3)_3$ ($G5_2$): AB9C3_cI208_220_c_3e_e 1405
$G7_1$ .....
1. Bastnäsite [$CeF(CO_3)$]:
ABCD3_hP36_190_h_g_af_hi 1230
$G7_2$ .....
1. Hambergite [$Be_2BO_3(OH)$, $G7_2$]:
AB2CD4_oP64_61_c_2c_c_4c 620
$G7_3$ .....
1. $G7_3$ [Northupite, $Na_3MgCl(CO_3)_2$] (*obsolete*):
A2BCD3E6_cF208_227_e_c_d_f_g 1475
$G7_4$ .....
1. Azurite [$Cu_3(CO_3)_2(OH)_2$, $G7_4$]:
A2B3C2D8_mP30_14_e_ce_e_4e 205
$G7_5$ .....
1. $G7_5$ ($PbCO_3\cdot PbCl_2$, Phosgenite) (*obsolete*):
AB2C3D2_tP16_90_c_f_ce_e 839
$H0_4$ .....
1. Scheelite ($CaWO_4$, $H0_4$): AB4C_tI24_88_b_f_a .. 837
$H0_5$ .....
1. High-Temperature Cubic $KClO_4$ ($H0_5$):
ABC4_cF24_216_b_a_e 1380
$H0_6$ .....
1. Huanzalaite ($MgWO_4$, $H0_6$):
AB4C_mP12_13_f_2g_e 199
$H0_8$ .....
1. $TlAlF_4$ ($H0_8$): AB4C_tP6_123_d_eh_a 910
$H0_9$ .....

°α-Potassium Nitrate (KNO_3) I, NH_4NO_3 III ($G0_{10}$), and Aragonite ($CaCO_3$, $G0_2$) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

□$KBrO_3$ ($G0_7$) and γ-Potassium Nitrate (KNO_3) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

1. $AgMnO_4$ ($H0_9$)[‖]: ABC4_mP24_14_e_e_4e 289

$H0_{10}$

1. $KICl_4{\cdot}H_2O$ ($H0_{10}$): A4BCD_mP28_14_4e_e_e_e 232

$H1_4$

1. $LiKSO_4$ ($H1_4$): ABC4D_hP14_173_a_b_bc_b 1180

$H1_5$

1. K_2PtCl_4 ($H1_5$): A4B2C_tP7_123_j_e_a 904

$H1_6$

1. Arcanite (K_2SO_4, $H1_6$): A2B4C_oP28_62_2c_2cd_c 635

$H1_7$

1. Thenardite [Na_2SO_4 (V), $H1_7$]: A2B4C_oF56_70_g_h_a 783

$H1_8$

1. Na_2CrO_4 ($H1_8$): AB2C4_oC28_63_c_bc_fg 755

$H2_2$

1. KH_2PO_4 ($H2_2$): A4BC4D_tI40_122_e_b_e_a 893

$H2_7$

1. Adamite [$Zn_2(AsO_4)(OH)$, $H2_7$]: ABC5D2_oP36_58_g_g_3gh_eg 586

$H2_8$

1. $BaAl_2O_4$ ($H2_8$): A2BC6_hP18_182_f_b_gh 1200

$H3_2$

1. Steklite [$KAl(SO_4)_2$, $H3_2$]: ABC8D2_hP12_150_b_a_dg_d 1060

$H4_1$

1. $K_2CuCl_4{\cdot}2H_2O$ ($H4_1$): A4BC4D2E2_tP26_136_fg_a_j_d_e 964

$H4_3$

1. $BeSO_4{\cdot}4H_2O$ ($H4_3$): AB8C8D_tI72_120_c_2i_2i_b 879

$H4_4$

1. Tutton salt [$Cu(NH_4)_2(SO_4)_2{\cdot}6H_2O$, $H4_4$]: AB20C2D14E2_mP78_14_a_10e_e_7e_e 255

$H4_5$

1. Retgersite (α-$NiSO_4{\cdot}6H_2O$, $H4_5$): A12BC10D_tP96_92_6b_a_5b_a 841

$H4_6$

1. Gypsum ($CaSO_4{\cdot}2H_2O$, $H4_6$): AB4C6D_mC48_15_e_2f_3f_e 366

$H4_7$

1. Bassanite [$CaSO_4(H_2O)_{0.5}$, $H4_7$]: A2B2C9D2_mC90_5_ab2c_3c_b13c_3c 81

$H4_8$

1. $Li_2SO_4{\cdot}H_2O$ ($H4_8$): A2B2C5D_mP20_4_2a_2a_5a_a 77

$H4_9$

1. $Pd(NH_3)_4Cl_2{\cdot}H_2O$ ($H4_9$): A2BC4D_tP16_127_h_d_i_a 929

$H4_{10}$

1. Chalcanthite ($CuSO_4{\cdot}5H_2O$, $H4_{10}$): AB10C9D_aP42_2_ae_10i_9i_i 58

$H4_{11}$

1. $Mg(ClO_4)_2{\cdot}6H_2O$ ($H4_{11}$): A2B6CD8_oP34_31_2a_2a2b_a_4a2b 435

$H4_{12}$

1. Morenosite ($NiSO_4{\cdot}7H_2O$, $H4_{12}$): A14BC11D_oP108_19_14a_a_11a_a 391

$H4_{13}$

1. α-Alum [$KAl(SO_4)_2 \cdot 12H_2O$, $H4_{13}$]: AB24CD28E2_cP224_205_a_4d_b_2c4d_c 1335

$H4_{14}$

1. β-Alum [$Al(NH_3CH_3)_2(SO_4)_2{\cdot}12H_2O$, $H4_{14}$]: AB2C36D2E20F2_cP252_205_a_c_6d_c_c3d_c 1344

$H4_{15}$

1. γ-Alum [$AlNa(SO_4)_2{\cdot}12H_2O$, $H4_{15}$]: AB24CD20E2_cP192_205_a_4d_b_c3d_c 1327

$H4_{16}$

1. 12-phosphotungstic acid [$H_3PW_{12}O_{40}{\cdot}5H_2O$ ($H4_{16}$)]: A5B40CD12_cP116_224_cd_e3k_a_k 1434

$H4_{17}$

1. $Ag_2SO_4{\cdot}4NH_3$ ($H4_{17}$): A2B12C4D4E_tP46_114_d_3e_e_e_a 869

$H4_{18}$

1. $LiClO_4{\cdot}3H_2O$ ($H4_{18}$): AB6CD7_hP30_186_b_d_a_b2c 1215

$H4_{19}$

1. $KAuBr_4{\cdot}2H_2O$ ($H4_{19}$): AB4C2D_mP32_14_e_4e_2e_e 268

$H4_{20}$

1. $(CdSO_4)_3{\cdot}8H_2O$ ($H4_{20}$): A3B16C20D3_mC168_15_ef_8f_10f_ef 342

$H4_{21}$

1. $H_3PW_{12}O_{40}{\cdot}29H_2O$ ($H4_{21}$): A29B40CD12_cF656_227_ae2fg_e3g_b_g 1467

$H4_{22}$

1. $BaNi(CN)_4{\cdot}4H_2O$ ($H4_{22}$): AB4C4D4E_mC56_15_e_2f_2f_2f_a 363

$H4_{23}$

1. Manganese-leonite [$K_2Mn(SO_4)_2{\cdot}4H_2O$, $H4_{23}$]: A8B2CD15E2_mC112_12_2i3j_j_ad_g4i5j_2i 167

$H5_6$

1. $H5_6$ [Tychite, $Na_6Mg_2SO_4(CO_3)_4$)] (*obsolete*): A4B2C6D16E_cF232_227_e_d_f_eg_a 1484

$H5_7$

1. Fluorapatite [$Ca_5F(PO_4)_3$, $H5_7$]: A5BC12D3_hP42_176_fh_a_2hi_h 1189

$H5_8$

[‖] γ-$Y_2Si_2O_7$, Ammonium Persulfate [$(NH_4)_2S_2O_8$, $K4_1$], Monazite ($LaPO_4$), and $AgMnO_4$ ($H0_9$) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

1. Sulphohalite [$Na_6ClF(SO_4)_2$, $H5_8$]:
ABC6D8E2_cF72_225_b_a_e_f_c 1453

***H5*$_9$** .....
1. $H5_9$ [Autunite, $Ca(UO_2)_2(PO_4)_2{\cdot}10\frac{1}{2}H_2O$] (*obsolete*)[⊗‡‡]:
AB2C2_tI10_139_a_d_e 994

***H5*$_{10}$** .....
1. Meta-autunite (I) [$Ca(UO_2)_2(PO_4)_2{\cdot}6H_2O$, $H5_{10}$]:
AB4C6DE_tP26_129_c_j_2ci_a_c 945

***H6*$_2$** .....
1. $K_2Sn(OH)_6$ ($H6_2$): A6B2C6D_hR15_148_f_c_f_a 1041

***H6*$_3$** .....
1. $K_2Pt(SCN)_6$ ($H6_3$)[××]: A2BC6_hP9_164_d_a_i .. 1099

***H6*$_4$** .....
1. $H6_4$ [$Ni(NO_3)_2(NH_3)_6$] (*obsolete*)[⊗⊗]:
A2B6CD6_cP60_205_c_d_a_d 1302

***I1*$_3$** .....
1. $I1_3$ ($SrCl_2{\cdot}(H_2O)_6$) (*obsolete*)[§§]:
A2B6C_hP9_162_d_k_a 1083

***I6*$_1$** .....
1. $Ni(H_2O)_6SnCl_6$ ($I6_1$):
A6B6CD_hR14_148_f_f_b_a 1044

***J1*$_5$** .....
1. $K_2OsO_2Cl_4$ ($J1_5$): A4B2C2D_tI18_139_h_d_e_a .. 980

***J1*$_6$** .....
1. Bararite (Trigonal $(NH_4)_2SiF_6$, $J1_6$)[××]:
A6B2C_hP9_164_i_d_a 1101

***J1*$_7$** .....
1. Bischofite ($MgCl_2{\cdot}6H_2O$, $J1_7$):
A2B12CD6_mC42_12_i_2i2j_a_ij 146

***J1*$_8$** .....
1. $RhCl_2(NH_3)_5Cl$ ($J1_8$):
A3B15C5D_oP96_62_cd_3c6d_3cd_c 658

***J1*$_9$** .....
1. $Ag[Co(NH_3)_2(NO_2)_4]$ ($J1_9$):
ABC4D2E8_tP32_126_a_b_h_e_k 926

***J1*$_{10}$** .....
1. $Zn(BrO_3)_2{\cdot}6H_2O$ ($J1_{10}$)[⊗⊗]:
A2B6C6D_cP60_205_c_d_d_a 1299

***J1*$_{11}$** .....
1. $NaSb(OH)_6$ ($J1_{11}$): AB6C_tP32_86_d_3g_c 819

***J1*$_{12}$** .....
1. $NaSbF_4(OH)_2$ ($J1_{12}$): A6BC_hP16_163_i_b_c ... 1087

***J1*$_{13}$** .....
1. K_2GeF_6 ($J1_{13}$)[××]: A6BC2_hP9_164_i_a_d 1103

***J2*$_1$** .....
1. $(NH_4)_3AlF_6$ ($J2_1$):
AB30C16D3_cF200_225_a_ej_2f_bc 1448

***J2*$_2$** .....
1. $CrCl_3(H_2O)_6$ ($J2_2$): A3BC6_hR20_167_e_b_f ... 1156

***J2*$_3$** .....
1. $Ca_3Al_2(OH)_{12}$ ($J2_3$):
A2B3C12D12_cI232_230_a_c_h_h 1492

***J2*$_4$** .....
1. $K_3Co(NO_2)_6$ ($J2_4$):
AB3C6D12_cF88_202_a_bc_e_h 1289

***J2*$_5$** .....
1. $Cu_3[Fe(CN)_6]_2{\cdot}xH_2O$ ($J2_5$, $x \approx 3$):
A6B9CD2E6_cF96_225_e_bf_a_c_e 1443

***J2*$_6$** .....
1. Cryolite (Na_3AlF_6, $J2_6$):
AB6C3_mP20_14_a_3e_de 282

***J3*$_1$** .....
1. $K_3TlCl_6{\cdot}2H_2O$ ($J3_1$):
A6B2C3D_tI168_139_egikl2m_ejn_bh2n_acf 984

***K0*$_1$** .....
1. $K_2S_2O_5$ ($K0_1$): A2B5C2_mP18_11_2e_e2f_2e 131

***K1*$_1$** .....
1. KSO_3 ($K1_1$): AB3C_hP30_150_ef_3g_c2d 1057

***K1*$_2$** .....
1. $CsSO_3$ ($K1_2$): AB3C_hP20_190_ac_i_f 1227

***K3*$_1$** .....
1. Cs_3CoCl_5 ($K3_1$): A5BC3_tI36_140_cl_b_ah 1012

***K3*$_3$** .....
1. Tl_2AlF_5 ($K3_3$): AB5C2_oC32_20_b_a2bc_c 413

***K3*$_4$** .....
1. $NH_4Pb_2Br_5$ ($K3_4$): A5BC2_tI32_140_bl_a_h 1010

***K3*$_5$** .....
1. Santite ($KB_5O_8{\cdot}4H_2O$, $K3_5$):
A5B8CD12_oC104_41_a2b_4b_a_6b 493

***K4*$_1$** .....

[⊗] Li_2CN_2 and $H5_9$ [Autunite, $Ca(UO_2)_2(PO_4)_2{\cdot}10\frac{1}{2}H_2O$] (*obsolete*) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[‡‡] $TlCo_2S_2$, $ThCr_2Si_2$, Li_2CN_2, and $H5_9$ [Autunite, $Ca(UO_2)_2(PO_4)_2{\cdot}10\frac{1}{2}H_2O$] (*obsolete*) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

[××] $K_2Pt(SCN)_6$ ($H6_3$), Bararite (Trigonal $(NH_4)_2SiF_6$, $J1_6$), and K_2GeF_6 ($J1_{13}$) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

[⊗⊗] $Zn(BrO_3)_2{\cdot}6H_2O$ ($J1_{10}$) and $H6_4$ [$Ni(NO_3)_2(NH_3)_6$] (*obsolete*) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

[§§] $I1_3$ ($SrCl_2{\cdot}(H_2O)_6$) (*obsolete*) and Rosiaite ($PbSb_2O_6$) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

1. Ammonium Persulfate [$(NH_4)_2S_2O_8$, $K4_1$][†‖]: AB4C_mP24_14_e_4e_e 274

$K5_1$ .....
1. $K_2S_3O_6$ ($K5_1$): A2B6C3_oP44_62_2c_2c2d_3c ... 648

$K6_1$ .....
1. ZrP_2O_7 High-Temperature ($K6_1$): A7B2C_cP40_205_bd_c_a 1319

$K6_2$ .....
1. K_2NbF_7 ($K6_2$): A7B2C_mP40_14_7e_2e_e 242

$K7_1$ .....
1. $K_3W_2Cl_9$ ($K7_1$): A9B3C2_hP28_176_hi_af_f 1195

$K7_2$ .....
1. $Cs_3Tl_2Cl_9$ ($K7_2$): A9B3C2_hR28_167_ef_e_c 1164

$K7_3$ .....
1. $Cs_3As_2Cl_9$ ($K7_3$): A2B9C3_hP14_150_d_eg_ad .. 1052

$K7_4$ .....
1. Mayenite ($12CaO{\cdot}7Al_2O_3$, $K7_4$, C12A7): A7B12C19_cI152_220_bc_2d_ace 1400

$K7_5$ .....
1. Chiolite ($Na_5Al_3F_{14}$, $K7_5$): A3B14C5_tP44_128_ac_ehi_bg 934

$K7_6$ .....
1. $AuCsCl_3$ ($K7_6$): AB3C_tI20_139_ab_eh_d 996

$L1_3(I)$ .....
1. Cubic CuPt ($L1_3$ (I), $D4$): AB_cF32_227_c_d ... 1488

$L1_a$ .....
1. $L1_a$ (disputed $CuPt_3$): AB7_cF32_225_b_ad 1451

$L2_a$ .....
1. δ-CuTi ($L2_a$): AB_tP2_123_a_d 912

$L'1_0$ .....
1. γ-Fe_4N ($L'1_0$): A4B_cP5_221_bc_a 1414

$L'2_0$ .....
1. "Martensite Type" FeC_x ($x \leq 0.06$) ($L'2_0$): AB_tI4_139_b_a 998

$L'3_0$ .....
1. Fe_2N (approximate, $L'3_0$)[∂∂]: AB_hP4_194_c_a .. 1272

$S0_1$ .....
1. Kyanite (Al_2SiO_5, $S0_1$): A2B5C_aP32_2_4i_10i_2i 44

$S0_2$ .....
1. Andalusite (Al_2SiO_5, $S0_2$): A2B5C_oP32_58_eg_3gh_g 576

$S0_3$ .....
1. Sillimanite (Al_2SiO_5, $S0_3$): A2B5C_oP32_62_bc_3cd_c 645

$S0_4$ .....
1. $S0_4$ (Staurolite, $Fe(OH)_2Al_4Si_2O_{10}$) (*obsolete*): A4BC12D2_oC76_63_eg_c_f3gh_g 746

$S0_5$ .....
1. Topaz ($Al_2SiO_4F_2$, $S0_5$): A2B2C4D_oP36_62_d_d_2cd_c 629

$S0_6$ .....
1. Titanite ($CaTiSiO_5$, $S0_6$): AB5CD_mC32_15_e_e2f_e_b 374

$S0_7$ .....
1. Norbergite [$Mg(F,OH)_2 \cdot Mg_2SiO_4$, $S0_7$]: A2B3C4D_oP40_62_d_cd_2cd_c 632

$S0_8$ .....
1. Zunyite [$Al_{13}(OH,F)_{18}Si_5O_{20}Cl$, $S0_8$]: A13BC18D20E5_cF228_216_dh_b_fh_2eh_ce 1363

$S1_5$ .....
1. Eulytine ($Bi_4(SiO_4)_3$, $S1_5$): A4B12C3_cI76_220_c_e_a 1397

$S2_2$ .....
1. Hemimorphite ($Zn_4Si_2O_7(OH)_2{\cdot}H_2O$, $S2_2$): A2B5CD2_oI40_44_2c_abcde_d_e 514

$S2_3$ .....
1. Vesuvianite ($Ca_{10}Al_4(Mg,Fe)_2Si_9O_{34}(OH)_4$, $S2_3$): A4B10C2D34E4F9_tP252_126_k_ce2k_f_h8k_k_d2k 916

$S3_3(I)$ .....
1. Crancrinite ($Na_6Ca_2Al_6Si_6O_{24}(CO_3)_2$, $S3_3$ (I)): A3BCD3E15F3_hP52_173_c_b_b_c_5c_c 1168

$S3_3(II)$ .....
1. Parawollastonite ($CaSiO_3$, $S3_3$ (II)): AB3C_mP60_14_3e_9e_3e 260

$S3_4(I)$ .....
1. Chabazite ($Ca_{1.4}Sr_{0.3}Al_{3.8}Si_{8.3}O_{24}{\cdot}13H_2O$, $S3_4$ (I)): A5B21C24D12_hR62_166_a2c_ehi_fg2h_i 1118

$S3_4(II)$ .....
1. $S3_4$ (II) (Catapleiite, $Na_2Zr(SiO_3)_3{\cdot}H_2O$) (*obsolete*): A3B2C9D3E_hP36_194_g_f_hk_h_a 1254

$S4_1$ .....
1. Diopside [$CaMg(SiO_3)_2$, $S4_1$]: ABC6D2_mC40_15_e_e_3f_f 378

$S4_2$ .....
1. Tremolite ($Ca_2Mg_5Si_8O_{22}(OH)_2$, $S4_2$): A2B2C5D24E8_mC82_12_h_i_agh_2i5j_2j 148

$S4_3$ .....
1. Enstatite ($MgSiO_3$, $S4_3$): AB3C_oP80_61_2c_6c_2c 623

$S4_4$ .....
1. Anthophyllite ($Mg_5Fe_2Si_8O_{22}(OH)_2$, $S4_4$): A2B5C22D2E8_oP156_62_d_c2d_2c10d_2c_4d 638

$S4_5$ .....
1. Chrysotile ($H_4Mg_3Si_2O_9$, $S4_5$): AB6C11D6E4_mC112_12_e_gi2j_i5j_2i2j_2j 178

$S4_6$ .....
1. Bertrandite ($Be_4Si_2O_7(OH)_2$, $S4_6$): A4B7C2D2_oC60_36_2b_a3b_2a_b 473

$S4_7$ .....

[†]Ammonium Persulfate [$(NH_4)_2S_2O_8$, $K4_1$] and Monazite ($LaPO_4$) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

1. Epididymite ($BeHNaO_8Si_3$, $S4_7$): ABCD8E3_oP112_62_d_2c_d_4c6d_3d 728
$S5_1$ .....
1. Muscovite ($KH_2Al_3Si_3O_{12}$, $S5_1$): A2BC10D2E4_mC76_15_f_e_5f_f_2f 329
$S5_2$ .....
1. Apophyllite ($KCa_4Si_8O_{20}F{\cdot}8H_2O$, $S5_2$): A4BC16DE28F8_tP116_128_h_a_2i_b_g3i_i 937
$S5_4$ .....
1. Nacrite [$Al_2Si_2O_5(OH)_4$, $S5_4$]: A2B4C9D2_mC68_9_2a_4a_9a_2a 119
$S5_5$ .....
1. Chlorite [$Mg_3(Mg_2Al)(Si_3Al)O_{10}(OH)_8$, $S5_5$]: A3B5C4D2_mC112_15_a3ef_5f_4f_2f 347
$S5_6$ .....
1. Pyrophyllite [$AlSi_2O_5(OH)$, $S5_6$]: AB5CD2_mC72_15_f_5f_f_2f 371
$S5_7$ .....
1. Cronstedtite {$Fe(Fe,Si)[(OH)_2,O]O_3$, $S5_7$}: AB3C2D_hR7_160_a_b_2a_a 1070
$S6_1$ .....
1. Analcime ($NaAlSi_2O_6{\cdot}H_2O$, $S6_1$): A2B2C3D12E4_tI184_142_f_f_be_3g_g 1019
$S6_2$ .....
1. Sodalite [$Na_4(AlSiO_4)_3Cl$, $S6_2$]: A3BC4D12E3_cP46_218_d_a_e_i_c 1394
$S6_3$ .....
1. Danburite ($CaB_2Si_2O_8$, $S6_3$): A2BC8D2_oP52_62_d_c_2c3d_d 651
$S6_4$ .....
1. Marialite Scapolite [$Na_4Cl(AlSi_3)_3O_{24}$, $S6_4$]: AB4C24D12_tI82_87_a_h_2h2i_hi 825
$S6_5$ .....
1. α-Carnegieite ($NaAlSiO_4$, $S6_5$): ABC4D_cP28_198_a_a_ab_a 1280
$S6_6$ .....
1. Na_2CaSiO_4 ($S6_6$): AB2C4D_cP32_198_a_2a_ab_a 1277
$S6_7$ .....
1. Sanidine ($KAlSi_3O_8$, $S6_7$): AB8C4_mC52_12_i_gi3j_2j 186
$S6_8$ .....
1. Albite ($NaAlSi_3O_8$, $S6_8$): ABC8D3_aP26_2_i_i_8i_3i 68
$S6_9$ .....
1. Hauyne [$(Na_{0.5}Ca_{0.3}K_{0.2})_8(Al_6Si_6O_{24})(SO_4)_{1.5}$, $S6_9$]: A3B4C4D4E16F4G3_cP76_218_c_e_e_e_ei_e_d .. 1389
$S6_{10}$ .....
1. Natrolite ($Na_2Al_2Si_3O_{10}{\cdot}2H_2O$, $S6_{10}$): A2B4C2D12E3_oF184_43_b_2b_b_6b_ab 499
None .....
1. $NaC_5H_{11}O_8S$: A5B11CD8E_aP26_1_5a_11a_a_8a_a 35
2. $Ni(NO_3)_2(H_2O)_6$: A12B2CD12_aP54_2_12i_2i_i_12i 38
3. $Co_2B_2O_5$: A2B2C5_aP18_2_2i_2i_5i 42
4. α-$Ho_2Si_2O_7$: A2B7C2_aP44_2_4i_14i_4i 47
5. $Co_3(SeO_3)_3{\cdot}H_2O$: A3B2C10D3_aP36_2_ah2i_2i_10i_3i 51
6. δ-WO_3: A3B_aP32_2_12i_4i 54
7. Wollastonite ($CaSiO_3$): AB3C_aP30_2_3i_9i_3i 65
8. TaTi (BCC SQS-16): AB_aP16_2_4i_4i 71
9. $W_2O_3(PO_4)_2$: A11B2C2_mP60_4_22a_4a_4a 73
10. Ca_3UO_6: A3B6C_mP20_4_3a_6a_a 79
11. $NbAs_2$: A2B_mC12_5_2c_c 85
12. $Rb_2CaCu_6(PO_4)_4O_2$: AB6C18D4E2_mC62_5_a_2b2c_9c_2c_c 89
13. $C2$ $(Ba,Ca)CO_3$: ABC3_mC10_5_b_a_ac 92
14. Ta_5Ti_{11} (BCC SQS-16): A5B11_mP16_6_2abc_2a3b3c 94
15. $Na_2Ca_6Si_4O_{15}$: A6B2C15D4_mP54_7_6a_2a_15a_4a 96
16. Low-Temperature Mo_8O_{23}: A8B23_mP124_7_16a_46a 100
17. Calaverite ($AuTe_2$): AB2_mP12_7_2a_4a 107
18. Monoclinic Co_4Al_{13}: A13B4_mC102_8_17a11b_8a2b 109
19. $TaTi_3$ (BCC SQS-16): AB3_mC32_8_4a_12a 113
20. $TaTi_3$ (BCC SQS-16): AB3_mC32_8_4a_4a4b 115
21. Chrysotile ($Mg_3Si_2O_5(OH)_4$): A3B5C4D2_mC56_9_3a_5a_4a_2a 122
22. $Cs_6W_{11}O_{36}$: A6B36C11_mC212_9_6a_36a_11a ... 125
23. $ZrSe_3$: A3B_mP8_11_3e_e 133
24. y-$Y_2Si_2O_7$: A7B2C2_mP22_11_3e2f_2e_ab 135
25. Barytocalcite ($BaCa(CO_3)_2$): AB2CD6_mP20_11_e_2e_e_2e2f 137
26. O(OH)Y: ABC_mP6_11_e_e_e 139
27. $Al_{13}Fe_4$: A13B4_mC102_12_dg8i5j_4ij 141
28. Os_4Al_{13}: A13B4_mC34_12_b6i_2i 144
29. β-Ga_2O_3: A2B3_mC20_12_2i_3i 151
30. $K_2Ti_2O_5$: A2B5C2_mC18_12_i_a2i_i 153
31. CaC_2-III: A2B_mC12_12_2i_i 155
32. Tolbachite ($CuCl_2$): A2B_mC6_12_i_a 157
33. Staurolite ($Al_5Fe_2O_{10}(OH)_2Si_2$): A5B2C10D2E2_mC84_12_acghj_bdi_5j_2i_j 163
34. Monoclinic $FeTlSe_2$: AB2C_mC16_12_g_2i_i 171
35. $NbTe_2$: AB2_mC18_12_ai_3i 173
36. Sr_2NiTeO_6: AB6C2D_mC40_12_ad_gh4i_j_bc 182
37. Ta_2PdSe_6: AB6C2_mC18_12_a_3i_i 184
38. $MnPS_3$: ABC3_mC20_12_g_i_ij 189
39. $AlNbO_4$: ABC4_mC24_12_i_i_4i 191
40. SiAs: AB_mC24_12_3i_3i 193
41. High-Temperature Mo_8O_{23}: A8B23_mP62_13_4g_c11g 195

42. $Cs_{11}O_3$: A11B3_mP56_14_11e_3e 201
43. $HgCl_2{\cdot}2HgO$: A2B3C2_mP14_14_e_ae_e 208
44. Monoclinic Cu_2OSeO_3: A2B4C_mP28_14_abe_4e_e 213
45. $Sb_4O_5Cl_2$: A2B5C4_mP22_14_e_c2e_2e 216
46. Ca_2UO_5: A2B5C_mP32_14_2e_5e_ab 218
47. Gd_2SiO_5 (RE_2SiO_5 X1): A2B5C_mP32_14_2e_5e_e 221
48. Sanguite ($KCuCl_3$): A3BC_mP20_14_3e_e_e 224
49. γ-WO_3: A3B_mP32_14_6e_2e 226
50. $K_2Ni(CN)_4$: A4B2C4D_mP22_14_2e_e_2e_a 230
51. γ-$Y_2Si_2O_7$[‖]: A4BC_mP24_14_4e_e_e 235
52. $K_2Pt(SCN)_6{\cdot}2H_2O$: A6B4C2D6E2FG6_mP54_14_3e_2e_e_3e_e_a_3e 238
53. Manganese-leonite 110 K $[K_2Mn(SO_4)_2{\cdot}4H_2O]$: A8B2CD12E2_mP100_14_8e_2e_ad_12e_2e 246
54. β-B_2H_6[*]: AB3_mP16_14_e_3e 264
55. B_2H_6 ($P2_1/c$)[*]: AB3_mP16_14_e_3e 266
56. Anhydrous $KAuBr_4$: AB4C_mP24_14_ab_4e_e 271
57. Monazite ($LaPO_4$)[†‖]: AB4C_mP24_14_e_4e_e 277
58. Sr_2MnTeO_6: AB6C2D_mP20_14_a_3e_e_d 280
59. KNO_2 III[‡]: ABC2_mP16_14_e_e_2e 285
60. Cu(OH)Cl: ABCD_mP16_14_e_e_e_e 295
61. α-ICl[§]: AB_mP16_14_2e_2e 299
62. LiAs[§]: AB_mP16_14_2e_2e 301
63. ϵ-1,2,3,4,5,6-Hexachlorocyclohexane (C_6Cl_6): AB_mP24_14_3e_3e 303
64. Pararealgar (AsS)[¶]: AB_mP32_14_4e_4e 306
65. Ag_2PbO_2: A2B2C_mC20_15_ad_f_e 312
66. Catapleiite ($Na_2ZrSi_3O_9{\cdot}2H_2O$): A2B3C9D3E_mC144_15_2f_bcdef_9f_3f_ae 314
67. Na_2PrO_3: A2B3C_mC48_15_aef_3f_2e 319
68. Eudidymite ($BeHNaO_8Si_3$): A2B4C2D17E6_mC124_15_f_2f_f_e8f_3f 322
69. ζ-Nb_2O_5 (B-Nb_2O_5): A2B5_mC28_15_f_e2f 327
70. $Rb_2C_2O_4{\cdot}H_2O$: A2BC4D2_mC36_15_f_e_2f_f 332
71. Alluaudite $[NaMnFe_2(PO_4)_3]$: A2BCD12E3_mC76_15_f_e_b_6f_ef 334
72. ThC_2 (C_g): A2B_mC12_15_f_e 338
73. Clinocervantite (β-Sb_2O_4): A2B_mC24_15_2f_ce 340
74. Y_2SiO_5 (RE_2SiO_5 X2): A5BC2_mC64_15_5f_f_2f 351
75. α-$Zn_2V_2O_7$: A7B2C2_mC44_15_e3f_f_f 354
76. Manganese-leonite 185 K $[K_2Mn(SO_4)_2{\cdot}4H_2O]$: A8B2CD12E2_mC200_15_8f_2f_ce_2e11f_2f 357
77. Ta_2NiSe_5: AB5C2_mC32_15_e_e2f_f 369
78. β-Ga (*obsolete*): A_mC4_15_e 381
79. $NaNbO_3$: ABC3_oP40_17_abcd_2e_abcd4e 383
80. γ-TeO_2: A2B_oP12_18_2c_c 386
81. Diamminetriamidodizinc Chloride $([Zn_2(NH_3)_2(NH_2)_3]Cl)$: AB12C5D2_oP40_18_a_6c_b2c_c 388
82. Ferroelectric $NH_4H_2PO_4$: A6BC4D_oP48_19_6a_a_4a_a 398
83. β-Arabinose $[(CH_2O)_{20}]$: AB2C_oP80_19_5a_10a_5a 401
84. $NaAlCl_4$: AB4C_oP24_19_a_4a_a 405
85. NaP: AB_oP16_19_2a_2a 407
86. $AlPO_4$ "low cristobalite type": AB4C_oC24_20_b_2c_a 411
87. $HoSb_2$: AB2_oC6_21_a_k 415
88. Predicted Phase IV $Cd_2Re_2O_7$: A2B7C2_oF88_22_k_bdefghij_k 417
89. Mercury (II) Azide $[Hg(N_3)_2]$: AB6_oP28_29_a_6a 420
90. Low-Temperature $(NH_3CH_3)Al(SO_4)_2{\cdot}12H_2O$: ABC30DE20F2_oP220_29_a_a_30a_a_20a_2a 422
91. Orthorhombic Co_4Al_{13}: A13B4_oP102_31_17a11b_8a2b 430
92. B_4SrO_7: A4B7C_oP24_31_2b_a3b_a 438
93. $Mo_{17}O_{47}$: A17B47_oP128_32_a8c_a23c 442
94. Possible δ-$Gd_2Si_2O_7$: A2B7C2_oP44_33_2a_7a_2a 447
95. CaB_2O_4 (III): A2BC4_oP84_33_6a_3a_12a 450
96. Cervantite (α-Sb_2O_4): A2B_oP24_33_4a_2a 455
97. CsB_4O_6F: A4BCD6_oP48_33_4a_a_a_6a 457
98. $LiGaO_2$: ABC2_oP16_33_a_a_2a 460
99. γ-$LiIO_3$: ABC3_oP20_33_a_a_3a 462
100. $MnF_{2-x}(OH)_x$: A2BC2_oP10_34_c_a_c 464
101. Si_2N_2O: A2BC2_oC20_36_b_a_b 466
102. Bi_2GeO_5: A2BC5_oC32_36_b_a_a2b 468
103. Ni_3Si_2: A3B2_oC80_36_4a4b_2a3b 470
104. MoP_2: AB2_oC12_36_a_2a 476
105. α-Potassium Nitrate (KNO_3) II: ABC3_oC80_36_2ab_2ab_2a5b 478
106. Ta_3Ti_{13} (BCC SQS-16): A3B13_oC32_38_ac_a2bcdef 481
107. Ta_3Ti_5 (BCC SQS-16): A3B5_oC32_38_abce_abcdf 483
108. $NaNb_6O_{15}F$: ABC6D15_oC46_38_b_b_2a2d_2ab4d2e 485
109. $Rb_2Mo_2O_7$: A2B7C2_oC88_40_abc_2b6c_a3b 488
110. Orthorhombic CrO_3: AB3_oC16_40_b_a2b 491

[*] β-B_2H_6 and B_2H_6 ($P2_1/c$) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[§] α-ICl and LiAs have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

111. Ag_2O_3[**]: A2B3_oF40_43_b_ab 497
112. Blossite (α-$Cu_2V_2O_7$): A2B7C2_oF88_43_b_a3b_b 504
113. Archerite (KH_2PO_4): A2BC4D_oF64_43_b_a_2b_a 507
114. Cs_2Se: A2B_oF24_43_b_a 510
115. Zr_2Al_3[**]: A3B2_oF40_43_ab_b 512
116. $Nb_2Zr_6O_{17}$: A2B17C6_oI100_46_ab_b8c_3c 523
117. Parkerite ($Ni_3Bi_2S_2$): AB2C_oP8_51_e_be_f 527
118. $LiNb_6O_{15}F$: ABC6D15_oP46_51_f_d_2e2i_aef4i2j 529
119. Carnallite [$Mg(H_2O)_6KCl_3$]: A3B12CDE6_oP276_52_d4e_18e_ce_de_2d8e 532
120. Orthorhombic $Sr_4Ru_3O_{10}$: A10B3C4_oP68_55_2e2fgh2i_adef_2e2f 548
121. $Nb_2Pd_3Se_8$: A2B3C8_oP26_55_h_ag_2g2h 551
122. $Ru_{11}B_8$: A8B11_oP38_55_g3h_a3g2h 556
123. $HoMn_2O_5$: AB2C5_oP32_55_g_fh_eghi 559
124. Calciborite (CaB_2O_4 II): A2BC4_oP56_56_2e_e_4e 562
125. $SrUO_4$: A4BC_oP24_57_cde_d_a 569
126. Lueshite ($NaNbO_3$): ABC3_oP40_57_cd_e_cd2e 571
127. Kotoite ($Mg_3(BO_3)_2$): A2B3C6_oP22_58_g_af_gh 574
128. Protoanthophyllite ($H_2Mg_7Si_8O_{24}$): A2B7C24D8_oP82_58_g_ae2f_2g5h_2h 579
129. In_4Se_3: A4B3_oP28_58_4g_3g 584
130. InS: AB_oP8_58_g_g 589
131. RuB_2: A2B_oP6_59_f_a 591
132. Shcherbinaite (V_2O_5) (*revised*): A5B2_oP14_59_a2f_f 595
133. ζ-Fe_2N[×]: A2B_oP12_60_d_c 600
134. α-PbO_2[×]: A2B_oP12_60_d_c 602
135. Cr_5O_{12}: A5B12_oP68_60_c2d_6d 604
136. Ca_2RuO_4: A2B4C_oP28_61_c_2c_a 611
137. $(TiCl_4{\cdot}POCl_3)_2$: A7BCD_oP80_61_7c_c_c_c 616
138. COCl: ABC_oP24_61_c_c_c 627
139. Cs_2Sb: A2B_oP24_62_4c_2c 656
140. Original β-WO_3 (*obsolete*): A3B_oP32_62_ab4c_2c 665
141. P_4Se_3: A4B3_oP112_62_8c4d_4c4d 668
142. Mo_4P_3: A4B3_oP56_62_8c_6c 673
143. $K_2SnCl_4{\cdot}H_2O$: A4BC2D_oP32_62_2cd_c_d_c 679
144. $VOSO_4$: A5BC_oP28_62_3cd_c_c 682
145. Possible δ-$Y_2Si_2O_7$: A7B2C2_oP44_62_3c2d_2c_d 684
146. $SbCl_5{\cdot}POCl_3$: A8BCD_oP44_62_4c2d_c_c_c 692
147. Autunite {$Ca[(UO_2)(PO_4)]_2(H_2O)_{11}$}: AB22C23D2E2_oP200_62_c_11d_3c10d_d_d 695
148. Atacamite ($Cu_2(OH)_3Cl$): AB2C3D3_oP36_62_c_ac_cd_cd 704
149. Chalcocyanite ($CuSO_4$): AB4C_oP24_62_a_2cd_c 711
150. Rynersonite (Orthorhombic $CaTa_2O_6$): AB6C2_oP36_62_c_2c2d_d 713
151. Copper (II) Azide [$Cu(N_3)_2$]: AB6_oP28_62_c_6c 716
152. α-Potassium Nitrate (KNO_3) I[°]: ABC3_oP20_62_c_c_cd 720
153. MnCuP: ABC_oP12_62_c_c_c 736
154. $Cu_2Pb(SeO_3)_2Br_2$: A2B2C6DE2_oC52_63_g_e_fh_c_f 740
155. Pd_5Pu_3: A5B3_oC32_63_cfg_ce 749
156. $ZrTe_5$: A5B_oC24_63_c2f_c 751
157. $ThFe_2SiC$: AB2CD_oC20_63_b_f_c_c 757
158. Re_3B: AB3_oC16_63_c_cf 761
159. Ta_2NiS_5[††]: AB5C2_oC32_63_c_c2f_f 763
160. V_3AsC: ABC3_oC20_63_c_b_cf 765
161. Si_{24} Clathrate: A_oC24_63_3f 767
162. Base-centered orthorhombic $Sr_4Ru_3O_{10}$: A10B3C4_oC68_64_2dfg_ad_2d 769
163. $Na_2Mo_2O_7$: A2B2C7_oC88_64_ef_df_3f2g 772
164. Li_2PrO_3: A2B3C_oC12_65_h_bh_a 775
165. Nb_3O_7F: A3B8_oC22_65_ag_bd2gh 779
166. High-Temperature Cryolite (Na_3AlF_6): AB6C3_oI20_71_a_in_cj 787
167. $CsFeS_2$ (100 K): ABC2_oI16_71_g_i_eh 789
168. CsO: AB_oI8_71_g_i 791
169. $CeCu_2$: AB2_oI12_74_e_h 797
170. $LiCuVO_4$: ABC4D_oI28_74_a_d_hi_e 799
171. Gwihabaite [NH_4NO_3 (V)]: A4B2C3_tP72_77_8d_ab2c2d_6d 801
172. Kesterite [$Cu_2(Zn,Fe)SnS_4$]: A2BCD4_tI16_82_ac_b_d_g 805
173. $MoPO_5$: AB5C_tP14_85_c_cg_b 810
174. $Nd_4Re_2O_{11}$: A4B11C2_tP68_86_2g_ab5g_g 815
175. β-$LiIO_3$: ABC3_tP40_86_g_g_3g 822
176. Sr_2NiWO_6: AB6C2D_tI20_87_a_eh_d_b 828
177. $Na_4Ge_9O_{20}$: A9B4C20_tI132_88_a2f_f_5f 830
178. Copper (I) Azide (CuN_3): AB3_tI32_88_d_cf 835
179. Paratellurite (α-TeO_2): A2B_tP12_92_b_a 846
180. Phase III $Cd_2Re_2O_7$: A2B7C2_tI44_98_f_bcde_f 848
181. VSe_2O_6: A6B2C_tP72_103_abc5d_2d_abc 855

[**] Ag_2O_3 and Zr_2Al_3 have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

[×] ζ-Fe_2N and α-PbO_2 have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

182. $BaNiSn_3$: ABC3_tI10_107_a_a_ab 859
183. Ammonium Chlorite (NH_4ClO_2): AB4CD2_tP16_113_c_f_a_e 863
184. $C_{19}Sc_{15}$: A19B15_tP68_114_bc4e_ac3e 865
185. Phase II $Cd_2Re_2O_7$: A2B7C2_tI44_119_i_bdefgh_i 874
186. Tetragonal $TlFeS_2$: AB2C_tI8_119_c_e_a 877
187. $SrCu_2(BO_3)_2$: A2B2C6D_tI44_121_i_i_ij_c 882
188. K_3CrO_8: AB3C8_tI24_121_a_bd_2i 886
189. α-V_3S: AB3_tI32_121_g_f2i 888
190. $NH_4H_2PO_4$: A8BC4D_tI56_122_2e_b_e_a 896
191. NaS_2: AB2_tI48_122_cd_2e 899
192. $CaO_2(H_2O)_8$: AB8C2_tP22_124_a_n_h 914
193. Phosgenite [$Pb_2Cl_2(CO_3)$]: AB2C3D2_tP32_127_g_eh_gk_k 931
194. $CaBe_2Ge_2$: A2BC2_tP10_129_ac_c_bc 943
195. LaOAgS: ABCD_tP8_129_b_c_a_c 950
196. $Sr(OH)_2(H_2O)_8$: A18B10C_tP116_130_2c4g_2c2g_a 952
197. α-WO_3: A3B_tP16_130_cf_c 957
198. Zr_3Al_2: A2B3_tP20_136_j_dfg 960
199. $ZrFe_4Si_2$: A4B2C_tP14_136_i_g_b 962
200. $Nd_2Fe_{14}B$: AB14C2_tP68_136_f_ce2j2k_fg 966
201. $Sr_4Ti_3O_{10}$: A10B4C3_tI34_139_c2eg_2e_ae 970
202. $TlCo_2S_2$[∂‡‡]: A2B2C_tI10_139_d_e_a 972
203. $ThCr_2Si_2$[∂‡‡]: A2B2C_tI10_139_d_e_a 974
204. Au_2Nb_3: A2B3_tI10_139_e_ae 976
205. K_2NiF_4: A4B2C_tI14_139_ce_e_a 982
206. $Sr_3Ti_2O_7$: A7B3C2_tI24_139_aeg_be_e 988
207. Li_2CN_2[⊗‡‡]: AB2C2_tI10_139_a_d_e 992
208. V_4SiSb_2: A2BC4_tI28_140_h_a_k 1000
209. $Pu_{31}Rh_{20}$: A31B20_tI204_140_b2gh3m_ac2fh3l .. 1004
210. $BaCd_{11}$: AB11_tI48_141_a_bdi 1016
211. Cd_3As_2: A2B3_tI160_142_deg_3g 1025
212. La_3BWO_9 ($P3$): AB3C9D_hP28_143_2a_2d_6d_bc 1030
213. $RbNO_3$ (IV): AB3C_hP45_144_3a_9a_3a 1033
214. Li_7TaO_6: A8B6C_hR15_148_cf_f_a 1047
215. $SrCl_2{\cdot}(H_2O)_6$: A2B12C6D_hP21_150_d_2g_ef_a 1049
216. Paralstonite ($BaCa(CO_3)_2$): AB2CD6_hP30_150_e_c2d_f_3g 1054
217. $KBe_2BO_3F_2$: AB2C2DE3_hR9_155_b_c_c_a_e .. 1063
218. SbI_3S_{24}: A3B24C_hR28_160_b_2b3c_a 1065
219. Fe_3PO_7: A3B7C_hR11_160_b_a2b_a 1068
220. Low-Temperature $GaMo_4S_8$: AB4C8_hR13_160_a_ab_2a2b 1072
221. γ-Potassium Nitrate (KNO_3)[□]: ABC3_hR5_160_a_a_b 1076
222. α-BaB_2O_4 (Low-Temperature): A2BC4_hR42_161_2b_b_4b 1078
223. Rosiaite ($PbSb_2O_6$)[§§]: A6BC2_hP9_162_k_a_d .. 1085
224. Colquiriite ($LiCaAlF_6$): ABC6D_hP18_163_d_b_i_c 1090
225. Predicted Li_2MgH_{16} 300 GPa: A16B2C_hP19_164_2d2i_d_b 1093
226. Ce_2O_2S[¶¶]: A2B2C_hP5_164_d_d_a 1095
227. Brucite [$Mg(OH)_2$][¶¶]: A2BC2_hP5_164_d_a_d .. 1097
228. Jacutingaite (Pt_2HgSe_3): AB2C3_hP12_164_d_ae_i 1105
229. Nevskite (BiSe): AB_hP12_164_c2d_c2d 1109
230. $MnBi_2Te_4$[°°]: A2BC4_hR7_166_c_a_2c 1114
231. Shandite ($Ni_3Pb_2S_2$): A3B2C2_hR7_166_d_ab_c 1116
232. Rhombohedral $CuTi_2S_4$: AB4C2_hR28_166_2c_2c2h_abh 1126
233. $CaCu_4P_2$[°°]: AB4C2_hR7_166_a_2c_c 1129
234. $CaUO_4$: AB4C_hR6_166_b_2c_a 1131
235. $TaTi_7$ (BCC SQS-16): AB7_hR16_166_c_c2h ... 1133
236. Rhombohedral Delafossite ($CuFeO_2$): ABC2_hR4_166_a_b_c 1136
237. β-Potassium Nitrate (KNO_3): ABC6_hR8_166_a_b_h 1138
238. $Zr_{21}Re_{25}$: A25B21_hR92_167_b2e3f_e3f 1142
239. β-BaB_2O_4 (High-Temperature): A2BC4_hR42_167_f_ac_2f 1151
240. Rinneite ($K_3NaFeCl_6$): A6BC3D_hR22_167_f_b_e_a 1161
241. La_3CuSiS_7: AB3C7D_hP24_173_a_c_b2c_b 1172
242. La_3BWO_9 ($P6_3$): AB3C9D_hP28_173_a_c_3c_b . 1175
243. α-$LiIO_3$: ABC3_hP10_173_b_a_c 1178
244. $Rh_{20}Si_{13}$: A10B7_hP34_176_c3h_b2h 1182
245. β-Si_3N_4: A4B3_hP14_176_ch_h 1187
246. Hg_2O_2NaI: A2BCD2_hP18_180_f_c_b_i 1198
247. $Zn_2Mo_3O_8$: A3B8C2_hP26_186_c_ab2c_2b 1204
248. Cr-233 Quasi-One-Dimensional Superconductor ($K_2Cr_3As_3$): A3B3C2_hP16_187_jk_jk_ck 1220
249. Cs_7O: A7B_hP24_187_ai2j2kn_j 1222
250. ZrNiAl: ABC_hP9_189_g_ad_f 1225

[∂] $TlCo_2S_2$ and $ThCr_2Si_2$ have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[¶¶] Ce_2O_2S and Brucite [$Mg(OH)_2$] have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

[°°] $MnBi_2Te_4$ and $CaCu_4P_2$ have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

251. Hexagonal WO_3: A3B_hP12_191_gl_f 1235
252. Ti_5Ga_4: A4B5_hP18_193_bg_dg 1241
253. Proposed 300 GPa HfH_{10}: A10B_hP22_194_bhj_c 1243
254. Magnetoplumbite ($PbFe_{12}O_{19}$): A12B19C_hP64_194_ab2fk_efh2k_d 1246
255. ReB_3: A3B_hP8_194_af_c 1257
256. $Cs_3Cr_2Cl_9$: A9B2C3_hP28_194_hk_f_bf 1259
257. $Na_{0.74}CoO_2$: AB2C2_hP10_194_a_bc_f 1261
258. $EuIn_2P_2$: AB2C2_hP10_194_a_f_f 1263
259. Lu_2CoGa_3: AB3C2_hP24_194_f_k_bh 1265
260. Hexagonal Delafossite ($CuAlO_2$): ABC2_hP8_194_a_c_f 1268
261. Cubic Cu_2OSeO_3: A2B4C_cP56_198_ab_2a2b_2a 1274
262. $Bi_3Ru_3O_{11}$: A3B11C3_cP68_201_be_efh_g 1285
263. $LaFe_4P_{12}$: A4BC12_cI34_204_c_a_g 1292
264. $NaMn_7O_{12}$: A7BC12_cI40_204_bc_a_g 1294
265. CaB_2O_4 (IV): A2BC4_cP84_205_d_ac_2d 1309
266. $NaSbF_6$: A6BC_cP32_205_d_b_a 1316
267. $NaCr(SO_4)_2{\cdot}12H_2O$ Alum: AB12CD8E2_cP96_205_a_2d_b_cd_c 1322
268. Al_2Mo_3C: A2BC3_cP24_213_c_a_d 1358
269. Mg_3Ru_2: A3B2_cP20_213_d_c 1361
270. Murataite $[(Y,Na)_6(Zn,Fe)_5Ti_{12}O_{29}(O,F)_{10}F_4]$: A16B40C12D6E5_cF316_216_eh_e2g2h_h_f_be .. 1367
271. $Sm_{11}Cd_{45}$: A45B11_cF448_216_bd4efg5h_ac2eh . 1371
272. Hg_2TiCu Inverse Heusler: AB2C_cF16_216_b_ad_c 1376
273. $GaMo_4S_8$: AB4C8_cF52_216_a_e_2e 1378
274. AlN (cF40): AB_cF40_216_de_ce 1382
275. Tennantite ($Cu_{12}As_4S_{13}$): A4B24C13_cI82_217_c_deg_ag 1384
276. AlN (cI16): AB_cI16_217_c_c 1387
277. AlN (cI24): AB_cI24_220_a_b 1412
278. Predicted High-Pressure $YCaH_{12}$: AB12C_cP14_221_a_h_b 1416
279. Dodecatungstophosphoric Acid Hexahydrate $[H_3PW_{12}O_{40}{\cdot}6H_2O]$: A27B52CD12_cP184_224_dl_eh3k_a_k 1420
280. $H_3PW_{12}O_{40}{\cdot}3H_2O$: A3B40CD12_cP112_224_d_e3k_a_k 1429
281. LaH_{10} High-T_c Superconductor: A10B_cF44_225_cf_b 1439
282. Double Perovskite (Ba_2MnWO_6): A2BC6D_cF40_225_c_a_e_b 1441
283. γ-Ga_2O_3: A11B4_cF120_227_acdf_e 1455
284. Predicted Li_2MgH_{16} High-Temperature Superconductor (250 GPa): A16B2C_cF152_227_eg_d_a 1458
285. $Mg_3Cr_2Al_{18}$: A18B2C3_cF184_227_fg_d_ac 1461
286. $Zn_{22}Zr$: A22B_cF184_227_cdfg_a 1464

Duplicate AFLOW Label Index

AB3_mP16_14_e_3e .....
1. β-B_2H_6 264
2. B_2H_6 ($P2_1/c$) 266

AB4C_mP24_14_e_4e_e .....
1. Ammonium Persulfate $[(NH_4)_2S_2O_8, K4_1]$ 274
2. Monazite ($LaPO_4$) 277

ABC2_mP16_14_e_e_2e .....
1. KNO_2 III 285
2. Manganite (γ-MnO(OH), $E0_6$) 287

AB_mP16_14_2e_2e .....
1. α-ICl 299
2. LiAs 301

AB_mP32_14_4e_4e .....
1. Pararealgar (AsS) 306
2. Realgar (AsS, B_l) 309

A2B_oP12_60_d_c .....
1. ζ-Fe_2N 600
2. α-PbO_2 602

ABC3_oP20_62_c_c_cd .....
1. α-Potassium Nitrate (KNO_3) I 720
2. NH_4NO_3 III ($G0_{10}$) 723
3. Aragonite ($CaCO_3$, $G0_2$) 726

A2B2C_tI10_139_d_e_a .....
1. $TlCo_2S_2$ 972
2. $ThCr_2Si_2$ 974

AB2C2_tI10_139_a_d_e .....
1. Li_2CN_2 992
2. $H5_9$ [Autunite, $Ca(UO_2)_2(PO_4)_2{\cdot}10\frac{1}{2}H_2O$] (*obsolete*) 994

ABC3_hR5_160_a_a_b .....
1. $KBrO_3$ ($G0_7$) 1074
2. γ-Potassium Nitrate (KNO_3) 1076

Similar AFLOW Label Index

A4BC_mP24_14_4e_e_e .....
1. γ-$Y_2Si_2O_7$: A4BC_mP24_14_4e_e_e 235
2. Ammonium Persulfate $[(NH_4)_2S_2O_8, K4_1]$: AB4C_mP24_14_e_4e_e 274
3. Monazite ($LaPO_4$): AB4C_mP24_14_e_4e_e 277
4. $AgMnO_4$ ($H0_9$): ABC4_mP24_14_e_e_4e 289

A2B3_oF40_43_b_ab .....
1. Ag_2O_3: A2B3_oF40_43_b_ab 497
2. Zr_2Al_3: A3B2_oF40_43_ab_b 512

A2B5C_oC32_63_f_c2f_c .....
1. Pseudobrookite (Fe_2TiO_5, $E4_1$): A2B5C_oC32_63_f_c2f_c 742
2. Ta_2NiS_5: AB5C2_oC32_63_c_c2f_f 763

A2B2C_tI10_139_d_e_a.TlCo2S2 .....
1. $TlCo_2S_2$: A2B2C_tI10_139_d_e_a 972
2. $ThCr_2Si_2$: A2B2C_tI10_139_d_e_a 974
3. Li_2CN_2: AB2C2_tI10_139_a_d_e 992

4. $H5_9$ [Autunite, $Ca(UO_2)_2(PO_4)_2{\cdot}10\frac{1}{2}H_2O$] (*obsolete*):
AB2C2_tI10_139_a_d_e 994
A2B6C_hP9_162_d_k_a .....
1. $I1_3$ ($SrCl_2{\cdot}(H_2O)_6$) (*obsolete*):
A2B6C_hP9_162_d_k_a 1083
2. Rosiaite ($PbSb_2O_6$): A6BC2_hP9_162_k_a_d 1085
A2B2C_hP5_164_d_d_a .....
1. Ce_2O_2S: A2B2C_hP5_164_d_d_a 1095
2. Brucite [$Mg(OH)_2$]: A2BC2_hP5_164_d_a_d 1097
A2BC6_hP9_164_d_a_i .....
1. $K_2Pt(SCN)_6$ ($H6_3$): A2BC6_hP9_164_d_a_i 1099
2. Bararite (Trigonal $(NH_4)_2SiF_6$, $J1_6$):
A6B2C_hP9_164_i_d_a 1101
3. K_2GeF_6 ($J1_{13}$): A6BC2_hP9_164_i_a_d 1103
A2BC4_hR7_166_c_a_2c .....
1. $MnBi_2Te_4$: A2BC4_hR7_166_c_a_2c 1114
2. $CaCu_4P_2$: AB4C2_hR7_166_a_2c_c 1129
AB_hP4_194_a_c .....
1. $LiZn_2$ (C_k): AB_hP4_194_a_c 1270
2. Fe_2N (approximate, $L'3_0$): AB_hP4_194_c_a 1272
A2B6C6D_cP60_205_c_d_d_a .....
1. $Zn(BrO_3)_2{\cdot}6H_2O$ ($J1_{10}$):
A2B6C6D_cP60_205_c_d_d_a 1299
2. $H6_4$ [$Ni(NO_3)_2(NH_3)_6$] (*obsolete*):
A2B6CD6_cP60_205_c_d_a_d 1302

CIF Index

1. $B30$ (MgZn?): AB_oI48_44_6d_ab2cde 1605
2. $C17$ (Fe_2B) (*obsolete*):
AB2_tI12_121_ab_i 1685
3. $C2$ (Ba,Ca)CO_3: ABC3_mC10_5_b_a_ac 1511
4. $C26_a$ (NO_2) (*obsolete*):
AB2_cI36_199_b_c 1773
5. $C27$ (CdI_2) (*questionable*):
AB2_hP6_186_b_ab 1757
6. $C53$ ($SrBr_2$) (*obsolete*):
A2B_oP12_62_2c_c 1633
7. $D0_{10}$ (WO_3) (*obsolete*):
A3B_oP16_57_a2d_d 1614
8. $D0_{13}$ ($AlCl_3$) (*obsolete*):
AB3_hP4_164_b_ad 1735
9. $D0_{15}$ ($AlCl_3$) (*obsolete*):
AB3_mC16_5_c_3c 1510
10. $D0_6$ (Tysonite, LaF_3) (*obsolete*):
A3B_hP24_193_ack_g 1762
11. $D0_7$ (CrO_3) (*obsolete*):
AB3_oC16_20_a_bc 1579
12. $D2_2$ ($MgZn_5$?) (*problematic*):
AB5_mC48_12_2i_ac5i2j 1531
13. $D6_2$ (Sb_2O_4) (*obsolete*):
A2B_cF96_227_abf_cd 1830
14. $D8_7$ (Shcherbinaite, V_2O_5) (*obsolete*):
A5B2_oP14_31_a2b_b 1587
15. $E2_3$ ($LiIO_3$) (*obsolete*):
ABC3_hP10_182_c_b_g 1754
16. $E3_1$ (β-Ag_2HgI_4) (*obsolete*):
A2BC4_tP7_111_f_a_n 1680
17. $E6_1$ ($Sr(OH)_2(H_2O)_8$) (*obsolete*):
A8B2C_tP11_123_r_f_a 1690
18. $E6_2$ [$SrO_2(H_2O)_8$] (*possibly obsolete*):
A8B2C_tP11_123_r_h_a 1690
19. $F5_{11}$ (KNO_2) (*obsolete*):
ABC2_mC8_8_a_a_b 1517
20. $F5_4$ (NH_4ClO_2) (*obsolete*):
ABC2_tP8_100_b_a_c 1678
21. $F6_1$ (Chalcopyrite, $CuFeS_2$) (*obsolete*):
ABC2_tP4_115_a_c_g 1682
22. $G7_3$ [Northupite, $Na_3MgCl(CO_3)_2$] (*obsolete*):
A2BCD3E6_cF208_227_e_c_d_f_g 1828
23. $G7_5$ ($PbCO_3 \cdot PbCl_2$, Phosgenite) (*obsolete*):
AB2C3D2_tP16_90_c_f_ce_e 1675
24. $H5_6$ [Tychite, $Na_6Mg_2SO_4(CO_3)_4$)] (*obsolete*):
A4B2C6D16E_cF232_227_e_d_f_eg_a 1833
25. $H5_9$ [Autunite, $Ca(UO_2)_2(PO_4)_2{\cdot}10\frac{1}{2}H_2O$] (*obsolete*)[⊗‡‡]:
AB2C2_tI10_139_a_d_e 1710
26. $H6_4$ $Ni(NO_3)_2(NH_3)_6$] (*obsolete*)[⊗⊗]:
A2B6CD6_cP60_205_c_d_a_d 1778
27. $I1_3$ ($SrCl_2{\cdot}(H_2O)_6$) (*obsolete*)[§§]:
A2B6C_hP9_162_d_k_a 1729
28. $L1_a$ (disputed $CuPt_3$): AB7_cF32_225_b_ad 1818
29. $S0_4$ (Staurolite, $Fe(OH)_2Al_4Si_2O_{10}$) (*obsolete*):
A4BC12D2_oC76_63_eg_c_f3gh_g 1653
30. $S3_4$ (II) (Catapleiite, $Na_2Zr(SiO_3)_3{\cdot}\ H_2O$) (*obsolete*):
A3B2C9D3E_hP36_194_g_f_hk_h_a 1766
31. α-AgI ($B23$): A21B_cI44_229_bdh_a 1835
32. α-Alum [$KAl(SO_4)_2 \cdot 12H_2O$, $H4_{13}$]:
AB24CD28E2_cP224_205_a_4d_b_2c4d_c 1785

[⊗] Li_2CN_2 and $H5_9$ [Autunite, $Ca(UO_2)_2(PO_4)_2{\cdot}10\frac{1}{2}H_2O$] (*obsolete*) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[‡‡] $TlCo_2S_2$, $ThCr_2Si_2$, Li_2CN_2, and $H5_9$ [Autunite, $Ca(UO_2)_2(PO_4)_2{\cdot}10\frac{1}{2}H_2O$] (*obsolete*) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

[⊗⊗] $Zn(BrO_3)_2{\cdot}6H_2O$ ($J1_{10}$) and $H6_4$ [$Ni(NO_3)_2(NH_3)_6$] (*obsolete*) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

[§§] $I1_3$ ($SrCl_2{\cdot}(H_2O)_6$) (*obsolete*) and Rosiaite ($PbSb_2O_6$) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

33. α-BaB_2O_4 (Low-Temperature): A2BC4_hR42_161_2b_b_4b 1729
34. α-Carnegieite ($NaAlSiO_4$, $S6_5$): ABC4D_cP28_198_a_a_ab_a 1772
35. α-$Ho_2Si_2O_7$: A2B7C2_aP44_2_4i_14i_4i 1501
36. α-ICl[§]: AB_mP16_14_2e_2e 1556
37. α-$LiIO_3$: ABC3_hP10_173_b_a_c 1749
38. α-PbO_2[×]: A2B_oP12_60_d_c 1622
39. α-Potassium Nitrate (KNO_3) I[°]: ABC3_oP20_62_c_c_cd 1648
40. α-Potassium Nitrate (KNO_3) II: ABC3_oC80_36_2ab_2ab_2a5b 1596
41. α-V_3S: AB3_tI32_121_g_f2i 1686
42. α-WO_3: A3B_tP16_130_cf_c 1701
43. α-$Zn_2V_2O_7$: A7B2C2_mC44_15_e3f_f_f 1567
44. β-Alum [$Al(NH_3CH_3)_2(SO_4)_2{\cdot}12H_2O$, $H4_{14}$]: AB2C36D2E20F2_cP252_205_a_c_6d_c_c3d_c ... 1787
45. β-Alumina (Al_2O_3, $D5_6$): A2B3_hP60_194_3fk_cdef2k 1765
46. β-Arabinose [$(CH_2O)_{20}$]: AB2C_oP80_19_5a_10a_5a 1577
47. β-B_2H_6[*]: AB3_mP16_14_e_3e 1549
48. β-BaB_2O_4 (High-Temperature): A2BC4_hR42_167_f_ac_2f 1744
49. β-Ga (*obsolete*): A_mC4_15_e 1573
50. β-Ga_2O_3: A2B3_mC20_12_2i_3i 1526
51. β-$LiIO_3$: ABC3_tP40_86_g_g_3g 1672
52. β-Potassium Nitrate (KNO_3): ABC6_hR8_166_a_b_h 1742
53. β-Si_3N_4: A4B3_hP14_176_ch_h 1751
54. δ-CuTi ($L2_a$): AB_tP2_123_a_d 1691
55. δ-Ni_3Sn_4 ($D7_a$): A3B4_mC14_12_ai_2i 1527
56. δ-WO_3: A3B_aP32_2_12i_4i 1503
57. ϵ-1,2,3,4,5,6-Hexachlorocyclohexane (C_6Cl_6): AB_mP24_14_3e_3e 1557
58. η-NiSi (B_d): AB_oP8_62_c_c 1651
59. γ-Alum [$AlNa(SO_4)_2{\cdot}12H_2O$, $H4_{15}$]: AB24CD20E2_cP192_205_a_4d_b_c3d_c 1783
60. γ-Fe_4N ($L'1_0$): A4B_cP5_221_bc_a 1804
61. γ-Ga_2O_3: A11B4_cF120_227_acdf_e 1820
62. γ-$LiIO_3$: ABC3_oP20_33_a_a_3a 1592
63. γ-Potassium Nitrate (KNO_3)[□]: ABC3_hR5_160_a_a_b 1728
64. γ-TeO_2: A2B_oP12_18_2c_c 1574
65. γ-WO_3: A3B_mP32_14_6e_2e 1542
66. γ-$Y_2Si_2O_7$[‖]: A4BC_mP24_14_4e_e_e 1543
67. ζ-Fe_2N[×]: A2B_oP12_60_d_c 1621
68. ζ-Nb_2O_5 (B-Nb_2O_5): A2B5_mC28_15_f_e2f 1561
69. y-$Y_2Si_2O_7$: A7B2C2_mP22_11_3e2f_2e_ab 1521
70. $(CdSO_4)_3{\cdot}8H_2O$ ($H4_{20}$): A3B16C20D3_mC168_15_ef_8f_10f_ef 1565
71. $(NH_4)_3AlF_6$ ($J2_1$): AB30C16D3_cF200_225_a_ej_2f_bc 1816
72. $(TiCl_4{\cdot}POCl_3)_2$: A7BCD_oP80_61_7c_c_c_c 1625
73. 12-phosphotungstic acid [$H_3PW_{12}O_{40}{\cdot}5H_2O$ ($H4_{16}$)]: A5B40CD12_cP116_224_cd_e3k_a_k 1810
74. Adamite [$Zn_2(AsO_4)(OH)$, $H2_7$]: ABC5D2_oP36_58_g_g_3gh_eg 1618
75. Ag_2O_3[**]: A2B3_oF40_43_b_ab 1600
76. Ag_2PbO_2: A2B2C_mC20_15_ad_f_e 1559
77. $Ag_2SO_4{\cdot}4NH_3$ ($H4_{17}$): A2B12C4D4E_tP46_114_d_3e_e_e_a 1682
78. $AgMnO_4$ ($H0_9$)[‖]: ABC4_mP24_14_e_e_4e 1554
79. $AgNO_2$ ($F5_{12}$): ABC2_oI8_44_a_a_d 1604
80. $Ag[Co(NH_3)_2(NO_2)_4]$ ($J1_9$): ABC4D2E8_tP32_126_a_b_h_e_k 1694
81. $Al_{13}Fe_4$: A13B4_mC102_12_dg8i5j_4ij 1523
82. Al_2Mo_3C: A2BC3_cP24_213_c_a_d 1789
83. $Al(PO_3)_3$ ($G5_2$): AB9C3_cI208_220_c_3e_e 1802
84. AlN (cF40): AB_cF40_216_de_ce 1797
85. AlN (cI16): AB_cI16_217_c_c 1798
86. AlN (cI24): AB_cI24_220_a_b 1804
87. $AlNbO_4$: ABC4_mC24_12_i_i_4i 1534
88. $AlPO_4$ "low cristobalite type": AB4C_oC24_20_b_2c_a 1580
89. Albite ($NaAlSi_3O_8$, $S6_8$): ABC8D3_aP26_2_i_i_8i_3i 1506
90. Alluaudite [$NaMnFe_2(PO_4)_3$]: A2BCD12E3_mC76_15_f_e_b_6f_ef 1563

[§] α-ICl and LiAs have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[×] ζ-Fe_2N and α-PbO_2 have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[°] α-Potassium Nitrate (KNO_3) I, NH_4NO_3 III ($G0_{10}$), and Aragonite ($CaCO_3$, $G0_2$) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[*] β-B_2H_6 and B_2H_6 ($P2_1/c$) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[□] $KBrO_3$ ($G0_7$) and γ-Potassium Nitrate (KNO_3) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[‖] γ-$Y_2Si_2O_7$, Ammonium Persulfate [$(NH_4)_2S_2O_8$, $K4_1$], Monazite ($LaPO_4$), and $AgMnO_4$ ($H0_9$) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

[**] Ag_2O_3 and Zr_2Al_3 have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

91. Ammonium Chlorite (NH_4ClO_2): AB4CD2_tP16_113_c_f_a_e 1680
92. Ammonium Persulfate [$(NH_4)_2S_2O_8$, $K4_1$][†‖]: AB4C_mP24_14_e_4e_e 1551
93. Analcime ($NaAlSi_2O_6{\cdot}H_2O$, $S6_1$): A2B2C3D12E4_tI184_142_f_f_be_3g_g 1716
94. Andalusite (Al_2SiO_5, $S0_2$): A2B5C_oP32_58_eg_3gh_g 1616
95. Anhydrous $KAuBr_4$: AB4C_mP24_14_ab_4e_e .. 1550
96. Anthophyllite ($Mg_5Fe_2Si_8O_{22}(OH)_2$, $S4_4$): A2B5C22D2E8_oP156_62_d_c2d_2c10d_2c_4d ... 1630
97. Apophyllite ($KCa_4Si_8O_{20}F{\cdot}8H_2O$, $S5_2$): A4BC16DE28F8_tP116_128_h_a_2i_b_g3i_i 1696
98. Aragonite ($CaCO_3$, $G0_2$)[°]: ABC3_oP20_62_c_c_cd 1649
99. Arcanite (K_2SO_4, $H1_6$): A2B4C_oP28_62_2c_2cd_c 1629
100. Archerite (KH_2PO_4): A2BC4D_oF64_43_b_a_2b_a 1602
101. Arsenopyrite (FeAsS, $E0_7$): ABC_mP12_14_e_e_e 1556
102. Atacamite ($Cu_2(OH)_3Cl$): AB2C3D3_oP36_62_c_ac_cd_cd 1644
103. Au_2Nb_3: A2B3_tI10_139_e_ae 1705
104. $AuCsCl_3$ ($K7_6$): AB3C_tI20_139_ab_eh_d 1711
105. Autunite $\{Ca[(UO_2)(PO_4)]_2(H_2O)_{11}\}$: AB22C23D2E2_oP200_62_c_11d_3c10d_d_d 1642
106. Azurite [$Cu_3(CO_3)_2(OH)_2$, $G7_4$]: A2B3C2D8_mP30_14_e_ce_e_4e 1537
107. $B_{13}C_2$ "B_4C" ($D1_g$): A13B2_hR15_166_b2h_c .. 1736
108. B_2H_6 ($P2_1/c$)[*]: AB3_mP16_14_e_3e 1549
109. B_4SrO_7: A4B7C_oP24_31_2b_a3b_a 1586
110. $BaAl_2O_4$ ($H2_8$): A2BC6_hP18_182_f_b_gh 1754
111. $BaCd_{11}$: AB11_tI48_141_a_bdi 1716
112. $BaNi(CN)_4{\cdot}4H_2O$ ($H4_{22}$): AB4C4D4E_mC56_15_e_2f_2f_2f_a 1569
113. $BaNiSn_3$: ABC3_tI10_107_a_a_ab 1680
114. Bararite (Trigonal $(NH_4)_2SiF_6$, $J1_6$)[××]: A6B2C_hP9_164_i_d_a 1733
115. Barytocalcite ($BaCa(CO_3)_2$): AB2CD6_mP20_11_e_2e_e_2e2f 1522
116. Base-centered orthorhombic $Sr_4Ru_3O_{10}$: A10B3C4_oC68_64_2dfg_ad_2d 1659
117. Bassanite [$CaSO_4(H_2O)_{0.5}$, $H4_7$]: A2B2C9D2_mC90_5_ab2c_3c_b13c_3c 1509
118. Bastnäsite [$CeF(CO_3)$]: ABCD3_hP36_190_h_g_af_hi 1760
119. $BeSO_4{\cdot}4H_2O$ ($H4_3$): AB8C8D_tI72_120_c_2i_2i_b 1684
120. Berthierite ($FeSb_2S_4$, $E3_3$): AB4C2_oP28_62_c_4c_2c 1645
121. Bertrandite ($Be_4Si_2O_7(OH)_2$, $S4_6$): A4B7C2D2_oC60_36_2b_a3b_2a_b 1595
122. Bi_2GeO_5: A2BC5_oC32_36_b_a_a2b 1593
123. $Bi_3Ru_3O_{11}$: A3B11C3_cP68_201_be_efh_g 1774
124. Bischofite ($MgCl_2{\cdot}6H_2O$, $J1_7$): A2B12CD6_mC42_12_i_2i2j_a_ij 1524
125. Blossite (α-$Cu_2V_2O_7$): A2B7C2_oF88_43_b_a3b_b 1601
126. Boric Acid (H_3BO_3, $G5_1$): AB3C3_aP28_2_2i_6i_6i 1504
127. Bromocarnallite ($KMg(H_2O)_6(Cl,Br)_3$, $E2_6$): A3B6CD_tP44_85_bcg_3g_ac_e 1669
128. Brucite [$Mg(OH)_2$][¶¶]: A2BC2_hP5_164_d_a_d .. 1732
129. $C_{19}Sc_{15}$: A19B15_tP68_114_bc4e_ac3e 1681
130. COCl: ABC_oP24_61_c_c_c 1627
131. Ca_2RuO_4: A2B4C_oP28_61_c_2c_a 1624
132. Ca_2UO_5: A2B5C_mP32_14_2e_5e_ab 1540
133. $Ca_3Al_2(OH)_{12}$ ($J2_3$): A2B3C12D12_cI232_230_a_c_h_h 1836
134. Ca_3UO_6: A3B6C_mP20_4_3a_6a_a 1508
135. CaB_2O_4 (III): A2BC4_oP84_33_6a_3a_12a 1589
136. CaB_2O_4 (IV): A2BC4_cP84_205_d_ac_2d 1780
137. CaB_2O_4 I ($E3_2$): A2BC4_oP28_60_d_c_2d 1621
138. $CaBe_2Ge_2$: A2BC2_tP10_129_ac_c_bc 1698
139. CaC_2-I ($C11_a$): A2B_tI6_139_e_a 1706
140. CaC_2-III: A2B_mC12_12_2i_i 1526
141. $CaCu_4P_2$[°°]: AB4C2_hR7_166_a_2c_c 1739
142. $CaO_2(H_2O)_8$: AB8C2_tP22_124_a_n_h 1692
143. $CaSi_2$ ($C12$): AB2_hR6_166_c_2c 1738
144. $CaUO_4$: AB4C_hR6_166_b_2c_a 1740
145. Calaverite ($AuTe_2$): AB2_mP12_7_2a_4a 1515
146. Calciborite (CaB_2O_4 II): A2BC4_oP56_56_2e_e_4e 1613
147. Carnallite [$Mg(H_2O)_6KCl_3$]: A3B12CDE6_oP276_52_d4e_18e_ce_de_2d8e 1607
148. Catapleiite ($Na_2ZrSi_3O_9{\cdot}2H_2O$): A2B3C9D3E_mC144_15_2f_bcdef_9f_3f_ae 1559

[†]Ammonium Persulfate [$(NH_4)_2S_2O_8$, $K4_1$] and Monazite ($LaPO_4$) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[××]$K_2Pt(SCN)_6$ ($H6_3$), Bararite (Trigonal $(NH_4)_2SiF_6$, $J1_6$), and K_2GeF_6 ($J1_{13}$) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

[¶¶]Ce_2O_2S and Brucite [$Mg(OH)_2$] have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

[°°]$MnBi_2Te_4$ and $CaCu_4P_2$ have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

149. Cd_3As_2: A2B3_tI160_142_deg_3g 1717
150. Cd(OH)Cl ($E0_3$): ABCD_hP8_186_b_b_a_a 1758
151. Ce_2O_2S[¶¶]: A2B2C_hP5_164_d_d_a 1732
152. $CeCu_2$: AB2_oI12_74_e_h 1666
153. Cervantite (α-Sb_2O_4): A2B_oP24_33_4a_2a 1590
154. Chabazite ($Ca_{1.4}Sr_{0.3}Al_{3.8}Si_{8.3}O_{24}{\cdot}13H_2O$, $S3_4$ (I)): A5B21C24D12_hR62_166_a2c_ehi_fg2h_i 1737
155. Chalcanthite ($CuSO_4{\cdot}5H_2O$, $H4_{10}$): AB10C9D_aP42_2_ae_10i_9i_i 1504
156. Chalcocyanite ($CuSO_4$): AB4C_oP24_62_a_2cd_c 1646
157. Chiolite ($Na_5Al_3F_{14}$, $K7_5$): A3B14C5_tP44_128_ac_ehi_bg 1696
158. Chlorite [$Mg_3(Mg_2Al)(Si_3Al)O_{10}(OH)_8$, $S5_5$]: A3B5C4D2_mC112_15_a3ef_5f_4f_2f 1566
159. Chrysotile ($H_4Mg_3Si_2O_9$, $S4_5$): AB6C11D6E4_mC112_12_e_gi2j_i5j_2i2j_2j 1531
160. Chrysotile ($Mg_3Si_2O_5(OH)_4$): A3B5C4D2_mC56_9_3a_5a_4a_2a 1518
161. Clinocervantite (β-Sb_2O_4): A2B_mC24_15_2f_ce 1564
162. Co_2Al_9 ($D8_d$): A9B2_mP22_14_a4e_e 1546
163. $Co_2B_2O_5$: A2B2C5_aP18_2_2i_2i_5i 1500
164. $Co_3(SeO_3)_3{\cdot}H_2O$: A3B2C10D3_aP36_2_ah2i_2i_10i_3i 1502
165. Co_9S_8 ($D8_9$): A9B8_cF68_225_af_ce 1815
166. Colquiriite ($LiCaAlF_6$): ABC6D_hP18_163_d_b_i_c 1731
167. Columbite ($FeNb_2O_6$, $E5_1$): AB2C6_oP36_60_c_d_3d 1623
168. Copper (I) Azide (CuN_3): AB3_tI32_88_d_cf 1674
169. Copper (II) Azide [$Cu(N_3)_2$]: AB6_oP28_62_c_6c 1647
170. Cr_5O_{12}: A5B12_oP68_60_c2d_6d 1622
171. Cr-233 Quasi-One-Dimensional Superconductor ($K_2Cr_3As_3$): A3B3C2_hP16_187_jk_jk_ck 1758
172. $CrCl_3(H_2O)_6$ ($J2_2$): A3BC6_hR20_167_e_b_f ... 1745
173. Crancrinite ($Na_6Ca_2Al_6Si_6O_{24}(CO_3)_2$, $S3_3$ (I)): A3BCD3E15F3_hP52_173_c_b_b_c_5c_c 1747
174. Cronstedtite {Fe(Fe,Si)[$(OH)_2$,O]O_3, $S5_7$}: AB3C2D_hR7_160_a_b_2a_a 1727
175. Cryolite (Na_3AlF_6, $J2_6$): AB6C3_mP20_14_a_3e_de 1552
176. $Cs_{11}O_3$: A11B3_mP56_14_11e_3e 1536
177. Cs_2Sb: A2B_oP24_62_4c_2c 1634
178. Cs_2Se: A2B_oF24_43_b_a 1602
179. $Cs_3As_2Cl_9$ ($K7_3$): A2B9C3_hP14_150_d_eg_ad .. 1723
180. Cs_3CoCl_5 ($K3_1$): A5BC3_tI36_140_cl_b_ah 1715
181. $Cs_3Cr_2Cl_9$: A9B2C3_hP28_194_hk_f_bf 1767
182. $Cs_3Tl_2Cl_9$ ($K7_2$): A9B3C2_hR28_167_ef_e_c 1747
183. $Cs_6W_{11}O_{36}$: A6B36C11_mC212_9_6a_36a_11a .. 1519
184. Cs_7O: A7B_hP24_187_ai2j2kn_j 1759
185. CsB_4O_6F: A4BCD6_oP48_33_4a_a_a_6a 1591
186. $CsFeS_2$ (100 K): ABC2_oI16_71_g_i_eh 1664
187. CsO: AB_oI8_71_g_i 1665
188. $CsSO_3$ ($K1_2$): AB3C_hP20_190_ac_i_f 1760
189. $Cu_2Pb(SeO_3)_2Br_2$: A2B2C6DE2_oC52_63_g_e_fh_c_f 1652
190. $Cu_3[Fe(CN)_6]_2{\cdot}xH_2O$ ($J2_5$, $x \approx 3$): A6B9CD2E6_cF96_225_e_bf_a_c_e 1813
191. Cu(OH)Cl: ABCD_mP16_14_e_e_e_e 1555
192. Cubic Cu_2OSeO_3: A2B4C_cP56_198_ab_2a2b_2a 1771
193. Cubic CuPt ($L1_3$ (I), $D4$): AB_cF32_227_c_d ... 1834
194. Danburite ($CaB_2Si_2O_8$, $S6_3$): A2BC8D2_oP52_62_d_c_2c3d_d 1632
195. Diamminetriamidodizinc Chloride ($[Zn_2(NH_3)_2(NH_2)_3]Cl$): AB12C5D2_oP40_18_a_6c_b2c_c 1574
196. Diaspore (AlOOH, $E0_2$): ABC2_oP16_62_c_c_2c 1647
197. Diopside [$CaMg(SiO_3)_2$, $S4_1$]: ABC6D2_mC40_15_e_e_3f_f 1572
198. Dodecatungstophosphoric Acid Hexahydrate [$H_3PW_{12}O_{40}{\cdot}6H_2O$]: A27B52CD12_cP184_224_dl_eh3k_a_k 1806
199. Dolomite [$MgCa(CO_3)_2$, $G1_1$]: A2BCD6_hR10_148_c_a_b_f 1720
200. Double Perovskite (Ba_2MnWO_6): A2BC6D_cF40_225_c_a_e_b 1812
201. Enstatite ($MgSiO_3$, $S4_3$): AB3C_oP80_61_2c_6c_2c 1627
202. Epididymite ($BeHNaO_8Si_3$, $S4_7$): ABCD8E3_oP112_62_d_2c_d_4c6d_3d 1649
203. Eriochalcite ($CuCl_2{\cdot}2H_2O$, $C45$): A2BC4D2_oP18_53_h_a_i_e 1609
204. $EuIn_2P_2$: AB2C2_hP10_194_a_f_f 1769
205. Eudidymite ($BeHNaO_8Si_3$): A2B4C2D17E6_mC124_15_f_2f_f_e8f_3f 1561
206. Eulytine ($Bi_4(SiO_4)_3$, $S1_5$): A4B12C3_cI76_220_c_e_a 1801
207. $Fe_2(CO)_9$ ($F4_1$): A9B2C9_hP40_176_hi_f_hi 1752
208. Fe_2N (approximate, $L'3_0$)[∂∂]: AB_hP4_194_c_a .. 1771
209. Fe_3PO_7: A3B7C_hR11_160_b_a2b_a 1726
210. Fe_8N ($D2_g$): A8B_tI18_139_deh_a 1709
211. FeF_3 ($D0_{12}$): A3B_hR8_167_e_b 1745
212. Ferroelectric $NH_4H_2PO_4$: A6BC4D_oP48_19_6a_a_4a_a 1577

[∂∂] $LiZn_2$ (C_k) and Fe_2N (approximate, $L'3_0$) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

213. Ferroelectric $NaNO_2$ ($F5_5$): ABC2_oI8_44_a_a_c 1604
214. Fluorapatite [$Ca_5F(PO_4)_3$, $H5_7$]: A5BC12D3_hP42_176_fh_a_2hi_h 1752
215. Ga_2Mg_5 ($D8_g$): A2B5_oI28_72_j_bfj 1665
216. $GaMo_4S_8$: AB4C8_cF52_216_a_e_2e 1795
217. Gd_2SiO_5 (RE_2SiO_5 X1): A2B5C_mP32_14_2e_5e_e 1540
218. Gwihabaite [NH_4NO_3 (V)]: A4B2C3_tP72_77_8d_ab2c2d_6d 1667
219. Gypsum ($CaSO_4{\cdot}2H_2O$, $H4_6$): AB4C6D_mC48_15_e_2f_3f_e 1570
220. $H_3PW_{12}O_{40}{\cdot}29H_2O$ ($H4_{21}$): A29B40CD12_cF656_227_ae2fg_e3g_b_g 1826
221. $H_3PW_{12}O_{40}{\cdot}3H_2O$: A3B40CD12_cP112_224_d_e3k_a_k 1808
222. Hambergite [$Be_2BO_3(OH)$, $G7_2$]: AB2CD4_oP64_61_c_2c_c_4c 1626
223. Hauyne [$(Na_{0.5}Ca_{0.3}K_{0.2})_8(Al_6Si_6O_{24})(SO_4)_{1.5}$, $S6_9$]: A3B4C4D4E16F4G3_cP76_218_c_e_e_e_ei_e_d .. 1799
224. Hemimorphite ($Zn_4Si_2O_7(OH)_2{\cdot}H_2O$, $S2_2$): A2B5CD2_oI40_44_2c_abcde_d_e 1603
225. Hexagonal Delafossite ($CuAlO_2$): ABC2_hP8_194_a_c_f 1770
226. Hexagonal WO_3: A3B_hP12_191_gl_f 1762
227. Hg_2O_2NaI: A2BCD2_hP18_180_f_c_b_i 1753
228. Hg_2TiCu Inverse Heusler: AB2C_cF16_216_b_ad_c 1794
229. $HgCl_2{\cdot}2HgO$: A2B3C2_mP14_14_e_ae_e 1538
230. High-Temperature Cryolite (Na_3AlF_6): AB6C3_oI20_71_a_in_cj 1664
231. High-Temperature Cubic $KClO_4$ ($H0_5$): ABC4_cF24_216_b_a_e 1796
232. High-Temperature Mo_8O_{23}: A8B23_mP62_13_4g_c11g 1535
233. $HoMn_2O_5$: AB2C5_oP32_55_g_fh_eghi 1613
234. $HoSb_2$: AB2_oC6_21_a_k 1581
235. Huanzalaite ($MgWO_4$, $H0_6$): AB4C_mP12_13_f_2g_e 1536
236. In_4Se_3: A4B3_oP28_58_4g_3g 1618
237. InS: AB_oP8_58_g_g 1619
238. Jacutingaite (Pt_2HgSe_3): AB2C3_hP12_164_d_ae_i 1734
239. $K_2CuCl_4{\cdot}2H_2O$ ($H4_1$): A4BC4D2E2_tP26_136_fg_a_j_d_e 1702
240. K_2GeF_6 ($J1_{13}$)[××]: A6BC2_hP9_164_i_a_d 1734
241. $K_2HgCl_4{\cdot}H_2O$ ($E3_4$): A4BCD2_oP32_55_ghi_f_e_gh 1612
242. K_2NbF_7 ($K6_2$): A7B2C_mP40_14_7e_2e_e 1545
243. $K_2Ni(CN)_4$: A4B2C4D_mP22_14_2e_e_2e_a 1542
244. K_2NiF_4: A4B2C_tI14_139_ce_e_a 1707
245. $K_2OsO_2Cl_4$ ($J1_5$): A4B2C2D_tI18_139_h_d_e_a . 1706
246. $K_2Pt(SCN)_6$ ($H6_3$)[××]: A2BC6_hP9_164_d_a_i .. 1733
247. $K_2Pt(SCN)_6{\cdot}2H_2O$: A6B4C2D6E2FG6_mP54_14_3e_2e_e_3e_e_a_3e . 1544
248. K_2PtCl_4 ($H1_5$): A4B2C_tP7_123_j_e_a 1689
249. $K_2S_2O_5$ ($K0_1$): A2B5C2_mP18_11_2e_e2f_2e ... 1521
250. $K_2S_3O_6$ ($K5_1$): A2B6C3_oP44_62_2c_2c2d_3c .. 1632
251. $K_2Sn(OH)_6$ ($H6_2$): A6B2C6D_hR15_148_f_c_f_a 1721
252. $K_2SnCl_4{\cdot}H_2O$: A4BC2D_oP32_62_2cd_c_d_c ... 1639
253. $K_2SnCl_4{\cdot}H_2O$ ($E3_5$): A4BC2D_oP32_62_2cd_b_2c_a 1638
254. $K_2Ti_2O_5$: A2B5C2_mC18_12_i_a2i_i 1526
255. $K_3Co(NO_2)_6$ ($J2_4$): AB3C6D12_cF88_202_a_bc_e_h 1774
256. K_3CrO_8: AB3C8_tI24_121_a_bd_2i 1686
257. $K_3TlCl_6{\cdot}2H_2O$ ($J3_1$): A6B2C3D_tI168_139_egikl2m_ejn_bh2n_acf 1708
258. $K_3W_2Cl_9$ ($K7_1$): A9B3C2_hP28_176_hi_af_f 1753
259. $K_4[Mo(CN)_8]{\cdot}2H_2O$ ($F2_1$): A8B4C4DE8F2_oP108_62_4c2d_2d_2cd_c_4c2d_d 1641
260. K(SH) ($B22$): AB_hR2_166_a_b 1742
261. $KAuBr_4{\cdot}2H_2O$ ($H4_{19}$): AB4C2D_mP32_14_e_4e_2e_e 1550
262. $KBe_2BO_3F_2$: AB2C2DE3_hR9_155_b_c_c_a_e .. 1725
263. $KBrO_3$ ($G0_7$)[□]: ABC3_hR5_160_a_a_b 1728
264. $KFeS_2$ ($F5_a$): ABC2_mC16_15_e_e_f 1572
265. KH_2PO_4 ($H2_2$): A4BC4D_tI40_122_e_b_e_a 1687
266. KHF_2 ($F5_2$): A2BC_tI16_140_h_d_a 1712
267. $KICl_4{\cdot}H_2O$ ($H0_{10}$): A4BCD_mP28_14_4e_e_e_e 1543
268. KNO_2 III[‡]: ABC2_mP16_14_e_e_2e 1553
269. KSO_3 ($K1_1$): AB3C_hP30_150_ef_3g_c2d 1724
270. Kesterite [$Cu_2(Zn,Fe)SnS_4$]: A2BCD4_tI16_82_ac_b_d_g 1668
271. Kotoite ($Mg_3(BO_3)_2$): A2B3C6_oP22_58_g_af_gh 1616
272. Kyanite (Al_2SiO_5, $S0_1$): A2B5C_aP32_2_4i_10i_2i 1501
273. La_3BWO_9 ($P3$): AB3C9D_hP28_143_2a_2d_6d_bc 1718
274. La_3BWO_9 ($P6_3$): AB3C9D_hP28_173_a_c_3c_b . 1749
275. La_3CuSiS_7: AB3C7D_hP24_173_a_c_b2c_b 1748
276. $LaFe_4P_{12}$: A4BC12_cI34_204_c_a_g 1775
277. LaH_{10} High-T_c Superconductor: A10B_cF44_225_cf_b 1811
278. LaOAgS: ABCD_tP8_129_b_c_a_c 1699
279. Lepidocrocite (γ-FeO(OH), $E0_4$): AB2C2_oC20_63_c_f_2c 1655
280. Li_2CN_2[⊗‡‡]: AB2C2_tI10_139_a_d_e 1710
281. Li_2PrO_3: A2B3C_oC12_65_h_bh_a 1660

[‡] KNO_2 III and Manganite (γ-MnO(OH), $E0_6$) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

282. $Li_2SO_4{\cdot}H_2O$ ($H4_8$): A2B2C5D_mP20_4_2a_2a_5a_a 1508
283. Li_7TaO_6: A8B6C_hR15_148_cf_f_a 1722
284. LiAs[§]: AB_mP16_14_2e_2e 1556
285. $LiClO_4{\cdot}3H_2O$ ($H4_{18}$): AB6CD7_hP30_186_b_d_a_b2c 1757
286. $LiCuVO_4$: ABC4D_oI28_74_a_d_hi_e 1667
287. $LiGaO_2$: ABC2_oP16_33_a_a_2a 1591
288. $LiKSO_4$ ($H1_4$): ABC4D_hP14_173_a_b_bc_b 1750
289. $LiNb_6O_{15}F$: ABC6D15_oP46_51_f_d_2e2i_aef4i2j 1606
290. $LiOH{\cdot}H_2O$ ($B36$): A3BC2_mC24_12_ij_h_gi 1528
291. $LiZn_2$ (C_k)[∂∂]: AB_hP4_194_a_c 1770
292. Low-Temperature $(NH_3CH_3)Al(SO_4)_2{\cdot}12H_2O$: ABC30DE20F2_oP220_29_a_a_30a_a_20a_2a 1582
293. Low-Temperature $GaMo_4S_8$: AB4C8_hR13_160_a_ab_2a2b 1727
294. Low-Temperature Mo_8O_{23}: A8B23_mP124_7_16a_46a 1513
295. Lu_2CoGa_3: AB3C2_hP24_194_f_k_bh 1769
296. Lueshite ($NaNbO_3$): ABC3_oP40_57_cd_e_cd2e 1615
297. Maghemite (γ-Fe_2O_3, $D5_7$): A2B3_cP60_212_bcd_ace 1788
298. Magnetoplumbite ($PbFe_{12}O_{19}$): A12B19C_hP64_194_ab2fk_efh2k_d 1764
299. Manganese-leonite 110 K $[K_2Mn(SO_4)_2{\cdot}4H_2O]$: A8B2CD12E2_mP100_14_8e_2e_ad_12e_2e 1545
300. Manganese-leonite 185 K $[K_2Mn(SO_4)_2{\cdot}4H_2O]$: A8B2CD12E2_mC200_15_8f_2f_ce_2e11f_2f 1568
301. Manganese-leonite $[K_2Mn(SO_4)_2{\cdot}4H_2O$, $H4_{23}]$: A8B2CD15E2_mC112_12_2i3j_j_ad_g4i5j_2i 1529
302. Manganite (γ-MnO(OH), $E0_6$)[‡]: ABC2_mP16_14_e_e_2e 1553
303. Marialite Scapolite $[Na_4Cl(AlSi_3)_3O_{24}$, $S6_4]$: AB4C24D12_tI82_87_a_h_2h2i_hi 1672
304. Mayenite ($12CaO{\cdot}7Al_2O_3$, $K7_4$, C12A7): A7B12C19_cI152_220_bc_2d_ace 1801
305. Mercury (II) Azide $[Hg(N_3)_2]$: AB6_oP28_29_a_6a 1582
306. Mercury Cyanide $[Hg(CN)_2$, $F1_1]$: A2BC2_tI40_122_e_d_e 1687
307. Meta-autunite (I) $[Ca(UO_2)_2(PO_4)_2{\cdot}6H_2O$, $H5_{10}]$: AB4C6DE_tP26_129_c_j_2ci_a_c 1698
308. Mg_2Cu (C_b): AB2_oF48_70_g_fg 1663
309. $Mg_3Cr_2Al_{18}$: A18B2C3_cF184_227_fg_d_ac 1823
310. Mg_3P_2 ($D5_5$): A3B2_cP10_224_d_b 1808
311. Mg_3Ru_2: A3B2_cP20_213_d_c 1790
312. $Mg(ClO_4)_2{\cdot}6H_2O$ ($H4_{11}$): A2B6CD8_oP34_31_2a_2a2b_a_4a2b 1586
313. $Mg(NH_3)_2Cl_2$ ($E1_3$): A2B8CD2_oC26_65_h_r_a_i 1661
314. $MgCuAl_2$ ($E1_a$): A2BC_oC16_63_f_c_c 1653
315. Mn_3As ($D0_d$): AB3_oC16_63_c_3c 1657
316. $MnBi_2Te_4$[°°]: A2BC4_hR7_166_c_a_2c 1736
317. MnCuP: ABC_oP12_62_c_c_c 1651
318. $MnF_{2-x}(OH)_x$: A2BC2_oP10_34_c_a_c 1592
319. $MnPS_3$: ABC3_mC20_12_g_i_ij 1534
320. $Mo_{17}O_{47}$: A17B47_oP128_32_a8c_a23c 1587
321. Mo_4P_3: A4B3_oP56_62_8c_6c 1637
322. MoP_2: AB2_oC12_36_a_2a 1595
323. $MoPO_5$: AB5C_tP14_85_c_cg_b 1669
324. Monazite ($LaPO_4$)[†‖]: AB4C_mP24_14_e_4e_e 1551
325. Monoclinic Co_4Al_{13}: A13B4_mC102_8_17a11b_8a2b 1515
326. Monoclinic Cu_2OSeO_3: A2B4C_mP28_14_abe_4e_e 1539
327. Monoclinic $FeTlSe_2$: AB2C_mC16_12_g_2i_i 1530
328. Morenosite ($NiSO_4{\cdot}7H_2O$, $H4_{12}$): A14BC11D_oP108_19_14a_a_11a_a 1575
329. Murataite $[(Y,Na)_6(Zn,Fe)_5Ti_{12}O_{29}(O,F)_{10}F_4]$: A16B40C12D6E5_cF316_216_eh_e2g2h_h_f_be 1791
330. Muscovite ($KH_2Al_3Si_3O_{12}$, $S5_1$): A2BC10D2E4_mC76_15_f_e_5f_f_2f 1562
331. NH_4Br ($B25$): AB4C_tP12_129_c_i_a 1699
332. NH_4CdCl_3 ($E2_4$): AB3C_oP20_62_c_3c_c 1645
333. NH_4ClBrI ($F5_{14}$): ABCD_oP16_62_c_c_c_c 1650
334. $NH_4H_2PO_2$ ($F5_7$): A2BC2D_oC24_67_m_a_n_g 1662
335. $NH_4H_2PO_4$: A8BC4D_tI56_122_2e_b_e_a 1688
336. NH_4HF_2 ($F5_8$): A2BC_oP16_53_eh_ab_g 1610
337. NH_4HgCl_3 ($E2_5$): A3BC_tP5_123_cg_a_d 1689
338. NH_4I_3 ($D0_{16}$): A3B_oP16_62_3c_c 1635
339. NH_4NO_3 I ($G0_8$): AB_cP2_221_a_b 1806
340. NH_4NO_3 II ($G0_9$): ABC3_tP10_100_b_a_bc 1678
341. NH_4NO_3 III ($G0_{10}$)[°]: ABC3_oP20_62_c_c_cd 1648
342. NH_4NO_3 IV ($G0_{11}$): A4B2C3_oP18_59_ef_ab_af 1620
343. $NH_4Pb_2Br_5$ ($K3_4$): A5BC2_tI32_140_bl_a_h 1714
344. NO_2 (Modern, $C26$): AB2_cI36_204_d_g 1777
345. $Na_{0.74}CoO_2$: AB2C2_hP10_194_a_bc_f 1768
346. $Na_2Ca_6Si_4O_{15}$: A6B2C15D4_mP54_7_6a_2a_15a_4a 1512
347. Na_2CaSiO_4 ($S6_6$): AB2C4D_cP32_198_a_2a_ab_a 1772
348. Na_2CrO_4 ($H1_8$): AB2C4_oC28_63_c_bc_fg 1656
349. $Na_2Mo_2O_7$: A2B2C7_oC88_64_ef_df_3f2g 1660
350. Na_2PrO_3: A2B3C_mC48_15_aef_3f_2e 1560
351. Na_2SO_3 ($G3_2$): A2B3C_hP12_147_abd_g_d 1720
352. $Na_4Ge_9O_{20}$: A9B4C20_tI132_88_a2f_f_5f 1673
353. $NaAlCl_4$: AB4C_oP24_19_a_4a_a 1578
354. $NaC_5H_{11}O_8S$: A5B11CD8E_aP26_1_5a_11a_a_8a_a 1499
355. $NaCr(SO_4)_2{\cdot}12H_2O$ Alum: AB12CD8E2_cP96_205_a_2d_b_cd_c 1782
356. $NaMn_7O_{12}$: A7BC12_cI40_204_bc_a_g 1776

357. $NaNb_6O_{15}F$: ABC6D15_oC46_38_b_b_2a2d_2ab4d2e 1597
358. $NaNbO_3$: ABC3_oP40_17_abcd_2e_abcd4e 1573
359. NaP: AB_oP16_19_2a_2a 1579
360. NaS_2: AB2_tI48_122_cd_2e 1688
361. $NaSb(OH)_6$ ($J1_{11}$): AB6C_tP32_86_d_3g_c 1671
362. $NaSbF_4(OH)_2$ ($J1_{12}$): A6BC_hP16_163_i_b_c ... 1730
363. $NaSbF_6$: A6BC_cP32_205_d_b_a 1781
364. Nacrite [$Al_2Si_2O_5(OH)_4$, $S5_4$]: A2B4C9D2_mC68_9_2a_4a_9a_2a 1518
365. Nahcolite ($NaHCO_3$, $G0_{12}$): ABCD3_mP24_14_e_e_e_3e 1554
366. Natrolite ($Na_2Al_2Si_3O_{10}{\cdot}2H_2O$, $S6_{10}$): A2B4C2D12E3_oF184_43_b_2b_b_6b_ab 1600
367. $Nb_2Pd_3Se_8$: A2B3C8_oP26_55_h_ag_2g2h 1611
368. $Nb_2Zr_6O_{17}$: A2B17C6_oI100_46_ab_b8c_3c 1605
369. Nb_3O_7F: A3B8_oC22_65_ag_bd2gh 1661
370. $NbAs_2$: A2B_mC12_5_2c_c 1510
371. $NbTe_2$: AB2_mC18_12_ai_3i 1530
372. $Nd_2Fe_{14}B$: AB14C2_tP68_136_f_ce2j2k_fg 1703
373. $Nd_4Re_2O_{11}$: A4B11C2_tP68_86_2g_ab5g_g 1670
374. $Nd(BrO_3)_3{\cdot}9H_2O$ ($G2_2$): A3B9CD9_hP44_186_c_3c_b_cd 1756
375. Nevskite (BiSe): AB_hP12_164_c2d_c2d 1735
376. Ni_3Si_2: A3B2_oC80_36_4a4b_2a3b 1594
377. $Ni(H_2O)_6SnCl_6$ ($I6_1$): A6B6CD_hR14_148_f_f_b_a 1721
378. $Ni(NO_3)_2(H_2O)_6$: A12B2CD12_aP54_2_12i_2i_i_12i 1499
379. Norbergite [$Mg(F,OH)_2 \cdot Mg_2SiO_4$, $S0_7$]: A2B3C4D_oP40_62_d_cd_2cd_c 1629
380. O(OH)Y: ABC_mP6_11_e_e_e 1523
381. Original β-WO_3 (*obsolete*): A3B_oP32_62_ab4c_2c 1636
382. Orpiment (As_2S_3, $D5_f$): A2B3_mP20_14_2e_3e . 1538
383. Orthorhombic Co_4Al_{13}: A13B4_oP102_31_17a11b_8a2b 1584
384. Orthorhombic CrO_3: AB3_oC16_40_b_a2b 1599
385. Orthorhombic $Sr_4Ru_3O_{10}$: A10B3C4_oP68_55_2e2fgh2i_adef_2e2f 1610
386. Os_4Al_{13}: A13B4_mC34_12_b6i_2i 1524
387. P_4Se_3: A4B3_oP112_62_8c4d_4c4d 1636
388. $PNCl_2$ ($E1_4$): A2BC_tP32_86_2g_g_g 1670
389. Paralstonite ($BaCa(CO_3)_2$): AB2CD6_hP30_150_e_c2d_f_3g 1723
390. Pararealgar (AsS)[¶]: AB_mP32_14_4e_4e 1557
391. Paratellurite (α-TeO_2): A2B_tP12_92_b_a 1677
392. Parawollastonite ($CaSiO_3$, $S3_3$ (II)): AB3C_mP60_14_3e_9e_3e 1548
393. Parkerite ($Ni_3Bi_2S_2$): AB2C_oP8_51_e_be_f 1606
394. $Pb(NO_3)_2$ ($G2_1$): A2B6C_cP36_205_c_d_a 1779
395. Pd_5Pu_3: A5B3_oC32_63_cfg_ce 1654
396. $Pd(NH_3)_4Cl_2{\cdot}H_2O$ ($H4_9$): A2BC4D_tP16_127_h_d_i_a 1695
397. Phase II $Cd_2Re_2O_7$: A2B7C2_tI44_119_i_bdefgh_i 1683
398. Phase III $Cd_2Re_2O_7$: A2B7C2_tI44_98_f_bcde_f 1677
399. Phosgenite [$Pb_2Cl_2(CO_3)$]: AB2C3D2_tP32_127_g_eh_gk_k 1695
400. Possible δ-$Gd_2Si_2O_7$: A2B7C2_oP44_33_2a_7a_2a 1589
401. Possible δ-$Y_2Si_2O_7$: A7B2C2_oP44_62_3c2d_2c_d 1640
402. Predicted High-Pressure $YCaH_{12}$: AB12C_cP14_221_a_h_b 1805
403. Predicted Li_2MgH_{16} 300 GPa: A16B2C_hP19_164_2d2i_d_b 1731
404. Predicted Li_2MgH_{16} High-Temperature Superconductor (250 GPa): A16B2C_cF152_227_eg_d_a 1822
405. Predicted Phase IV $Cd_2Re_2O_7$: A2B7C2_oF88_22_k_bdefghij_k 1581
406. Proposed 300 GPa HfH_{10}: A10B_hP22_194_bhj_c 1763
407. Protoanthophyllite ($H_2Mg_7Si_8O_{24}$): A2B7C24D8_oP82_58_g_ae2f_2g5h_2h 1617
408. Pseudobrookite (Fe_2TiO_5, $E4_1$)[††]: A2B5C_oC32_63_f_c2f_c 1652
409. Pt_2Sn_3 ($D5_b$): A2B3_hP10_194_f_bf 1765
410. $Pu_{31}Rh_{20}$: A31B20_tI204_140_b2gh3m_ac2fh3l .. 1713
411. Pyrophyllite [$AlSi_2O_5(OH)$, $S5_6$]: AB5CD2_mC72_15_f_5f_f_2f 1571
412. $Rb_2C_2O_4{\cdot}H_2O$: A2BC4D2_mC36_15_f_e_2f_f 1563
413. $Rb_2CaCu_6(PO_4)_4O_2$: AB6C18D4E2_mC62_5_a_2b2c_9c_2c_c 1511
414. $Rb_2Mo_2O_7$: A2B7C2_oC88_40_abc_2b6c_a3b ... 1598
415. $RbNO_3$ (IV): AB3C_hP45_144_3a_9a_3a 1719
416. Re_3B: AB3_oC16_63_c_cf 1657
417. ReB_3: A3B_hP8_194_af_c 1767
418. Realgar (AsS, B_l)[¶]: AB_mP32_14_4e_4e 1558
419. Retgersite (α-$NiSO_4{\cdot}6H_2O$, $H4_5$): A12BC10D_tP96_92_6b_a_5b_a 1676

[¶]Pararealgar (AsS) and Realgar (AsS, B_l) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[††]Pseudobrookite (Fe_2TiO_5, $E4_1$) and Ta_2NiS_5 have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

420. $Rh_{20}Si_{13}$: A10B7_hP34_176_c3h_b2h 1750
421. $RhCl_2(NH_3)_5Cl$ ($J1_8$): A3B15C5D_oP96_62_cd_3c6d_3cd_c 1634
422. Rhombohedral $CuTi_2S_4$: AB4C2_hR28_166_2c_2c2h_abh 1739
423. Rhombohedral Delafossite ($CuFeO_2$): ABC2_hR4_166_a_b_c 1741
424. Rinneite ($K_3NaFeCl_6$): A6BC3D_hR22_167_f_b_e_a 1746
425. Rosiaite ($PbSb_2O_6$)[§§]: A6BC2_hP9_162_k_a_d .. 1730
426. $Ru_{11}B_8$: A8B11_oP38_55_g3h_a3g2h 1612
427. RuB_2: A2B_oP6_59_f_a 1620
428. Rynersonite (Orthorhombic $CaTa_2O_6$): AB6C2_oP36_62_c_2c2d_d 1646
429. Sanguite ($KCuCl_3$): A3BC_mP20_14_3e_e_e 1541
430. Sanidine ($KAlSi_3O_8$, $S6_7$): AB8C4_mC52_12_i_gi3j_2j 1533
431. Santite ($KB_5O_8{\cdot}4H_2O$, $K3_5$): A5B8CD12_oC104_41_a2b_4b_a_6b 1599
432. $Sb_4O_5Cl_2$: A2B5C4_mP22_14_e_c2e_2e 1539
433. $SbCl_5{\cdot}POCl_3$: A8BCD_oP44_62_4c2d_c_c_c 1642
434. SbI_3S_{24}: A3B24C_hR28_160_b_2b3c_a 1725
435. Scheelite ($CaWO_4$, $H0_4$): AB4C_tI24_88_b_f_a . 1675
436. Senarmontite (Sb_2O_3, $D6_1$): A3B2_cF80_227_f_e 1831
437. Shandite ($Ni_3Pb_2S_2$): A3B2C2_hR7_166_d_ab_c 1737
438. Shcherbinaite (V_2O_5) (*revised*): A5B2_oP14_59_a2f_f 1620
439. Si_{24} Clathrate: A_oC24_63_3f 1659
440. Si_2N_2O: A2BC2_oC20_36_b_a_b 1593
441. SiAs: AB_mC24_12_3i_3i 1535
442. Sillimanite (Al_2SiO_5, $S0_3$): A2B5C_oP32_62_bc_3cd_c 1631
443. $Sm_{11}Cd_{45}$: A45B11_cF448_216_bd4efg5h_ac2eh . 1793
444. SnI_4 ($D1_1$): A4B_cP40_205_cd_c 1780
445. Sodalite [$Na_4(AlSiO_4)_3Cl$, $S6_2$]: A3BC4D12E3_cP46_218_d_a_e_i_c 1800
446. Sr_2MnTeO_6: AB6C2D_mP20_14_a_3e_e_d 1552
447. Sr_2NiTeO_6: AB6C2D_mC40_12_ad_gh4i_j_bc ... 1532
448. Sr_2NiWO_6: AB6C2D_tI20_87_a_eh_d_b 1673
449. $Sr_3Ti_2O_7$: A7B3C2_tI24_139_aeg_be_e 1709
450. $Sr_4Ti_3O_{10}$: A10B4C3_tI34_139_c2eg_2e_ae 1704
451. $Sr(OH)_2(H_2O)_8$: A18B10C_tP116_130_2c4g_2c2g_a 1700
452. $SrCl_2{\cdot}(H_2O)_6$: A2B12C6D_hP21_150_d_2g_ef_a 1722
453. $SrCu_2(BO_3)_2$: A2B2C6D_tI44_121_i_i_ij_c 1685
454. $SrUO_4$: A4BC_oP24_57_cde_d_a 1615
455. Staurolite ($Al_5Fe_2O_{10}(OH)_2Si_2$): A5B2C10D2E2_mC84_12_acghj_bdi_5j_2i_j 1528
456. Steklite [$KAl(SO_4)_2$, $H3_2$]: ABC8D2_hP12_150_b_a_dg_d 1724
457. Sulphohalite [$Na_6ClF(SO_4)_2$, $H5_8$]: ABC6D8E2_cF72_225_b_a_e_f_c 1819
458. Swedenborgite ($NaBe_4SbO_7$, $E9_2$): A4BC7D_hP26_186_ac_b_a2c_b 1756
459. Ta_2NiS_5[††]: AB5C2_oC32_63_c_c2f_f 1658
460. Ta_2NiSe_5: AB5C2_mC32_15_e_e2f_f 1570
461. Ta_2PdSe_6: AB6C2_mC18_12_a_3i_i 1533
462. Ta_3Ti_{13} (BCC SQS-16): A3B13_oC32_38_ac_a2bcdef 1596
463. Ta_3Ti_5 (BCC SQS-16): A3B5_oC32_38_abce_abcdf 1597
464. Ta_5Ti_{11} (BCC SQS-16): A5B11_mP16_6_2abc_2a3b3c 1512
465. TaTi (BCC SQS-16): AB_aP16_2_4i_4i 1506
466. $TaTi_3$ (BCC SQS-16): AB3_mC32_8_4a_12a 1516
467. $TaTi_3$ (BCC SQS-16): AB3_mC32_8_4a_4a4b ... 1517
468. $TaTi_7$ (BCC SQS-16): AB7_hR16_166_c_c2h ... 1741
469. Tellurite (β-TeO_2, $C52$): A2B_oP24_61_2c_c 1624
470. Tennantite ($Cu_{12}As_4S_{13}$): A4B24C13_cI82_217_c_deg_ag 1798
471. Tetragonal $TlFeS_2$: AB2C_tI8_119_c_e_a 1683
472. Th_7S_{12} ($D8_k$): A3B2_hP20_176_2h_ah 1751
473. ThC_2 (C_g): A2B_mC12_15_f_e 1564
474. $ThCr_2Si_2$[∂‡‡]: A2B2C_tI10_139_d_e_a 1705
475. $ThFe_2SiC$: AB2CD_oC20_63_b_f_c_c 1656
476. Thenardite [Na_2SO_4 (V), $H1_7$]: A2B4C_oF56_70_g_h_a 1662
477. Ti_5Ga_4: A4B5_hP18_193_bg_dg 1763
478. $TiBe_{12}$ (approximate, $D2_a$): A12B_hP13_191_cdei_a 1761
479. Titanite ($CaTiSiO_5$, $S0_6$): AB5CD_mC32_15_e_e2f_e_b 1571
480. Tl_2AlF_5 ($K3_3$): AB5C2_oC32_20_b_a2bc_c 1580
481. $TlAlF_4$ ($H0_8$): AB4C_tP6_123_d_eh_a 1691
482. $TlCo_2S_2$[∂‡‡]: A2B2C_tI10_139_d_e_a 1704
483. Tolbachite ($CuCl_2$): A2B_mC6_12_i_a 1527
484. Topaz ($Al_2SiO_4F_2$, $S0_5$): A2B2C4D_oP36_62_d_d_2cd_c 1628
485. Tremolite ($Ca_2Mg_5Si_8O_{22}(OH)_2$, $S4_2$): A2B2C5D24E8_mC82_12_h_i_agh_2i5j_2j 1525
486. Tutton salt [$Cu(NH_4)_2(SO_4)_2{\cdot}6H_2O$, $H4_4$]: AB20C2D14E2_mP78_14_a_10e_e_7e_e 1547
487. U_6Mn ($D2_c$): AB6_tI28_140_a_hk 1715
488. V_3AsC: ABC3_oC20_63_c_b_cf 1658
489. V_4SiSb_2: A2BC4_tI28_140_h_a_k 1712
490. $VOSO_4$: A5BC_oP28_62_3cd_c_c 1639
491. VSe_2O_6: A6B2C_tP72_103_abc5d_2d_abc 1679

[∂] $TlCo_2S_2$ and $ThCr_2Si_2$ have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

492. Vesuvianite ($Ca_{10}Al_4(Mg,Fe)_2Si_9O_{34}(OH)_4$, $S2_3$): A4B10C2D34E4F9_tP252_126_k_ce2k_f_h8k_k_d2k 1692
493. $W_2O_3(PO_4)_2$: A11B2C2_mP60_4_22a_4a_4a 1507
494. Wülfingite (ϵ-$Zn(OH)_2$, $C31$): A2B2C_oP20_19_2a_2a_a 1576
495. Wollastonite ($CaSiO_3$): AB3C_aP30_2_3i_9i_3i 1505
496. Y_2SiO_5 (RE_2SiO_5 X2): A5BC2_mC64_15_5f_f_2f 1567
497. $Zn_{22}Zr$: A22B_cF184_227_cdfg_a 1825
498. $Zn_2Mo_3O_8$: A3B8C2_hP26_186_c_ab2c_2b 1755
499. $Zn(BrO_3)_2{\cdot}6H_2O$ ($J1_{10}$)[⊗⊗]: A2B6C6D_cP60_205_c_d_d_a 1777
500. $Zn(NH_3)_2Cl_2$ ($E1_2$): A2B6C2D_oI44_74_h_ij_i_e 1666
501. $Zr_{21}Re_{25}$: A25B21_hR92_167_b2e3f_e3f 1743
502. Zr_2Al_3[**]: A3B2_oF40_43_ab_b 1603
503. Zr_3Al_2: A2B3_tP20_136_j_dfg 1701
504. $ZrFe_4Si_2$: A4B2C_tP14_136_i_g_b 1702
505. ZrNiAl: ABC_hP9_189_g_ad_f 1759
506. ZrP_2O_7 High-Temperature ($K6_1$): A7B2C_cP40_205_bd_c_a 1782
507. $ZrSe_3$: A3B_mP8_11_3e_e 1521
508. $ZrTe_5$: A5B_oC24_63_c2f_c 1655
509. Zunyite [$Al_{13}(OH,F)_{18}Si_5O_{20}Cl$, $S0_8$]: A13BC18D20E5_cF228_216_dh_b_fh_2eh_ce 1790
510. "Martensite Type" FeC_x ($x \leq 0.06$) ($L'2_0$): AB_tI4_139_b_a 1711

POSCAR Index

1. $B30$ (MgZn?): AB_oI48_44_6d_ab2cde 1605
2. $C17$ (Fe_2B) (*obsolete*): AB2_tI12_121_ab_i 1685
3. $C2$ $(Ba,Ca)CO_3$: ABC3_mC10_5_b_a_ac 1512
4. $C26_a$ (NO_2) (*obsolete*): AB2_cI36_199_b_c 1773
5. $C27$ (CdI_2) (*questionable*): AB2_hP6_186_b_ab 1757
6. $C53$ ($SrBr_2$) (*obsolete*): A2B_oP12_62_2c_c 1633
7. $D0_{10}$ (WO_3) (*obsolete*): A3B_oP16_57_a2d_d 1615
8. $D0_{13}$ ($AlCl_3$) (*obsolete*): AB3_hP4_164_b_ad 1735
9. $D0_{15}$ ($AlCl_3$) (*obsolete*): AB3_mC16_5_c_3c 1510
10. $D0_6$ (Tysonite, LaF_3) (*obsolete*): A3B_hP24_193_ack_g 1763
11. $D0_7$ (CrO_3) (*obsolete*): AB3_oC16_20_a_bc 1580
12. $D2_2$ ($MgZn_5$?) (*problematic*): AB5_mC48_12_2i_ac5i2j 1531
13. $D6_2$ (Sb_2O_4) (*obsolete*): A2B_cF96_227_abf_cd 1831
14. $D8_7$ (Shcherbinaite, V_2O_5) (*obsolete*): A5B2_oP14_31_a2b_b 1587
15. $E2_3$ ($LiIO_3$) (*obsolete*): ABC3_hP10_182_c_b_g 1755
16. $E3_1$ (β-Ag_2HgI_4) (*obsolete*): A2BC4_tP7_111_f_a_n 1680
17. $E6_1$ ($Sr(OH)_2(H_2O)_8$) (*obsolete*): A8B2C_tP11_123_r_f_a 1690
18. $E6_2$ [$SrO_2(H_2O)_8$] (*possibly obsolete*): A8B2C_tP11_123_r_h_a 1691
19. $F5_{11}$ (KNO_2) (*obsolete*): ABC2_mC8_8_a_a_b 1518
20. $F5_4$ (NH_4ClO_2) (*obsolete*): ABC2_tP8_100_b_a_c 1678
21. $F6_1$ (Chalcopyrite, $CuFeS_2$) (*obsolete*): ABC2_tP4_115_a_c_g 1683
22. $G7_3$ [Northupite, $Na_3MgCl(CO_3)_2$] (*obsolete*): A2BCD3E6_cF208_227_e_c_d_f_g 1829
23. $G7_5$ ($PbCO_3 \cdot PbCl_2$, Phosgenite) (*obsolete*): AB2C3D2_tP16_90_c_f_ce_e 1676
24. $H5_6$ [Tychite, $Na_6Mg_2SO_4(CO_3)_4$)] (*obsolete*): A4B2C6D16E_cF232_227_e_d_f_eg_a 1834
25. $H5_9$ [Autunite, $Ca(UO_2)_2(PO_4)_2{\cdot}10\frac{1}{2}H_2O$] (*obsolete*)[⊗‡‡]: AB2C2_tI10_139_a_d_e 1711
26. $H6_4$ [$Ni(NO_3)_2(NH_3)_6$] (*obsolete*)[⊗⊗]: A2B6CD6_cP60_205_c_d_a_d 1779
27. $I1_3$ ($SrCl_2{\cdot}(H_2O)_6$) (*obsolete*)[§§]: A2B6C_hP9_162_d_k_a 1730
28. $L1_a$ (disputed $CuPt_3$): AB7_cF32_225_b_ad 1819
29. $S0_4$ (Staurolite, $Fe(OH)_2Al_4Si_2O_{10}$) (*obsolete*): A4BC12D2_oC76_63_eg_c_f3gh_g 1654
30. $S3_4$ (II) (Catapleiite, $Na_2Zr(SiO_3)_3{\cdot}H_2O$) (*obsolete*): A3B2C9D3E_hP36_194_g_f_hk_h_a 1767
31. α-AgI ($B23$): A21B_cI44_229_bdh_a 1836

[⊗] Li_2CN_2 and $H5_9$ [Autunite, $Ca(UO_2)_2(PO_4)_2{\cdot}10\frac{1}{2}H_2O$] (*obsolete*) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[‡‡] $TlCo_2S_2$, $ThCr_2Si_2$, Li_2CN_2, and $H5_9$ [Autunite, $Ca(UO_2)_2(PO_4)_2{\cdot}10\frac{1}{2}H_2O$] (*obsolete*) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

[⊗⊗] $Zn(BrO_3)_2{\cdot}6H_2O$ ($J1_{10}$) and $H6_4$ [$Ni(NO_3)_2(NH_3)_6$] (*obsolete*) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

[§§] $I1_3$ ($SrCl_2{\cdot}(H_2O)_6$) (*obsolete*) and Rosiaite ($PbSb_2O_6$) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

32. α-Alum [$KAl(SO_4)_2 \cdot 12H_2O$, $H4_{13}$]: AB24CD28E2_cP224_205_a_4d_b_2c4d_c 1785
33. α-BaB_2O_4 (Low-Temperature): A2BC4_hR42_161_2b_b_4b 1729
34. α-Carnegieite ($NaAlSiO_4$, $S6_5$): ABC4D_cP28_198_a_a_ab_a 1773
35. α-$Ho_2Si_2O_7$: A2B7C2_aP44_2_4i_14i_4i 1502
36. α-ICl[§]: AB_mP16_14_2e_2e 1556
37. α-$LiIO_3$: ABC3_hP10_173_b_a_c 1749
38. α-PbO_2[×]: A2B_oP12_60_d_c 1622
39. α-Potassium Nitrate (KNO_3) I[°]: ABC3_oP20_62_c_c_cd 1648
40. α-Potassium Nitrate (KNO_3) II: ABC3_oC80_36_2ab_2ab_2a5b 1596
41. α-V_3S: AB3_tI32_121_g_f2i 1686
42. α-WO_3: A3B_tP16_130_cf_c 1701
43. α-$Zn_2V_2O_7$: A7B2C2_mC44_15_e3f_f_f 1568
44. β-Alum [$Al(NH_3CH_3)_2(SO_4)_2 \cdot 12H_2O$, $H4_{14}$]: AB2C36D2E20F2_cP252_205_a_c_6d_c_c3d_c ... 1787
45. β-Alumina (Al_2O_3, $D5_6$): A2B3_hP60_194_3fk_cdef2k 1766
46. β-Arabinose [$(CH_2O)_{20}$]: AB2C_oP80_19_5a_10a_5a 1578
47. β-B_2H_6[*]: AB3_mP16_14_e_3e 1549
48. β-BaB_2O_4 (High-Temperature): A2BC4_hR42_167_f_ac_2f 1744
49. β-Ga (*obsolete*): A_mC4_15_e 1573
50. β-Ga_2O_3: A2B3_mC20_12_2i_3i 1526
51. β-$LiIO_3$: ABC3_tP40_86_g_g_3g 1672
52. β-Potassium Nitrate (KNO_3): ABC6_hR8_166_a_b_h 1742
53. β-Si_3N_4: A4B3_hP14_176_ch_h 1751
54. δ-CuTi ($L2_a$): AB_tP2_123_a_d 1692
55. δ-Ni_3Sn_4 ($D7_a$): A3B4_mC14_12_ai_2i 1528
56. δ-WO_3: A3B_aP32_2_12i_4i 1503
57. ϵ-1,2,3,4,5,6-Hexachlorocyclohexane (C_6Cl_6): AB_mP24_14_3e_3e 1557
58. η-NiSi (B_d): AB_oP8_62_c_c 1652
59. γ-Alum [$AlNa(SO_4)_2 \cdot 12H_2O$, $H4_{15}$]: AB24CD20E2_cP192_205_a_4d_b_c3d_c 1784
60. γ-Fe_4N ($L'1_0$): A4B_cP5_221_bc_a 1805
61. γ-Ga_2O_3: A11B4_cF120_227_acdf_e 1821
62. γ-$LiIO_3$: ABC3_oP20_33_a_a_3a 1592
63. γ-Potassium Nitrate (KNO_3)[□]: ABC3_hR5_160_a_a_b 1728
64. γ-TeO_2: A2B_oP12_18_2c_c 1574
65. γ-WO_3: A3B_mP32_14_6e_2e 1542
66. γ-$Y_2Si_2O_7$[‖]: A4BC_mP24_14_4e_e_e 1544
67. ζ-Fe_2N[×]: A2B_oP12_60_d_c 1622
68. ζ-Nb_2O_5 (B-Nb_2O_5): A2B5_mC28_15_f_e2f 1562
69. y-$Y_2Si_2O_7$: A7B2C2_mP22_11_3e2f_2e_ab 1522
70. $(CdSO_4)_3 \cdot 8H_2O$ ($H4_{20}$): A3B16C20D3_mC168_15_ef_8f_10f_ef 1565
71. $(NH_4)_3AlF_6$ ($J2_1$): AB30C16D3_cF200_225_a_ej_2f_bc 1817
72. $(TiCl_4 \cdot POCl_3)_2$: A7BCD_oP80_61_7c_c_c_c 1625
73. 12-phosphotungstic acid [$H_3PW_{12}O_{40} \cdot 5H_2O$ ($H4_{16}$)]: A5B40CD12_cP116_224_cd_e3k_a_k 1810
74. Adamite [$Zn_2(AsO_4)(OH)$, $H2_7$]: ABC5D2_oP36_58_g_g_3gh_eg 1619
75. Ag_2O_3[**]: A2B3_oF40_43_b_ab 1600
76. Ag_2PbO_2: A2B2C_mC20_15_ad_f_e 1559
77. $Ag_2SO_4 \cdot 4NH_3$ ($H4_{17}$): A2B12C4D4E_tP46_114_d_3e_e_e_a 1682
78. $AgMnO_4$ ($H0_9$)[‖]: ABC4_mP24_14_e_e_4e 1554
79. $AgNO_2$ ($F5_{12}$): ABC2_oI8_44_a_a_d 1605
80. $Ag[Co(NH_3)_2(NO_2)_4]$ ($J1_9$): ABC4D2E8_tP32_126_a_b_h_e_k 1694
81. $Al_{13}Fe_4$: A13B4_mC102_12_dg8i5j_4ij 1523
82. Al_2Mo_3C: A2BC3_cP24_213_c_a_d 1790
83. $Al(PO_3)_3$ ($G5_2$): AB9C3_cI208_220_c_3e_e 1803
84. AlN (cF40): AB_cF40_216_de_ce 1798
85. AlN (cI16): AB_cI16_217_c_c 1799
86. AlN (cI24): AB_cI24_220_a_b 1804
87. $AlNbO_4$: ABC4_mC24_12_i_i_4i 1535
88. $AlPO_4$ "low cristobalite type": AB4C_oC24_20_b_2c_a 1580
89. Albite ($NaAlSi_3O_8$, $S6_8$): ABC8D3_aP26_2_i_i_8i_3i 1506
90. Alluaudite [$NaMnFe_2(PO_4)_3$]: A2BCD12E3_mC76_15_f_e_b_6f_ef 1564

[§] α-ICl and LiAs have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[×] ζ-Fe_2N and α-PbO_2 have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[°] α-Potassium Nitrate (KNO_3) I, NH_4NO_3 III ($G0_{10}$), and Aragonite ($CaCO_3$, $G0_2$) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[*] β-B_2H_6 and B_2H_6 ($P2_1/c$) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[□] $KBrO_3$ ($G0_7$) and γ-Potassium Nitrate (KNO_3) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[‖] γ-$Y_2Si_2O_7$, Ammonium Persulfate [$(NH_4)_2S_2O_8$, $K4_1$], Monazite ($LaPO_4$), and $AgMnO_4$ ($H0_9$) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

[**] Ag_2O_3 and Zr_2Al_3 have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

91. Ammonium Chlorite (NH_4ClO_2): AB4CD2_tP16_113_c_f_a_e 1681
92. Ammonium Persulfate [$(NH_4)_2S_2O_8$, $K4_1$][†‖]: AB4C_mP24_14_e_4e_e 1551
93. Analcime ($NaAlSi_2O_6{\cdot}H_2O$, $S6_1$): A2B2C3D12E4_tI184_142_f_f_be_3g_g 1717
94. Andalusite (Al_2SiO_5, $S0_2$): A2B5C_oP32_58_eg_3gh_g 1617
95. Anhydrous $KAuBr_4$: AB4C_mP24_14_ab_4e_e .. 1551
96. Anthophyllite ($Mg_5Fe_2Si_8O_{22}(OH)_2$, $S4_4$): A2B5C22D2E8_oP156_62_d_c2d_2c10d_2c_4d ... 1630
97. Apophyllite ($KCa_4Si_8O_{20}F{\cdot}8H_2O$, $S5_2$): A4BC16DE28F8_tP116_128_h_a_2i_b_g3i_i 1697
98. Aragonite ($CaCO_3$, $G0_2$)[°]: ABC3_oP20_62_c_c_cd 1649
99. Arcanite (K_2SO_4, $H1_6$): A2B4C_oP28_62_2c_2cd_c 1630
100. Archerite (KH_2PO_4): A2BC4D_oF64_43_b_a_2b_a 1602
101. Arsenopyrite (FeAsS, $E0_7$): ABC_mP12_14_e_e_e 1556
102. Atacamite ($Cu_2(OH)_3Cl$): AB2C3D3_oP36_62_c_ac_cd_cd 1644
103. Au_2Nb_3: A2B3_tI10_139_e_ae 1706
104. $AuCsCl_3$ ($K7_6$): AB3C_tI20_139_ab_eh_d 1711
105. Autunite $\{Ca[(UO_2)(PO_4)]_2(H_2O)_{11}\}$: AB22C23D2E2_oP200_62_c_11d_3c10d_d_d 1643
106. Azurite [$Cu_3(CO_3)_2(OH)_2$, $G7_4$]: A2B3C2D8_mP30_14_e_ce_e_4e 1538
107. $B_{13}C_2$ "B_4C" ($D1_g$): A13B2_hR15_166_b2h_c .. 1736
108. B_2H_6 ($P2_1/c$)[*]: AB3_mP16_14_e_3e 1549
109. B_4SrO_7: A4B7C_oP24_31_2b_a3b_a 1587
110. $BaAl_2O_4$ ($H2_8$): A2BC6_hP18_182_f_b_gh 1754
111. $BaCd_{11}$: AB11_tI48_141_a_bdi 1716
112. $BaNi(CN)_4{\cdot}4H_2O$ ($H4_{22}$): AB4C4D4E_mC56_15_e_2f_2f_2f_a 1569
113. $BaNiSn_3$: ABC3_tI10_107_a_a_ab 1680
114. Bararite (Trigonal $(NH_4)_2SiF_6$, $J1_6$)[××]: A6B2C_hP9_164_i_d_a 1733
115. Barytocalcite ($BaCa(CO_3)_2$): AB2CD6_mP20_11_e_2e_e_2e2f 1522
116. Base-centered orthorhombic $Sr_4Ru_3O_{10}$: A10B3C4_oC68_64_2dfg_ad_2d 1660
117. Bassanite [$CaSO_4(H_2O)_{0.5}$, $H4_7$]: A2B2C9D2_mC90_5_ab2c_3c_b13c_3c 1509
118. Bastnäsite [$CeF(CO_3)$]: ABCD3_hP36_190_h_g_af_hi 1761
119. $BeSO_4{\cdot}4H_2O$ ($H4_3$): AB8C8D_tI72_120_c_2i_2i_b 1684
120. Berthierite ($FeSb_2S_4$, $E3_3$): AB4C2_oP28_62_c_4c_2c 1645
121. Bertrandite ($Be_4Si_2O_7(OH)_2$, $S4_6$): A4B7C2D2_oC60_36_2b_a3b_2a_b 1595
122. Bi_2GeO_5: A2BC5_oC32_36_b_a_a2b 1594
123. $Bi_3Ru_3O_{11}$: A3B11C3_cP68_201_be_efh_g 1774
124. Bischofite ($MgCl_2{\cdot}6H_2O$, $J1_7$): A2B12CD6_mC42_12_i_2i2j_a_ij 1525
125. Blossite (α-$Cu_2V_2O_7$): A2B7C2_oF88_43_b_a3b_b 1602
126. Boric Acid (H_3BO_3, $G5_1$): AB3C3_aP28_2_2i_6i_6i 1505
127. Bromocarnallite ($KMg(H_2O)_6(Cl,Br)_3$, $E2_6$): A3B6CD_tP44_85_bcg_3g_ac_e 1669
128. Brucite [$Mg(OH)_2$][¶¶]: A2BC2_hP5_164_d_a_d .. 1733
129. $C_{19}Sc_{15}$: A19B15_tP68_114_bc4e_ac3e 1681
130. COCl: ABC_oP24_61_c_c_c 1628
131. Ca_2RuO_4: A2B4C_oP28_61_c_2c_a 1624
132. Ca_2UO_5: A2B5C_mP32_14_2e_5e_ab 1540
133. $Ca_3Al_2(OH)_{12}$ ($J2_3$): A2B3C12D12_cI232_230_a_c_h_h 1837
134. Ca_3UO_6: A3B6C_mP20_4_3a_6a_a 1509
135. CaB_2O_4 (III): A2BC4_oP84_33_6a_3a_12a 1590
136. CaB_2O_4 (IV): A2BC4_cP84_205_d_ac_2d 1780
137. CaB_2O_4 I ($E3_2$): A2BC4_oP28_60_d_c_2d 1621
138. $CaBe_2Ge_2$: A2BC2_tP10_129_ac_c_bc 1698
139. CaC_2-I ($C11_a$): A2B_tI6_139_e_a 1706
140. CaC_2-III: A2B_mC12_12_2i_i 1527
141. $CaCu_4P_2$[°°]: AB4C2_hR7_166_a_2c_c 1740
142. $CaO_2(H_2O)_8$: AB8C2_tP22_124_a_n_h 1692
143. $CaSi_2$ ($C12$): AB2_hR6_166_c_2c 1739
144. $CaUO_4$: AB4C_hR6_166_b_2c_a 1741
145. Calaverite ($AuTe_2$): AB2_mP12_7_2a_4a 1515
146. Calciborite (CaB_2O_4 II): A2BC4_oP56_56_2e_e_4e 1614
147. Carnallite [$Mg(H_2O)_6KCl_3$]: A3B12CDE6_oP276_52_d4e_18e_ce_de_2d8e 1608
148. Catapleiite ($Na_2ZrSi_3O_9{\cdot}2H_2O$): A2B3C9D3E_mC144_15_2f_bcdef_9f_3f_ae 1560

[†]Ammonium Persulfate [$(NH_4)_2S_2O_8$, $K4_1$] and Monazite ($LaPO_4$) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[××]$K_2Pt(SCN)_6$ ($H6_3$), Bararite (Trigonal $(NH_4)_2SiF_6$, $J1_6$), and K_2GeF_6 ($J1_{13}$) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

[¶¶]Ce_2O_2S and Brucite [$Mg(OH)_2$] have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

[°°]$MnBi_2Te_4$ and $CaCu_4P_2$ have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

149. Cd_3As_2: A2B3_tI160_142_deg_3g 1718
150. Cd(OH)Cl ($E0_3$): ABCD_hP8_186_b_b_a_a 1758
151. Ce_2O_2S[¶¶]: A2B2C_hP5_164_d_d_a 1732
152. $CeCu_2$: AB2_oI12_74_e_h 1667
153. Cervantite (α-Sb_2O_4): A2B_oP24_33_4a_2a 1591
154. Chabazite ($Ca_{1.4}Sr_{0.3}Al_{3.8}Si_{8.3}O_{24}{\cdot}13H_2O$, $S3_4$ (I)): A5B21C24D12_hR62_166_a2c_ehi_fg2h_i 1738
155. Chalcanthite ($CuSO_4{\cdot}5H_2O$, $H4_{10}$): AB10C9D_aP42_2_ae_10i_9i_i 1504
156. Chalcocyanite ($CuSO_4$): AB4C_oP24_62_a_2cd_c 1646
157. Chiolite ($Na_5Al_3F_{14}$, $K7_5$): A3B14C5_tP44_128_ac_ehi_bg 1696
158. Chlorite [$Mg_3(Mg_2Al)(Si_3Al)O_{10}(OH)_8$, $S5_5$]: A3B5C4D2_mC112_15_a3ef_5f_4f_2f 1566
159. Chrysotile ($H_4Mg_3Si_2O_9$, $S4_5$): AB6C11D6E4_mC112_12_e_gi2j_i5j_2i2j_2j 1532
160. Chrysotile ($Mg_3Si_2O_5(OH)_4$): A3B5C4D2_mC56_9_3a_5a_4a_2a 1519
161. Clinocervantite (β-Sb_2O_4): A2B_mC24_15_2f_ce 1565
162. Co_2Al_9 ($D8_d$): A9B2_mP22_14_a4e_e 1547
163. $Co_2B_2O_5$: A2B2C5_aP18_2_2i_2i_5i 1501
164. $Co_3(SeO_3)_3{\cdot}H_2O$: A3B2C10D3_aP36_2_ah2i_2i_10i_3i 1503
165. Co_9S_8 ($D8_9$): A9B8_cF68_225_af_ce 1816
166. Colquiriite ($LiCaAlF_6$): ABC6D_hP18_163_d_b_i_c 1731
167. Columbite ($FeNb_2O_6$, $E5_1$): AB2C6_oP36_60_c_d_3d 1624
168. Copper (I) Azide (CuN_3): AB3_tI32_88_d_cf 1675
169. Copper (II) Azide [$Cu(N_3)_2$]: AB6_oP28_62_c_6c 1647
170. Cr_5O_{12}: A5B12_oP68_60_c2d_6d 1623
171. Cr-233 Quasi-One-Dimensional Superconductor ($K_2Cr_3As_3$): A3B3C2_hP16_187_jk_jk_ck 1759
172. $CrCl_3(H_2O)_6$ ($J2_2$): A3BC6_hR20_167_e_b_f ... 1745
173. Crancrinite ($Na_6Ca_2Al_6Si_6O_{24}(CO_3)_2$, $S3_3$ (I)): A3BCD3E15F3_hP52_173_c_b_b_c_5c_c 1748
174. Cronstedtite {Fe(Fe,Si)[$(OH)_2$,O]O_3, $S5_7$}: AB3C2D_hR7_160_a_b_2a_a 1727
175. Cryolite (Na_3AlF_6, $J2_6$): AB6C3_mP20_14_a_3e_de 1553
176. $Cs_{11}O_3$: A11B3_mP56_14_11e_3e 1537
177. Cs_2Sb: A2B_oP24_62_4c_2c 1634
178. Cs_2Se: A2B_oF24_43_b_a 1603
179. $Cs_3As_2Cl_9$ ($K7_3$): A2B9C3_hP14_150_d_eg_ad .. 1723
180. Cs_3CoCl_5 ($K3_1$): A5BC3_tI36_140_cl_b_ah 1715
181. $Cs_3Cr_2Cl_9$: A9B2C3_hP28_194_hk_f_bf 1768
182. $Cs_3Tl_2Cl_9$ ($K7_2$): A9B3C2_hR28_167_ef_e_c 1747
183. $Cs_6W_{11}O_{36}$: A6B36C11_mC212_9_6a_36a_11a .. 1520
184. Cs_7O: A7B_hP24_187_ai2j2kn_j 1759
185. CsB_4O_6F: A4BCD6_oP48_33_4a_a_a_6a 1591
186. $CsFeS_2$ (100 K): ABC2_oI16_71_g_i_eh 1665
187. CsO: AB_oI8_71_g_i 1665
188. $CsSO_3$ ($K1_2$): AB3C_hP20_190_ac_i_f 1760
189. $Cu_2Pb(SeO_3)_2Br_2$: A2B2C6DE2_oC52_63_g_e_fh_c_f 1652
190. $Cu_3[Fe(CN)_6]_2{\cdot}xH_2O$ ($J2_5$, $x \approx 3$): A6B9CD2E6_cF96_225_e_bf_a_c_e 1814
191. Cu(OH)Cl: ABCD_mP16_14_e_e_e_e 1555
192. Cubic Cu_2OSeO_3: A2B4C_cP56_198_ab_2a2b_2a 1771
193. Cubic CuPt ($L1_3$ (I), $D4$): AB_cF32_227_c_d ... 1835
194. Danburite ($CaB_2Si_2O_8$, $S6_3$): A2BC8D2_oP52_62_d_c_2c3d_d 1633
195. Diamminetriamidodizinc Chloride ($[Zn_2(NH_3)_2(NH_2)_3]Cl$): AB12C5D2_oP40_18_a_6c_b2c_c 1575
196. Diaspore (AlOOH, $E0_2$): ABC2_oP16_62_c_c_2c 1648
197. Diopside [$CaMg(SiO_3)_2$, $S4_1$]: ABC6D2_mC40_15_e_e_3f_f 1573
198. Dodecatungstophosphoric Acid Hexahydrate [$H_3PW_{12}O_{40}{\cdot}6H_2O$]: A27B52CD12_cP184_224_dl_eh3k_a_k 1807
199. Dolomite [$MgCa(CO_3)_2$, $G1_1$]: A2BCD6_hR10_148_c_a_b_f 1721
200. Double Perovskite (Ba_2MnWO_6): A2BC6D_cF40_225_c_a_e_b 1813
201. Enstatite ($MgSiO_3$, $S4_3$): AB3C_oP80_61_2c_6c_2c 1627
202. Epididymite ($BeHNaO_8Si_3$, $S4_7$): ABCD8E3_oP112_62_d_2c_d_4c6d_3d 1650
203. Eriochalcite ($CuCl_2{\cdot}2H_2O$, $C45$): A2BC4D2_oP18_53_h_a_i_e 1610
204. $EuIn_2P_2$: AB2C2_hP10_194_a_f_f 1769
205. Eudidymite ($BeHNaO_8Si_3$): A2B4C2D17E6_mC124_15_f_2f_f_e8f_3f 1561
206. Eulytine ($Bi_4(SiO_4)_3$, $S1_5$): A4B12C3_cI76_220_c_e_a 1801
207. $Fe_2(CO)_9$ ($F4_1$): A9B2C9_hP40_176_hi_f_hi 1753
208. Fe_2N (approximate, $L'3_0$)[∂∂]: AB_hP4_194_c_a .. 1771
209. Fe_3PO_7: A3B7C_hR11_160_b_a2b_a 1726
210. Fe_8N ($D2_g$): A8B_tI18_139_deh_a 1710
211. FeF_3 ($D0_{12}$): A3B_hR8_167_e_b 1746
212. Ferroelectric $NH_4H_2PO_4$: A6BC4D_oP48_19_6a_a_4a_a 1577

[∂∂] $LiZn_2$ (C_k) and Fe_2N (approximate, $L'3_0$) have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

213. Ferroelectric $NaNO_2$ ($F5_5$): ABC2_oI8_44_a_a_c 1604
214. Fluorapatite [$Ca_5F(PO_4)_3$, $H5_7$]: A5BC12D3_hP42_176_fh_a_2hi_h 1752
215. Ga_2Mg_5 ($D8_g$): A2B5_oI28_72_j_bfj 1666
216. $GaMo_4S_8$: AB4C8_cF52_216_a_e_2e 1796
217. Gd_2SiO_5 (RE_2SiO_5 X1): A2B5C_mP32_14_2e_5e_e 1541
218. Gwihabaite [NH_4NO_3 (V)]: A4B2C3_tP72_77_8d_ab2c2d_6d 1668
219. Gypsum ($CaSO_4{\cdot}2H_2O$, $H4_6$): AB4C6D_mC48_15_e_2f_3f_e 1570
220. $H_3PW_{12}O_{40}{\cdot}29H_2O$ ($H4_{21}$): A29B40CD12_cF656_227_ae2fg_e3g_b_g 1827
221. $H_3PW_{12}O_{40}{\cdot}3H_2O$: A3B40CD12_cP112_224_d_e3k_a_k 1809
222. Hambergite [$Be_2BO_3(OH)$, $G7_2$]: AB2CD4_oP64_61_c_2c_c_4c 1626
223. Hauyne [$(Na_{0.5}Ca_{0.3}K_{0.2})_8(Al_6Si_6O_{24})(SO_4)_{1.5}$, $S6_9$]: A3B4C4D4E16F4G3_cP76_218_c_e_e_e_ei_e_d .. 1800
224. Hemimorphite ($Zn_4Si_2O_7(OH)_2{\cdot}H_2O$, $S2_2$): A2B5CD2_oI40_44_2c_abcde_d_e 1604
225. Hexagonal Delafossite ($CuAlO_2$): ABC2_hP8_194_a_c_f 1770
226. Hexagonal WO_3: A3B_hP12_191_gl_f 1762
227. Hg_2O_2NaI: A2BCD2_hP18_180_f_c_b_i 1754
228. Hg_2TiCu Inverse Heusler: AB2C_cF16_216_b_ad_c 1795
229. $HgCl_2{\cdot}2HgO$: A2B3C2_mP14_14_e_ae_e 1538
230. High-Temperature Cryolite (Na_3AlF_6): AB6C3_oI20_71_a_in_cj 1664
231. High-Temperature Cubic $KClO_4$ ($H0_5$): ABC4_cF24_216_b_a_e 1797
232. High-Temperature Mo_8O_{23}: A8B23_mP62_13_4g_c11g 1536
233. $HoMn_2O_5$: AB2C5_oP32_55_g_fh_eghi 1613
234. $HoSb_2$: AB2_oC6_21_a_k 1581
235. Huanzalaite ($MgWO_4$, $H0_6$): AB4C_mP12_13_f_2g_e 1536
236. In_4Se_3: A4B3_oP28_58_4g_3g 1618
237. InS: AB_oP8_58_g_g 1619
238. Jacutingaite (Pt_2HgSe_3): AB2C3_hP12_164_d_ae_i 1734
239. $K_2CuCl_4{\cdot}2H_2O$ ($H4_1$): A4BC4D2E2_tP26_136_fg_a_j_d_e 1703
240. K_2GeF_6 ($J1_{13}$)[××]: A6BC2_hP9_164_i_a_d 1734
241. $K_2HgCl_4{\cdot}H_2O$ ($E3_4$): A4BCD2_oP32_55_ghi_f_e_gh 1612
242. K_2NbF_7 ($K6_2$): A7B2C_mP40_14_7e_2e_e 1545
243. $K_2Ni(CN)_4$: A4B2C4D_mP22_14_2e_e_2e_a 1543
244. K_2NiF_4: A4B2C_tI14_139_ce_e_a 1707
245. $K_2OsO_2Cl_4$ ($J1_5$): A4B2C2D_tI18_139_h_d_e_a . 1707
246. $K_2Pt(SCN)_6$ ($H6_3$)[××]: A2BC6_hP9_164_d_a_i .. 1733
247. $K_2Pt(SCN)_6{\cdot}2H_2O$: A6B4C2D6E2FG6_mP54_14_3e_2e_e_3e_e_a_3e . 1544
248. K_2PtCl_4 ($H1_5$): A4B2C_tP7_123_j_e_a 1690
249. $K_2S_2O_5$ ($K0_1$): A2B5C2_mP18_11_2e_e2f_2e ... 1521
250. $K_2S_3O_6$ ($K5_1$): A2B6C3_oP44_62_2c_2c2d_3c .. 1632
251. $K_2Sn(OH)_6$ ($H6_2$): A6B2C6D_hR15_148_f_c_f_a 1721
252. $K_2SnCl_4{\cdot}H_2O$: A4BC2D_oP32_62_2cd_c_d_c ... 1639
253. $K_2SnCl_4{\cdot}H_2O$ ($E3_5$): A4BC2D_oP32_62_2cd_b_2c_a 1638
254. $K_2Ti_2O_5$: A2B5C2_mC18_12_i_a2i_i 1526
255. $K_3Co(NO_2)_6$ ($J2_4$): AB3C6D12_cF88_202_a_bc_e_h 1775
256. K_3CrO_8: AB3C8_tI24_121_a_bd_2i 1686
257. $K_3TlCl_6{\cdot}2H_2O$ ($J3_1$): A6B2C3D_tI168_139_egikl2m_ejn_bh2n_acf 1708
258. $K_3W_2Cl_9$ ($K7_1$): A9B3C2_hP28_176_hi_af_f 1753
259. $K_4[Mo(CN)_8]{\cdot}2H_2O$ ($F2_1$): A8B4C4DE8F2_oP108_62_4c2d_2d_2cd_c_4c2d_d 1641
260. K(SH) ($B22$): AB_hR2_166_a_b 1743
261. $KAuBr_4{\cdot}2H_2O$ ($H4_{19}$): AB4C2D_mP32_14_e_4e_2e_e 1550
262. $KBe_2BO_3F_2$: AB2C2DE3_hR9_155_b_c_c_a_e .. 1725
263. $KBrO_3$ ($G0_7$)[□]: ABC3_hR5_160_a_a_b 1728
264. $KFeS_2$ ($F5_a$): ABC2_mC16_15_e_e_f 1572
265. KH_2PO_4 ($H2_2$): A4BC4D_tI40_122_e_b_e_a 1688
266. KHF_2 ($F5_2$): A2BC_tI16_140_h_d_a 1713
267. $KICl_4{\cdot}H_2O$ ($H0_{10}$): A4BCD_mP28_14_4e_e_e_e 1543
268. KNO_2 III[‡]: ABC2_mP16_14_e_e_2e 1553
269. KSO_3 ($K1_1$): AB3C_hP30_150_ef_3g_c2d 1724
270. Kesterite [$Cu_2(Zn,Fe)SnS_4$]: A2BCD4_tI16_82_ac_b_d_g 1668
271. Kotoite ($Mg_3(BO_3)_2$): A2B3C6_oP22_58_g_af_gh 1616
272. Kyanite (Al_2SiO_5, $S0_1$): A2B5C_aP32_2_4i_10i_2i 1501
273. La_3BWO_9 ($P3$): AB3C9D_hP28_143_2a_2d_6d_bc 1719
274. La_3BWO_9 ($P6_3$): AB3C9D_hP28_173_a_c_3c_b . 1749
275. La_3CuSiS_7: AB3C7D_hP24_173_a_c_b2c_b 1748
276. $LaFe_4P_{12}$: A4BC12_cI34_204_c_a_g 1776
277. LaH_{10} High-T_c Superconductor: A10B_cF44_225_cf_b 1812
278. LaOAgS: ABCD_tP8_129_b_c_a_c 1699
279. Lepidocrocite (γ-FeO(OH), $E0_4$): AB2C2_oC20_63_c_f_2c 1655
280. Li_2CN_2[⊗‡‡]: AB2C2_tI10_139_a_d_e 1710
281. Li_2PrO_3: A2B3C_oC12_65_h_bh_a 1661

[‡] KNO_2 III and Manganite (γ-MnO(OH), $E0_6$) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

282. $Li_2SO_4{\cdot}H_2O$ ($H4_8$): A2B2C5D_mP20_4_2a_2a_5a_a 1508
283. Li_7TaO_6: A8B6C_hR15_148_cf_f_a 1722
284. LiAs[§]: AB_mP16_14_2e_2e 1557
285. $LiClO_4{\cdot}3H_2O$ ($H4_{18}$): AB6CD7_hP30_186_b_d_a_b2c 1758
286. $LiCuVO_4$: ABC4D_oI28_74_a_d_hi_e 1667
287. $LiGaO_2$: ABC2_oP16_33_a_a_2a 1592
288. $LiKSO_4$ ($H1_4$): ABC4D_hP14_173_a_b_bc_b 1750
289. $LiNb_6O_{15}F$: ABC6D15_oP46_51_f_d_2e2i_aef4i2j 1607
290. $LiOH{\cdot}H_2O$ ($B36$): A3BC2_mC24_12_ij_h_gi 1528
291. $LiZn_2$ (C_k)[∂∂]: AB_hP4_194_a_c 1771
292. Low-Temperature $(NH_3CH_3)Al(SO_4)_2{\cdot}12H_2O$: ABC30DE20F2_oP220_29_a_a_30a_a_20a_2a 1583
293. Low-Temperature $GaMo_4S_8$: AB4C8_hR13_160_a_ab_2a2b 1728
294. Low-Temperature Mo_8O_{23}: A8B23_mP124_7_16a_46a 1514
295. Lu_2CoGa_3: AB3C2_hP24_194_f_k_bh 1769
296. Lueshite ($NaNbO_3$): ABC3_oP40_57_cd_e_cd2e 1616
297. Maghemite (γ-Fe_2O_3, $D5_7$): A2B3_cP60_212_bcd_ace 1789
298. Magnetoplumbite ($PbFe_{12}O_{19}$): A12B19C_hP64_194_ab2fk_efh2k_d 1764
299. Manganese-leonite 110 K $[K_2Mn(SO_4)_2{\cdot}4H_2O]$: A8B2CD12E2_mP100_14_8e_2e_ad_12e_2e 1546
300. Manganese-leonite 185 K $[K_2Mn(SO_4)_2{\cdot}4H_2O]$: A8B2CD12E2_mC200_15_8f_2f_ce_2e11f_2f 1568
301. Manganese-leonite $[K_2Mn(SO_4)_2{\cdot}4H_2O$, $H4_{23}]$: A8B2CD15E2_mC112_12_2i3j_j_ad_g4i5j_2i 1529
302. Manganite (γ-MnO(OH), $E0_6$)[‡]: ABC2_mP16_14_e_e_2e 1554
303. Marialite Scapolite $[Na_4Cl(AlSi_3)_3O_{24}$, $S6_4]$: AB4C24D12_tI82_87_a_h_2h2i_hi 1673
304. Mayenite ($12CaO{\cdot}7Al_2O_3$, $K7_4$, C12A7): A7B12C19_cI152_220_bc_2d_ace 1802
305. Mercury (II) Azide $[Hg(N_3)_2]$: AB6_oP28_29_a_6a 1582
306. Mercury Cyanide $[Hg(CN)_2$, $F1_1]$: A2BC2_tI40_122_e_d_e 1687
307. Meta-autunite (I) $[Ca(UO_2)_2(PO_4)_2{\cdot}6H_2O$, $H5_{10}]$: AB4C6DE_tP26_129_c_j_2ci_a_c 1698
308. Mg_2Cu (C_b): AB2_oF48_70_g_fg 1663
309. $Mg_3Cr_2Al_{18}$: A18B2C3_cF184_227_fg_d_ac 1824
310. Mg_3P_2 ($D5_5$): A3B2_cP10_224_d_b 1808
311. Mg_3Ru_2: A3B2_cP20_213_d_c 1790
312. $Mg(ClO_4)_2{\cdot}6H_2O$ ($H4_{11}$): A2B6CD8_oP34_31_2a_2a2b_a_4a2b 1586
313. $Mg(NH_3)_2Cl_2$ ($E1_3$): A2B8CD2_oC26_65_h_r_a_i 1661
314. $MgCuAl_2$ ($E1_a$): A2BC_oC16_63_f_c_c 1653
315. Mn_3As ($D0_d$): AB3_oC16_63_c_3c 1657
316. $MnBi_2Te_4$[°°]: A2BC4_hR7_166_c_a_2c 1737
317. MnCuP: ABC_oP12_62_c_c_c 1651
318. $MnF_{2-x}(OH)_x$: A2BC2_oP10_34_c_a_c 1593
319. $MnPS_3$: ABC3_mC20_12_g_i_ij 1534
320. $Mo_{17}O_{47}$: A17B47_oP128_32_a8c_a23c 1588
321. Mo_4P_3: A4B3_oP56_62_8c_6c 1638
322. MoP_2: AB2_oC12_36_a_2a 1596
323. $MoPO_5$: AB5C_tP14_85_c_cg_b 1670
324. Monazite ($LaPO_4$)[†‖]: AB4C_mP24_14_e_4e_e 1552
325. Monoclinic Co_4Al_{13}: A13B4_mC102_8_17a11b_8a2b 1516
326. Monoclinic Cu_2OSeO_3: A2B4C_mP28_14_abe_4e_e 1539
327. Monoclinic $FeTlSe_2$: AB2C_mC16_12_g_2i_i 1530
328. Morenosite ($NiSO_4{\cdot}7H_2O$, $H4_{12}$): A14BC11D_oP108_19_14a_a_11a_a 1576
329. Muratatite $[(Y,Na)_6(Zn,Fe)_5Ti_{12}O_{29}(O,F)_{10}F_4]$: A16B40C12D6E5_cF316_216_eh_e2g2h_h_f_be 1792
330. Muscovite ($KH_2Al_3Si_3O_{12}$, $S5_1$): A2BC10D2E4_mC76_15_f_e_5f_f_2f 1562
331. NH_4Br ($B25$): AB4C_tP12_129_c_i_a 1699
332. NH_4CdCl_3 ($E2_4$): AB3C_oP20_62_c_3c_c 1645
333. NH_4ClBrI ($F5_{14}$): ABCD_oP16_62_c_c_c_c 1651
334. $NH_4H_2PO_2$ ($F5_7$): A2BC2D_oC24_67_m_a_n_g 1662
335. $NH_4H_2PO_4$: A8BC4D_tI56_122_2e_b_e_a 1688
336. NH_4HF_2 ($F5_8$): A2BC_oP16_53_eh_ab_g 1610
337. NH_4HgCl_3 ($E2_5$): A3BC_tP5_123_cg_a_d 1689
338. NH_4I_3 ($D0_{16}$): A3B_oP16_62_3c_c 1636
339. NH_4NO_3 I ($G0_8$): AB_cP2_221_a_b 1806
340. NH_4NO_3 II ($G0_9$): ABC3_tP10_100_b_a_bc 1679
341. NH_4NO_3 III ($G0_{10}$)[°]: ABC3_oP20_62_c_c_cd 1649
342. NH_4NO_3 IV ($G0_{11}$): A4B2C3_oP18_59_ef_ab_af 1620
343. $NH_4Pb_2Br_5$ ($K3_4$): A5BC2_tI32_140_bl_a_h 1714
344. NO_2 (Modern, $C26$): AB2_cI36_204_d_g 1777
345. $Na_{0.74}CoO_2$: AB2C2_hP10_194_a_bc_f 1768
346. $Na_2Ca_6Si_4O_{15}$: A6B2C15D4_mP54_7_6a_2a_15a_4a 1513
347. Na_2CaSiO_4 ($S6_6$): AB2C4D_cP32_198_a_2a_ab_a 1772
348. Na_2CrO_4 ($H1_8$): AB2C4_oC28_63_c_bc_fg 1656
349. $Na_2Mo_2O_7$: A2B2C7_oC88_64_ef_df_3f2g 1660
350. Na_2PrO_3: A2B3C_mC48_15_aef_3f_2e 1560
351. Na_2SO_3 ($G3_2$): A2B3C_hP12_147_abd_g_d 1720
352. $Na_4Ge_9O_{20}$: A9B4C20_tI132_88_a2f_f_5f 1674
353. $NaAlCl_4$: AB4C_oP24_19_a_4a_a 1579
354. $NaC_5H_{11}O_8S$: A5B11CD8E_aP26_1_5a_11a_a_8a_a 1499
355. $NaCr(SO_4)_2{\cdot}12H_2O$ Alum: AB12CD8E2_cP96_205_a_2d_b_cd_c 1783
356. $NaMn_7O_{12}$: A7BC12_cI40_204_bc_a_g 1777

357. $NaNb_6O_{15}F$: ABC6D15_oC46_38_b_b_2a2d_2ab4d2e 1598
358. $NaNbO_3$: ABC3_oP40_17_abcd_2e_abcd4e 1574
359. NaP: AB_oP16_19_2a_2a 1579
360. NaS_2: AB2_tI48_122_cd_2e 1689
361. $NaSb(OH)_6$ ($J1_{11}$): AB6C_tP32_86_d_3g_c 1671
362. $NaSbF_4(OH)_2$ ($J1_{12}$): A6BC_hP16_163_i_b_c ... 1731
363. $NaSbF_6$: A6BC_cP32_205_d_b_a 1782
364. Nacrite [$Al_2Si_2O_5(OH)_4$, $S5_4$]: A2B4C9D2_mC68_9_2a_4a_9a_2a 1518
365. Nahcolite ($NaHCO_3$, $G0_{12}$): ABCD3_mP24_14_e_e_e_3e 1555
366. Natrolite ($Na_2Al_2Si_3O_{10}{\cdot}2H_2O$, $S6_{10}$): A2B4C2D12E3_oF184_43_b_2b_b_6b_ab 1601
367. $Nb_2Pd_3Se_8$: A2B3C8_oP26_55_h_ag_2g2h 1611
368. $Nb_2Zr_6O_{17}$: A2B17C6_oI100_46_ab_b8c_3c 1606
369. Nb_3O_7F: A3B8_oC22_65_ag_bd2gh 1662
370. $NbAs_2$: A2B_mC12_5_2c_c 1510
371. $NbTe_2$: AB2_mC18_12_ai_3i 1531
372. $Nd_2Fe_{14}B$: AB14C2_tP68_136_f_ce2j2k_fg 1703
373. $Nd_4Re_2O_{11}$: A4B11C2_tP68_86_2g_ab5g_g 1671
374. $Nd(BrO_3)_3{\cdot}9H_2O$ ($G2_2$): A3B9CD9_hP44_186_c_3c_b_cd 1756
375. Nevskite (BiSe): AB_hP12_164_c2d_c2d 1735
376. Ni_3Si_2: A3B2_oC80_36_4a4b_2a3b 1594
377. $Ni(H_2O)_6SnCl_6$ ($I6_1$): A6B6CD_hR14_148_f_f_b_a 1722
378. $Ni(NO_3)_2(H_2O)_6$: A12B2CD12_aP54_2_12i_2i_i_12i 1500
379. Norbergite [$Mg(F,OH)_2 \cdot Mg_2SiO_4$, $S0_7$]: A2B3C4D_oP40_62_d_cd_2cd_c 1629
380. O(OH)Y: ABC_mP6_11_e_e_e 1523
381. Original β-WO_3 (*obsolete*): A3B_oP32_62_ab4c_2c 1636
382. Orpiment (As_2S_3, $D5_f$): A2B3_mP20_14_2e_3e . 1539
383. Orthorhombic Co_4Al_{13}: A13B4_oP102_31_17a11b_8a2b 1585
384. Orthorhombic CrO_3: AB3_oC16_40_b_a2b 1599
385. Orthorhombic $Sr_4Ru_3O_{10}$: A10B3C4_oP68_55_2e2fgh2i_adef_2e2f 1611
386. Os_4Al_{13}: A13B4_mC34_12_b6i_2i 1524
387. P_4Se_3: A4B3_oP112_62_8c4d_4c4d 1637
388. $PNCl_2$ ($E1_4$): A2BC_tP32_86_2g_g_g 1670
389. Paralstonite ($BaCa(CO_3)_2$): AB2CD6_hP30_150_e_c2d_f_3g 1724
390. Pararealgar (AsS)[¶]: AB_mP32_14_4e_4e 1558
391. Paratellurite (α-TeO_2): A2B_tP12_92_b_a 1677
392. Parawollastonite ($CaSiO_3$, $S3_3$ (II)): AB3C_mP60_14_3e_9e_3e 1548
393. Parkerite ($Ni_3Bi_2S_2$): AB2C_oP8_51_e_be_f 1606
394. $Pb(NO_3)_2$ ($G2_1$): A2B6C_cP36_205_c_d_a 1779
395. Pd_5Pu_3: A5B3_oC32_63_cfg_ce 1654
396. $Pd(NH_3)_4Cl_2{\cdot}H_2O$ ($H4_9$): A2BC4D_tP16_127_h_d_i_a 1695
397. Phase II $Cd_2Re_2O_7$: A2B7C2_tI44_119_i_bdefgh_i 1683
398. Phase III $Cd_2Re_2O_7$: A2B7C2_tI44_98_f_bcde_f 1678
399. Phosgenite [$Pb_2Cl_2(CO_3)$]: AB2C3D2_tP32_127_g_eh_gk_k 1696
400. Possible δ-$Gd_2Si_2O_7$: A2B7C2_oP44_33_2a_7a_2a 1589
401. Possible δ-$Y_2Si_2O_7$: A7B2C2_oP44_62_3c2d_2c_d 1640
402. Predicted High-Pressure $YCaH_{12}$: AB12C_cP14_221_a_h_b 1805
403. Predicted Li_2MgH_{16} 300 GPa: A16B2C_hP19_164_2d2i_d_b 1732
404. Predicted Li_2MgH_{16} High-Temperature Superconductor (250 GPa): A16B2C_cF152_227_eg_d_a 1823
405. Predicted Phase IV $Cd_2Re_2O_7$: A2B7C2_oF88_22_k_bdefghij_k 1582
406. Proposed 300 GPa HfH_{10}: A10B_hP22_194_bhj_c 1764
407. Protoanthophyllite ($H_2Mg_7Si_8O_{24}$): A2B7C24D8_oP82_58_g_ae2f_2g5h_2h 1617
408. Pseudobrookite (Fe_2TiO_5, $E4_1$)[††]: A2B5C_oC32_63_f_c2f_c 1653
409. Pt_2Sn_3 ($D5_b$): A2B3_hP10_194_f_bf 1765
410. $Pu_{31}Rh_{20}$: A31B20_tI204_140_b2gh3m_ac2fh3l .. 1713
411. Pyrophyllite [$AlSi_2O_5(OH)$, $S5_6$]: AB5CD2_mC72_15_f_5f_f_2f 1571
412. $Rb_2C_2O_4{\cdot}H_2O$: A2BC4D2_mC36_15_f_e_2f_f 1563
413. $Rb_2CaCu_6(PO_4)_4O_2$: AB6C18D4E2_mC62_5_a_2b2c_9c_2c_c 1511
414. $Rb_2Mo_2O_7$: A2B7C2_oC88_40_abc_2b6c_a3b ... 1598
415. $RbNO_3$ (IV): AB3C_hP45_144_3a_9a_3a 1719
416. Re_3B: AB3_oC16_63_c_cf 1657
417. ReB_3: A3B_hP8_194_af_c 1767
418. Realgar (AsS, B_l)[¶]: AB_mP32_14_4e_4e 1558
419. Retgersite (α-$NiSO_4{\cdot}6H_2O$, $H4_5$): A12BC10D_tP96_92_6b_a_5b_a 1676

[¶]Pararealgar (AsS) and Realgar (AsS, B_l) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

[††]Pseudobrookite (Fe_2TiO_5, $E4_1$) and Ta_2NiS_5 have similar AFLOW prototype labels (*i.e.*, same symmetry and set of Wyckoff positions with different stoichiometry labels due to alphabetic ordering of atomic species). They are generated by the same symmetry operations with different sets of parameters.

420. $Rh_{20}Si_{13}$: A10B7_hP34_176_c3h_b2h 1750
421. $RhCl_2(NH_3)_5Cl$ ($J1_8$): A3B15C5D_oP96_62_cd_3c6d_3cd_c 1635
422. Rhombohedral $CuTi_2S_4$: AB4C2_hR28_166_2c_2c2h_abh 1739
423. Rhombohedral Delafossite ($CuFeO_2$): ABC2_hR4_166_a_b_c 1742
424. Rinneite ($K_3NaFeCl_6$): A6BC3D_hR22_167_f_b_e_a 1746
425. Rosiaite ($PbSb_2O_6$)[§§]: A6BC2_hP9_162_k_a_d .. 1730
426. $Ru_{11}B_8$: A8B11_oP38_55_g3h_a3g2h 1613
427. RuB_2: A2B_oP6_59_f_a 1620
428. Rynersonite (Orthorhombic $CaTa_2O_6$): AB6C2_oP36_62_c_2c2d_d 1646
429. Sanguite ($KCuCl_3$): A3BC_mP20_14_3e_e_e 1541
430. Sanidine ($KAlSi_3O_8$, $S6_7$): AB8C4_mC52_12_i_gi3j_2j 1533
431. Santite ($KB_5O_8{\cdot}4H_2O$, $K3_5$): A5B8CD12_oC104_41_a2b_4b_a_6b 1600
432. $Sb_4O_5Cl_2$: A2B5C4_mP22_14_e_c2e_2e 1540
433. $SbCl_5{\cdot}POCl_3$: A8BCD_oP44_62_4c2d_c_c_c 1642
434. SbI_3S_{24}: A3B24C_hR28_160_b_2b3c_a 1726
435. Scheelite ($CaWO_4$, $H0_4$): AB4C_tI24_88_b_f_a . 1675
436. Senarmontite (Sb_2O_3, $D6_1$): A3B2_cF80_227_f_e 1832
437. Shandite ($Ni_3Pb_2S_2$): A3B2C2_hR7_166_d_ab_c 1737
438. Shcherbinaite (V_2O_5) (*revised*): A5B2_oP14_59_a2f_f 1621
439. Si_{24} Clathrate: A_oC24_63_3f 1659
440. Si_2N_2O: A2BC2_oC20_36_b_a_b 1593
441. SiAs: AB_mC24_12_3i_3i 1535
442. Sillimanite (Al_2SiO_5, $S0_3$): A2B5C_oP32_62_bc_3cd_c 1632
443. $Sm_{11}Cd_{45}$: A45B11_cF448_216_bd4efg5h_ac2eh . 1794
444. SnI_4 ($D1_1$): A4B_cP40_205_cd_c 1781
445. Sodalite [$Na_4(AlSiO_4)_3Cl$, $S6_2$]: A3BC4D12E3_cP46_218_d_a_e_i_c 1800
446. Sr_2MnTeO_6: AB6C2D_mP20_14_a_3e_e_d 1552
447. Sr_2NiTeO_6: AB6C2D_mC40_12_ad_gh4i_j_bc ... 1532
448. Sr_2NiWO_6: AB6C2D_tI20_87_a_eh_d_b 1673
449. $Sr_3Ti_2O_7$: A7B3C2_tI24_139_aeg_be_e 1709
450. $Sr_4Ti_3O_{10}$: A10B4C3_tI34_139_c2eg_2e_ae 1704
451. $Sr(OH)_2(H_2O)_8$: A18B10C_tP116_130_2c4g_2c2g_a 1700
452. $SrCl_2{\cdot}(H_2O)_6$: A2B12C6D_hP21_150_d_2g_ef_a 1723
453. $SrCu_2(BO_3)_2$: A2B2C6D_tI44_121_i_i_ij_c 1685
454. $SrUO_4$: A4BC_oP24_57_cde_d_a 1615
455. Staurolite ($Al_5Fe_2O_{10}(OH)_2Si_2$): A5B2C10D2E2_mC84_12_acghj_bdi_5j_2i_j 1529
456. Steklite [$KAl(SO_4)_2$, $H3_2$]: ABC8D2_hP12_150_b_a_dg_d 1725
457. Sulphohalite [$Na_6ClF(SO_4)_2$, $H5_8$]: ABC6D8E2_cF72_225_b_a_e_f_c 1820
458. Swedenborgite ($NaBe_4SbO_7$, $E9_2$): A4BC7D_hP26_186_ac_b_a2c_b 1757
459. Ta_2NiS_5[††]: AB5C2_oC32_63_c_c2f_f 1658
460. Ta_2NiSe_5: AB5C2_mC32_15_e_e2f_f 1570
461. Ta_2PdSe_6: AB6C2_mC18_12_a_3i_i 1533
462. Ta_3Ti_{13} (BCC SQS-16): A3B13_oC32_38_ac_a2bcdef 1597
463. Ta_3Ti_5 (BCC SQS-16): A3B5_oC32_38_abce_abcdf 1597
464. Ta_5Ti_{11} (BCC SQS-16): A5B11_mP16_6_2abc_2a3b3c 1512
465. TaTi (BCC SQS-16): AB_aP16_2_4i_4i 1507
466. $TaTi_3$ (BCC SQS-16): AB3_mC32_8_4a_12a 1517
467. $TaTi_3$ (BCC SQS-16): AB3_mC32_8_4a_4a4b ... 1517
468. $TaTi_7$ (BCC SQS-16): AB7_hR16_166_c_c2h ... 1741
469. Tellurite (β-TeO_2, $C52$): A2B_oP24_61_2c_c 1625
470. Tennantite ($Cu_{12}As_4S_{13}$): A4B24C13_cI82_217_c_deg_ag 1798
471. Tetragonal $TlFeS_2$: AB2C_tI8_119_c_e_a 1684
472. Th_7S_{12} ($D8_k$): A3B2_hP20_176_2h_ah 1751
473. ThC_2 (C_g): A2B_mC12_15_f_e 1564
474. $ThCr_2Si_2$[∂‡‡]: A2B2C_tI10_139_d_e_a 1705
475. $ThFe_2SiC$: AB2CD_oC20_63_b_f_c_c 1656
476. Thenardite [Na_2SO_4 (V), $H1_7$]: A2B4C_oF56_70_g_h_a 1663
477. Ti_5Ga_4: A4B5_hP18_193_bg_dg 1763
478. $TiBe_{12}$ (approximate, $D2_a$): A12B_hP13_191_cdei_a 1761
479. Titanite ($CaTiSiO_5$, $S0_6$): AB5CD_mC32_15_e_e2f_e_b 1572
480. Tl_2AlF_5 ($K3_3$): AB5C2_oC32_20_b_a2bc_c 1581
481. $TlAlF_4$ ($H0_8$): AB4C_tP6_123_d_eh_a 1691
482. $TlCo_2S_2$[∂‡‡]: A2B2C_tI10_139_d_e_a 1705
483. Tolbachite ($CuCl_2$): A2B_mC6_12_i_a 1527
484. Topaz ($Al_2SiO_4F_2$, $S0_5$): A2B2C4D_oP36_62_d_d_2cd_c 1628
485. Tremolite ($Ca_2Mg_5Si_8O_{22}(OH)_2$, $S4_2$): A2B2C5D24E8_mC82_12_h_i_agh_2i5j_2j 1525
486. Tutton salt [$Cu(NH_4)_2(SO_4)_2{\cdot}6H_2O$, $H4_4$]: AB20C2D14E2_mP78_14_a_10e_e_7e_e 1547
487. U_6Mn ($D2_c$): AB6_tI28_140_a_hk 1716
488. V_3AsC: ABC3_oC20_63_c_b_cf 1659
489. V_4SiSb_2: A2BC4_tI28_140_h_a_k 1712
490. $VOSO_4$: A5BC_oP28_62_3cd_c_c 1640
491. VSe_2O_6: A6B2C_tP72_103_abc5d_2d_abc 1679

[∂] $TlCo_2S_2$ and $ThCr_2Si_2$ have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters.

492. Vesuvianite ($Ca_{10}Al_4(Mg,Fe)_2Si_9O_{34}(OH)_4$, $S2_3$): A4B10C2D34E4F9_tP252_126_k_ce2k_f_h8k_k_d2k 1693
493. $W_2O_3(PO_4)_2$: A11B2C2_mP60_4_22a_4a_4a 1507
494. Wülfingite (ϵ-$Zn(OH)_2$, $C31$): A2B2C_oP20_19_2a_2a_a 1577
495. Wollastonite ($CaSiO_3$): AB3C_aP30_2_3i_9i_3i 1505
496. Y_2SiO_5 (RE_2SiO_5 X2): A5BC2_mC64_15_5f_f_2f 1567
497. $Zn_{22}Zr$: A22B_cF184_227_cdfg_a 1826
498. $Zn_2Mo_3O_8$: A3B8C2_hP26_186_c_ab2c_2b 1755
499. $Zn(BrO_3)_2{\cdot}6H_2O$ ($J1_{10}$)[⊗⊗]: A2B6C6D_cP60_205_c_d_d_a 1778
500. $Zn(NH_3)_2Cl_2$ ($E1_2$): A2B6C2D_oI44_74_h_ij_i_e 1666
501. $Zr_{21}Re_{25}$: A25B21_hR92_167_b2e3f_e3f 1743
502. Zr_2Al_3[**]: A3B2_oF40_43_ab_b 1603
503. Zr_3Al_2: A2B3_tP20_136_j_dfg 1702
504. $ZrFe_4Si_2$: A4B2C_tP14_136_i_g_b 1702
505. ZrNiAl: ABC_hP9_189_g_ad_f 1760
506. ZrP_2O_7 High-Temperature ($K6_1$): A7B2C_cP40_205_bd_c_a 1782
507. $ZrSe_3$: A3B_mP8_11_3e_e 1521
508. $ZrTe_5$: A5B_oC24_63_c2f_c 1655
509. Zunyite [$Al_{13}(OH,F)_{18}Si_5O_{20}Cl$, $S0_8$]: A13BC18D20E5_cF228_216_dh_b_fh_2eh_ce 1791
510. "Martensite Type" FeC_x ($x \leq 0.06$) ($L'2_0$): AB_tI4_139_b_a 1712